

# IBM SPSS Statistics 19 Command Syntax Reference



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# Introduction: A Guide to Command Syntax

The *Command Syntax Reference* is arranged alphabetically by command name to provide quick access to detailed information about each command in the syntax command language. This introduction groups commands into broad functional areas. Some commands are listed more than once because they perform multiple functions, and some older commands that have been deprecated in favor of newer and better alternatives (but are still supported) are not included here. Changes to the command syntax language (since version 12.0), including modifications to existing commands and addition of new commands, are provided in the section [Release History](#).

## Core System

The Core system contains the core functionality plus a number of charting procedures. There are also numerous add-on modules that contain specialized functionality.

## Getting Data

You can read in a variety of data formats, including data files saved in IBM® SPSS® Statistics format, SAS datasets, database tables from many database sources, Excel and other spreadsheets, and text data files with both simple and complex structures.

Command	Description	Page Number
<b>SPSS Statistics Data Files</b>		
Get	Reads SPSS Statistics data files.	on p. 832
Import	Reads portable data files created with the Export command.	on p. 968
Add Files	Combines multiple data files by adding cases.	on p. 137
Match Files	Combines multiple data files by adding variables.	on p. 1119
Update	Replaces values in a master file with updated values.	on p. 2094
<b>Data Files Created by Other Applications</b>		
Get Translate	Reads spreadsheet and dBASE files.	on p. 856
Get Data	Reads Excel files, text data files, and database tables.	on p. 839
<b>Database Tables</b>		
Get Data	Reads Excel files, text data files, and database tables.	on p. 839
Get Capture	Reads database tables.	on p. 836
<b>SAS and Stata Data Files</b>		
Get SAS	Reads SAS dataset and SAS transport files.	on p. 848
Get Stata	Reads Stata data files.	on p. 853

Command	Description	Page Number
<b>Text Data Files</b>		
Get Data	Reads Excel files, text data files, and database tables.	on p. 839
Data List	Reads text data files.	on p. 550
Begin Data-End Data	Used with Data List to read inline text data.	on p. 227
<b>Complex (nested, mixed, grouped, etc.) Text Data Files</b>		
File Type	Defines mixed, nested, and grouped data structures.	on p. 726
Record Type	Used with File Type to read complex text data files.	on p. 1691
Input Program	Generates case data and/or reads complex data files.	on p. 975
End Case	Used with Input Program to define cases.	on p. 676
End File	Used with Input Program to indicate end of file.	on p. 682
Repeating Data	Used with Input Program to read input cases whose records contain repeating groups of data.	on p. 1735
Reread	Used with Input Program to reread a record.	on p. 1785
Keyed Data List	Reads data from nonsequential files: <ul style="list-style-type: none"> <li>■ Direct-access files, which provide direct access by a record number.</li> <li>■ Keyed files, which provide access by a record key.</li> </ul>	on p. 983
Point	Used with Keyed Data to establish the location at which sequential access begins (or resumes) in a keyed file.	on p. 1542
<b>Working with Multiple Data Sources</b>		
Dataset Name	Provides the ability to have multiple data sources open at the same time.	on p. 583
Dataset Activate	Makes the named dataset the active dataset.	on p. 572

### ***Saving and Exporting Data***

You can save data in numerous formats, including SPSS Statistics data file, Excel spreadsheet, database table, delimited text, and fixed-format text.

Command	Description	Page Number
<b>Saving Data in SPSS Statistics Format</b>		
Save	Saves the active dataset in SPSS Statistics format.	on p. 1807
Xsave	Saves data in SPSS Statistics format without requiring a separate data pass.	on p. 2183
Export	Saves data in portable format.	on p. 696
Save Data Collection	Saves a data file in SPSS Statistics format and a metadata file in IBM® SPSS® Data Collection MDD format for use in Data Collection applications.	on p. 1814
<b>Saving Data as Text</b>		
Write	Saves data as fixed-format text.	on p. 2159



Command	Description	Page Number
Save Translate	Saves data as tab-delimited text and comma-delimited (CSV) text.	on p. 1822
<b>Saving Data in Spreadsheet Format</b>		
Save Translate	Saves data in Excel and other spreadsheet formats and dBASE format.	on p. 1822
<b>Writing Data Back to a Database Table</b>		
Save Translate	Replaces or appends to existing database tables or creates new database tables.	on p. 1822

### **Statistics Adapter**

Command	Description	Page Number
Repository Attributes	Sets attributes for an object in a IBM® SPSS® Collaboration and Deployment Services Repository.	on p. 1774
Repository Connect	Establishes a connection to a IBM SPSS Collaboration and Deployment Services Repository and logs in the user.	on p. 1779
Repository Copy	Copies an arbitrary file from the local file system to a IBM SPSS Collaboration and Deployment Services Repository or copies a file from a IBM SPSS Collaboration and Deployment Services Repository to the local file system.	on p. 1782

### **Data Definition**

SPSS Statistics data files can contain more than simply data values. The **dictionary** can contain a variety of metadata attributes, including measurement level, display format, descriptive variable and value labels, and special codes for missing values.

Command	Description	Page Number
Apply Dictionary	Applies variable and file-based dictionary information from an external SPSS Statistics data file.	on p. 211
Datafile Attribute	Creates user-defined attributes that can be saved with the data file.	on p. 569
Variable Attribute	Creates user-defined variable attributes that can be saved with variables in the data file.	on p. 2127
Variable Labels	Assigns descriptive labels to variables.	on p. 2130
Value Labels	Assigns descriptive labels to data values.	on p. 2116
Add Value Labels	Assigns descriptive labels to data values.	on p. 144
Variable Level	Specifies the level of measurement (nominal, ordinal, or scale).	on p. 2132
Missing Values	Specifies values to be treated as missing.	on p. 1197
Rename	Changes variable names.	on p. 1733
Formats	Changes variable print and write formats.	on p. 750
Print Formats	Changes variable print formats.	on p. 1597
Write Formats	Changes variable write formats.	on p. 2164

Command	Description	Page Number
Variable Alignment	Specifies the alignment of data values in the Data Editor.	on p. 2126
Variable Width	Specifies the column width for display of variables in the Data Editor.	on p. 2135
Mrsets	Defines and saves multiple response set information.	on p. 1255

### **Data Transformations**

You can perform data transformations ranging from simple tasks, such as collapsing categories for analysis, to more advanced tasks, such as creating new variables based on complex equations and conditional statements.

Command	Description	Page Number
Autorecode	Recodes the values of string and numeric variables to consecutive integers.	on p. 219
Compute	Creates new numeric variables or modifies the values of existing string or numeric variables.	on p. 333
Count	Counts occurrences of the same value across a list of variables.	on p. 366
Create	Produces new series as a function of existing series.	on p. 381
Date	Generates date identification variables.	on p. 586
Leave	Suppresses reinitialization and retains the current value of the specified variable or variables when the program reads the next case.	on p. 1014
Numeric	Declares new numeric variables that can be referred to before they are assigned values.	on p. 1395
Rank	Produces new variables containing ranks, normal scores, and Savage and related scores for numeric variables.	on p. 1655
Recode	Changes, rearranges, or consolidates the values of an existing variable.	on p. 1685
RMV	Replaces missing values with estimates computed by one of several methods.	on p. 1797
Shift Values	Creates new variables that contain the values of existing variables from preceding or subsequent cases.	on p. 1884
String	Declares new string variables.	on p. 1933
Temporary	Signals the beginning of temporary transformations that are in effect only for the next procedure.	on p. 1959
TMS Begin	Indicates the beginning of a block of transformations to be exported to a file in PMML format (with SPSS Statistics extensions).	on p. 1965
TMS End	Marks the end of a block of transformations to be exported as PMML.	on p. 1971
TMS Import	Converts a PMML file containing ADP transformations into command syntax.	on p. 1973
TMS Merge	Merges a PMML file containing exported transformations with a PMML model file.	on p. 1976

### **File Information**

You can add descriptive information to a data file and display file and data attributes for the active dataset or any selected SPSS Statistics data file.

<b>Command</b>	<b>Description</b>	<b>Page Number</b>
Add Documents	Creates a block of text of any length in the active dataset.	on p. 135
Display	Displays information from the dictionary of the active dataset.	on p. 648
Document	Creates a block of text of any length in the active dataset.	on p. 672
Drop Documents	Deletes all text added with Document or Add Documents.	on p. 674
Sysfile Info	Displays complete dictionary information for all variables in a SPSS Statistics data file.	on p. 1956

### **File Transformations**

Data files are not always organized in the ideal form for your specific needs. You may want to combine data files, sort the data in a different order, select a subset of cases, or change the unit of analysis by grouping cases together. A wide range of file transformation capabilities is available.

<b>Command</b>	<b>Description</b>	<b>Page Number</b>
Delete Variables	Deletes variables from the data file.	on p. 614
Sort Cases	Reorders the sequence of cases based on the values of one or more variables.	on p. 1893
Weight	Case replication weights based on the value of a specified variable.	on p. 2151
<b>Select Subsets of Cases</b>		
Filter	Excludes cases from analysis without deleting them from the file.	on p. 740
N of Cases	Deletes all but the first $n$ cases in the data file.	on p. 1310
Sample	Selects a random sample of cases from the data file, deleting unselected cases.	on p. 1805
Select If	Selects cases based on logical conditions, deleting unselected cases.	on p. 1848
Split File	Splits the data into separate analysis groups based on values of one or more split variables.	on p. 1930
Use	Designates a range of observations for time series procedures.	on p. 2101
<b>Change File Structure</b>		
Aggregate	Aggregates groups of cases or creates new variables containing aggregated values.	on p. 160
Casestovars	Restructures complex data that has multiple rows for a case.	on p. 253
Varstocases	Restructures complex data structures in which information about a variable is stored in more than one column.	on p. 2136
Flip	Transposes rows (cases) and columns (variables).	on p. 746
<b>Merge Data Files</b>		
Add Files	Combines multiple SPSS Statistics data files or open datasets by adding cases.	on p. 137

Command	Description	Page Number
Match Files	Combines multiple SPSS Statistics data files or open datasets by adding variables.	on p. 1119
Update	Replaces values in a master file with updated values.	on p. 2094

### **Programming Structures**

As with other programming languages, the command syntax contains standard programming structures that can be used to do many things. These include the ability to perform actions only if some condition is true (if/then/else processing), repeat actions, create an array of elements, and use loop structures.

Command	Description	Page Number
Break	Used with Loop and Do If-Else If to control looping that cannot be fully controlled with conditional clauses.	on p. 241
Do If-Else If	Conditionally executes one or more transformations based on logical expressions.	on p. 656
Do Repeat	Repeats the same transformations on a specified set of variables.	on p. 666
If	Conditionally executes a single transformation based on logical conditions.	on p. 961
Loop	Performs repeated transformations specified by the commands within the loop until they reach a specified cutoff.	on p. 1056
Vector	Associates a vector name with a set of variables or defines a vector of new variables.	on p. 2143

### **Programming Utilities**

Command	Description	Page Number
Define	Defines a program macro.	on p. 595
Echo	Displays a specified text string as text output.	on p. 675
Execute	Forces the data to be read and executes the transformations that precede it in the command sequence.	on p. 695
Host	Executes external commands at the operating system level.	on p. 957
Include	Includes commands from the specified file.	on p. 972
Insert	Includes commands from the specified file.	on p. 979
Script	Runs the specified script file.	on p. 1841

### **General Utilities**

Command	Description	Page Number
Cache	Creates a copy of the data in temporary disk space for faster processing.	on p. 242
Clear Transformations	Discards all data transformation commands that have accumulated since the last procedure.	on p. 304
Erase	Deletes the specified file.	on p. 684
File Handle	Assigns a unique file handle to the specified file.	on p. 722
New File	Creates a blank, new active dataset.	on p. 1322

Command	Description	Page Number
Permissions	Changes the read/write permissions for the specified file.	on p. 1520
Preserve	Stores current Set command specifications that can later be restored by the Restore command.	on p. 1578
Print	Prints the values of the specified variables as text output.	on p. 1588
Print Eject	Displays specified information at the top of a new page of the output.	on p. 1594
Print Space	Displays blank lines in the output.	on p. 1600
Restore	Restores Set specifications that were stored by Preserve.	on p. 1796
Set	Customizes program default settings.	on p. 1860
Show	Displays current settings, many of which are set by the Set command.	on p. 1887
Subtitle	Inserts a subtitle on each page of output.	on p. 1935
Title	Inserts a title on each page of output.	on p. 1963

### ***Matrix Operations***

Command	Description	Page Number
Matrix	Using matrix programs, you can write your own statistical routines in the compact language of matrix algebra.	on p. 1128
Matrix Data	Reads raw matrix materials and converts them to a matrix data file that can be read by procedures that handle matrix materials.	on p. 1171
Mconvert	Converts covariance matrix materials to correlation matrix materials or vice versa.	on p. 1189

### ***Output Management System***

The Output Management System (OMS) provides the ability to automatically write selected categories of output to different output files in different formats, including SPSS Statistics data file format, HTML, XML, and text.

Command	Description	Page Number
OMS	Controls the routing and format of output. Output can be routed to external files in XML, HTML, text, and SAV (SPSS Statistics data file) formats.	on p. 1403
OMSEnd	Ends active OMS commands.	on p. 1434
OMSInfo	Displays a table of all active OMS commands.	on p. 1436
OMSLog	Creates a log of OMS activity.	on p. 1437

### ***Output Documents***

These commands control Viewer windows and files.

Command	Description	Page Number
Output Activate	Controls the routing of output to Viewer output documents.	on p. 1460
Output Close	Closes the specified Viewer document.	on p. 1463
Output Display	Displays a table of all open Viewer documents.	on p. 1465

Command	Description	Page Number
Output Export	Exports output to external files in various formats (e.g., Word, Excel, PDF, HTML, text).	on p. 1466
Output Name	Assigns a name to the active Viewer document. The name is used to refer to the output document in subsequent Output commands.	on p. 1486
Output New	Creates a new Viewer output document, which becomes the active output document.	on p. 1488
Output Open	Opens a Viewer document, which becomes the active output document. You can use this command to append output to an existing output document.	on p. 1486
Output Save	Saves the contents of an open output document to a file.	on p. 1494

### Charts

Command	Description	Page Number
Caseplot	Casewise plots of sequence and time series variables.	on p. 243
GGraph	Bar charts, pie charts, line charts, scatterplots, custom charts.	on p. 864
Pplot	Probability plots of sequence and time series variables.	on p. 1546
Spchart	Control charts, including X-Bar, r, s, individuals, moving range, and u.	on p. 1897

### Time Series

The Core system provides some basic time series functionality, including a number of time series chart types. Extensive time series analysis features are provided in the Forecasting option. [For more information, see the topic Add-On Modules on p. 9.](#)

Command	Description	Page Number
ACF	Displays and plots the sample autocorrelation function of one or more time series.	on p. 128
CCF	Displays and plots the cross-correlation functions of two or more time series.	on p. 294
		on p. 541
PACF	Displays and plots the sample partial autocorrelation function of one or more time series.	on p. 1506
Tsplot	Plot of one or more time series or sequence variables.	on p. 2050
Fit	Displays a variety of descriptive statistics computed from the residual series for evaluating the goodness of fit of models.	on p. 743
Predict	Specifies the observations that mark the beginning and end of the forecast period.	on p. 1556
Tset	Sets global parameters to be used by procedures that analyze time series and sequence variables.	on p. 2020
Tshow	Displays a list of all of the current specifications on the Tset, Use, Predict, and Date commands.	on p. 2023
Verify	Produces a report on the status of the most current Date, Use, and Predict specifications.	on p. 2149

## Add-On Modules

Add-on modules are not included with the Core system. The commands available to you will depend on your software license.

### Statistics Base

Command	Description	Page Number
ALSCAL	Multidimensional scaling (MDS) and multidimensional unfolding (MDU) using an alternating least-squares algorithm.	on p. 178
Cluster	Hierarchical clusters of items based on distance measures of dissimilarity or similarity. The items being clustered are usually cases, although variables can also be clustered.	on p. 305
Codebook	Reports the dictionary information — such as variable names, variable labels, value labels, missing values — and summary statistics for all or specified variables and multiple response sets in the active dataset.	on p. 323
Correlations	Pearson correlations with significance levels, univariate statistics, covariances, and cross-product deviations.	on p. 349
Crosstabs	Crosstabulations (contingency tables) and measures of association.	on p. 390
Curvefit	Fits selected curves to a line plot.	on p. 541
Descriptives	Univariate statistics, including the mean, standard deviation, and range.	on p. 615
Discriminant	Classifies cases into one of several mutually exclusive groups based on their values for a set of predictor variables.	on p. 629
Examine	Descriptive statistics, stem-and-leaf plots, histograms, boxplots, normal plots, robust estimates of location, and tests of normality.	on p. 685
Factor	Identifies underlying variables, or factors, that explain the pattern of correlations within a set of observed variables.	on p. 704
Frequencies	Tables of counts and percentages and univariate statistics, including the mean, median, and mode.	on p. 753
Graph	Bar charts, pie charts, line charts, histograms, scatterplots, etc.	on p. 926
KNN	Classifies and predicts cases based upon the values “nearest neighboring” cases.	on p. 998
Linear	Creates a predictive model for a continuous target.	on p. 1017
List	Individual case listing.	on p. 1024
Means	Group means and related univariate statistics for dependent variables within categories of one or more independent variables.	on p. 1192
Mult Response	Frequency tables and crosstabulations for multiple-response data.	on p. 1259
Nonparametric	Collection of one-sample, independent samples, and related samples nonparametric tests.	on p. 1360
Nonpar Corr	Rank-order correlation coefficients: Spearman’s rho and Kendall’s tau-b, with significance levels.	on p. 1354
Npar Tests	Collection of one-sample, independent samples, and related samples nonparametric tests.	on p. 1376

<b>Command</b>	<b>Description</b>	<b>Page Number</b>
OLAP Cubes	Summary statistics for scale variables within categories defined by one or more categorical grouping variables.	on p. 1397
Oneway	One-way analysis of variance.	on p. 1439
Partial Corr	Partial correlation coefficients between two variables, adjusting for the effects of one or more additional variables.	on p. 1512
Plum	Analyzes the relationship between a polytomous ordinal dependent variable and a set of predictors.	on p. 1534
Proximities	Measures of similarity, dissimilarity, or distance between pairs of cases or pairs of variables.	on p. 1614
Quick Cluster	When the desired number of clusters is known, this procedure groups cases efficiently into clusters.	on p. 1648
Ratio Statistics	Descriptive statistics for the ratio between two variables.	on p. 1661
Regression	Multiple regression equations and associated statistics and plots.	on p. 1700
Reliability	Estimates reliability statistics for the components of multiple-item additive scales.	on p. 1724
Report	Individual case listing and group summary statistics.	on p. 1749
ROC	Receiver operating characteristic (ROC) curve and an estimate of the area under the curve.	on p. 1801
Summarize	Individual case listing and group summary statistics.	on p. 1937
TTest	One sample, independent samples, and paired samples <i>t</i> tests.	on p. 2060
Twostep Cluster	Groups observations into clusters based on a nearness criterion. The procedure uses a hierarchical agglomerative clustering procedure in which individual cases are successively combined to form clusters whose centers are far apart.	on p. 2065
Unianova	Regression analysis and analysis of variance for one dependent variable by one or more factors and/or variables.	on p. 2073
Xgraph	Creates 3-D bar charts, population pyramids, and dot plots.	on p. 2167

### **Advanced Statistics**

<b>Command</b>	<b>Description</b>	<b>Page Number</b>
Coxreg	Cox proportional hazards regression for analysis of survival times.	on p. 368
Genlin	Generalized Linear Model. Genlin allows you to fit a broad spectrum of “generalized” models in which the distribution of the error term need not be normal and the relationship between the dependent variable and predictors need only be linear through a specified transformation.	on p. 762
Genlinmixed	Generalized linear mixed models extend the linear model so that the target is linearly related to the factors and covariates via a specified link function, the target can have a non-normal distribution, and the observations can be correlated. Generalized linear mixed models cover a wide variety of models, from simple linear regression to complex multilevel models for non-normal longitudinal data.	on p. 798



Command	Description	Page Number
Genlog	A general procedure for model fitting, hypothesis testing, and parameter estimation for any model that has categorical variables as its major components.	on p. 820
GLM	General Linear Model. A general procedure for analysis of variance and covariance, as well as regression.	on p. 883
Hiloglinear	Fits hierarchical loglinear models to multidimensional contingency tables using an iterative proportional-fitting algorithm.	on p. 940
KM	Kaplan-Meier (product-limit) technique to describe and analyze the length of time to the occurrence of an event.	on p. 989
Mixed	The mixed linear model expands the general linear model used in the GLM procedure in that the data are permitted to exhibit correlation and non-constant variability.	on p. 1200
Survival	Actuarial life tables, plots, and related statistics.	on p. 1943
Varcomp	Estimates variance components for mixed models.	on p. 2119

### **Regression**

Command	Description	Page Number
Logistic Regression	Regresses a dichotomous dependent variable on a set of independent variables.	on p. 1028
Nomreg	Fits a multinomial logit model to a polytomous nominal dependent variable.	on p. 1340
NLR, CNLR	Nonlinear regression is used to estimate parameter values and regression statistics for models that are not linear in their parameters.	on p. 1323
WLS	Weighted Least Squares. Estimates regression models with different weights for different cases.	on p. 2153
2SLS	Two-stage least-squares regression.	on p. 123

### **Custom Tables**

Command	Description	Page Number
Ctables	Produces tables in one, two, or three dimensions and provides a great deal of flexibility for organizing and displaying the contents.	on p. 510

### **Decision Trees**

Command	Description	Page Number
Tree	Tree-based classification models.	on p. 1978

### **Categories**

Command	Description	Page Number
Catreg	Categorical regression with optimal scaling using alternating least squares.	on p. 278

<b>Command</b>	<b>Description</b>	<b>Page Number</b>
CatPCA	Principal components analysis.	on p. 261
Overals	Nonlinear canonical correlation analysis on two or more sets of variables.	on p. 1497
Correspondence	Displays the relationships between rows and columns of a two-way table graphically by a scatterplot matrix.	on p. 355
Multiple Correspondence	Quantifies nominal (categorical) data by assigning numerical values to the cases (objects) and categories, such that objects within the same category are close together and objects in different categories are far apart.	on p. 1269
Proxscal	Multidimensional scaling of proximity data to find a least-squares representation of the objects in a low-dimensional space.	on p. 1631

### ***Complex Samples***

<b>Command</b>	<b>Description</b>	<b>Page Number</b>
CSPlan	Creates a complex sample design or analysis specification.	on p. 476
CSSelect	Selects complex, probability-based samples from a population.	on p. 495
CSDescriptives	Estimates means, sums, and ratios, and computes their standard errors, design effects, confidence intervals, and hypothesis tests.	on p. 424
CSTabulate	Frequency tables and crosstabulations, and associated standard errors, design effects, confidence intervals, and hypothesis tests.	on p. 504
CSGLM	Linear regression analysis, and analysis of variance and covariance.	on p. 430
CSLogistic	Logistic regression analysis on a binary or multinomial dependent variable using the generalized link function.	on p. 444
CSOrdinal	Fits a cumulative odds model to an ordinal dependent variable for data that have been collected according to a complex sampling design.	on p. 459

### ***Neural Networks***

<b>Command</b>	<b>Description</b>	<b>Page Number</b>
MLP	Fits flexible predictive model for one or more target variables, which can be categorical or scale, based upon the values of factors and covariates.	on p. 1223
RBF	Fits flexible predictive model for one or more target variables, which can be categorical or scale, based upon the values of factors and covariates. Generally trains faster than MLP at the slight cost of some model flexibility.	on p. 1666

**Forecasting**

Command	Description	Page Number
Season	Estimates multiplicative or additive seasonal factors.	on p. 1843
Spectra	Periodogram and spectral density function estimates for one or more series.	on p. 1922
Tsapply	Loads existing time series models from an external file and applies them to data.	on p. 2006
Tsmodel	Estimates exponential smoothing, univariate Autoregressive Integrated Moving Average (ARIMA), and multivariate ARIMA (or transfer function models) models for time series, and produces forecasts.	on p. 2024

**Conjoint**

Command	Description	Page Number
Conjoint	Analyzes score or rank data from full-concept conjoint studies.	on p. 339
Orthoplan	Orthogonal main-effects plan for a full-concept conjoint analysis.	on p. 1455
Plancards	Full-concept profiles, or cards, from a plan file for conjoint analysis.	on p. 1521

**Bootstrapping**

Command	Description	Page Number
Bootstrap	Bootstrapping is an alternative to parametric estimates when the assumptions of those methods are in doubt, or where parametric inference is impossible or requires very complicated formulas for the calculation of standard errors.	on p. 234

**Missing Values**

Command	Description	Page Number
Multiple Imputation	Performs multiple imputations of missing values. Many other procedures can analyze a multiply-imputed dataset to produce pooled results which are more accurate than the singly-imputed datasets produced by MVA.	on p. 1283
MVA	Missing Value Analysis. Describes missing value patterns and estimates (imputes) missing values.	on p. 1296

**Data Preparation**

Command	Description	Page Number
Adp	Automatically prepares data for modeling.	on p. 147
Detectanomaly	Searches for unusual cases based on deviations from the norms of their cluster groups.	on p. 620

Command	Description	Page Number
Validatedata	Identifies suspicious and invalid cases, variables, and data values in the active dataset.	on p. 2104
Optimal Binning	Discretizes scale “binning input” variables to produce categories that are “optimal” with respect to the relationship of each binning input variable with a specified categorical guide variable.	on p. 1449

## ***Release History***

This section details changes to the command syntax language occurring after release 12.0. Information is organized alphabetically by command and changes for a given command are grouped by release. For commands introduced after 12.0, the introductory release is noted. Additions of new functions (used for instance with `COMPUTE`) and changes to existing functions are detailed under the heading *Functions*, located at the end of this section.

### ***ADP***

Release 18

- Command introduced.

### ***AGGREGATE***

Release 13.0

- `MODE` keyword introduced.
- `OVERWRITE` keyword introduced.

Release 17.0

- `AGGREGATE` runs without a break variable.

### ***ALTER TYPE***

Release 16.0

- Command introduced.

### ***APPLY DICTIONARY***

Release 14.0

- `ATTRIBUTES` keyword introduced on `FILEINFO` and `VARINFO` subcommands.

Release 18

- `ROLE` keyword introduced on `VARINFO` subcommands.

### ***AUTORECODE***

Release 13.0

- `BLANK` subcommand introduced.
- `GROUP` subcommand introduced.
- `APPLY TEMPLATE` and `SAVE TEMPLATE` subcommands introduced.

### ***BEGIN GPL***

Release 14.0

- Command introduced.

### ***BEGIN PROGRAM***

Release 14.0

- Command introduced.

### ***BOOTSTRAP***

Release 18

- Command introduced.

### ***CASEPLOT***

Release 14.0

- For plots with one variable, new option to specify a value with the `REFERENCE` keyword on the `FORMAT` subcommand.

### ***CATPCA***

Release 13.0

- `NDIM` keyword introduced on `PLOT` subcommand.
- The maximum label length on the `PLOT` subcommand is increased to 64 for variable names, 255 for variable labels, and 60 for value labels (previous value was 20).

### ***CATREG***

Release 13.0

- The maximum category label length on the `PLOT` subcommand is increased to 60 (previous value was 20).

Release 17.0

- `MULTISTART` and `FIXSIGNS` keywords added to `INITIAL` subcommand.
- `REGULARIZATION` subcommand added.
- `RESAMPLE` subcommand added.
- `REGU` keyword added to `PRINT` subcommand.

- REGU keyword added to PLOT subcommand.
- SUPPLEMENTARY categories not occurring in data used to create the model are now interpolated.

***CD***

Release 13.0

- Command introduced.

***CODEBOOK***

Release 17.0

- Command introduced.

Release 18

- ROLE keyword added to VARINFO subcommand.

***CORRESPONDENCE***

Release 13.0

- For the NDIM keyword on the PLOT subcommand, the default is changed to all dimensions.
- The maximum label length on the PLOT subcommand is increased to 60 (previous value was 20).

***CROSSTABS***

Release 19.0

- HIDESMALLCOUNTS subcommand introduced.
- SHOWDIM subcommand introduced.
- PROP and BPROP keywords introduced on the CELLS subcommand.

***CSGLM***

Release 13.0

- Command introduced.

***CSLOGISTIC***

Release 13.0

- Command introduced.

Release 17.0

- Added support for SET THREADS.

**CSORDINAL**

Release 15.0

- Command introduced.

Release 17.0

- Added support for SET THREADS.

**CTABLES**

Release 13.0

- HSUBTOTAL keyword introduced on the CATEGORIES subcommand.

Release 14.0

- INCLUDEMRSETS keyword introduced on the SIGTEST and COMPARETEST subcommands.
- CATEGORIES keyword introduced on the SIGTEST and COMPARETEST subcommands.
- MEANSVARIANCE keyword introduced on the COMPARETEST subcommand.

Release 18.0

- MERGE keyword introduced on the COMPARETEST subcommand.
- PCOMPUTE and PPROPERTIES subcommands introduced.

Release 19.0

- HIDESMALLCOUNTS subcommand introduced.

**CURVEFIT**

Release 19.0

- TEMPLATE subcommand introduced.

**DATA LIST**

Release 16.0

- ENCODING subcommand added for Unicode support.

**DATAFILE ATTRIBUTE**

Release 14.0

- Command introduced.

**DATASET ACTIVATE**

Release 14.0

- Command introduced.

***DATASET CLOSE***

Release 14.0

- Command introduced.

***DATASET COPY***

Release 14.0

- Command introduced.

***DATASET DECLARE***

Release 14.0

- Command introduced.

***DATASET DISPLAY***

Release 14.0

- Command introduced.

***DATASET NAME***

Release 14.0

- Command introduced.

***DEFINE-!ENDDFINE***

Release 14.0

- For syntax processed in interactive mode, modifications to the macro facility may affect macro calls occurring at the end of a command. [For more information, see the topic Overview on p. 596.](#)

***DETECTANOMALY***

Release 14.0

- Command introduced.

***DISPLAY***

Release 14.0

- ATTRIBUTES keyword introduced.

Release 15.0

- @ATTRIBUTES keyword introduced.



**DO REPEAT-END REPEAT**

Release 14.0

- ALL keyword introduced.

**EXTENSION**

Release 16.0

- Command introduced.

**FILE HANDLE**

Release 13.0

- The NAME subcommand is modified to accept a path and/or file.

Release 16.0

- ENCODING subcommand added for Unicode support.

**FILE TYPE**

Release 16.0

- ENCODING subcommand added for Unicode support.

**GENLIN**

Release 15.0

- Command introduced.

Release 16.0

- Added multinomial and tweedie distributions; added MLE estimation option for ancillary parameter of negative binomial distribution (MODEL subcommand, DISTRIBUTION keyword). Notes related to the addition of the new distributions added throughout.
- Added cumulative Cauchit, cumulative complementary log-log, cumulative logit, cumulative negative log-log, and cumulative probit link functions (MODEL subcommand, LINK keyword).
- Added likelihood-ratio chi-square statistics as an alternative to Wald statistics (CRITERIA subcommand, ANALYSISTYPE keyword).
- Added profile likelihood confidence intervals as an alternative to Wald confidence intervals (CRITERIA subcommand, CTYPE keyword).
- Added option to specify initial value for ancillary parameter of negative binomial distribution (CRITERIA subcommand, INITIAL keyword).
- Changed default display of the likelihood function for GEEs to show the full value instead of the kernel (CRITERIA subcommand, LIKELIHOOD keyword).

**GENLINMIXED**

Release 19

- Command introduced.

**GET CAPTURE**

Release 15.0

- UNENCRYPTED subcommand introduced.

**GET DATA**

Release 13.0

- ASSUMEDSTRWIDTH subcommand introduced for TYPE=ODBC.

Release 14.0

- ASSUMEDSTRWIDTH subcommand extended to TYPE=XLS.
- TYPE=OLEDB introduced.

Release 15.0

- ASSUMEDSTRWIDTH subcommand extended to TYPE=OLEDB.

Release 16.0

- TYPE=XLSX and TYPE=XLSM introduced.

Release 17.0

- ENCRYPTED subcommand introduced.

**GET SAS**

Release 19

- ENCODING subcommand introduced.

**GET STATA**

Release 14.0

- Command introduced.

Release 19

- ENCODING subcommand introduced.

**GGRAPH**

Release 14.0

- Command introduced.

**Release 15.0**

- `RENAME` syntax qualifier deprecated.
- `COUNTCI`, `MEDIANCI`, `MEANCI`, `MEANSD`, and `MEANSE` functions introduced.

**Release 17.0**

- Added `SOURCE=VIZTEMPLATE` to support visualization templates.
- Added `VIZSTYLESHEET` keyword to support visualization stylesheets.

**Release 19.0**

- Added `LOCATION=FILE` to support visualization templates stored in an arbitrary location on the file system.

***GLM*****Release 17.0**

- `POSTHOC` subcommand: `T2`, `T3`, `GH`, and `C` keywords are not valid when multiple factors in the model.
- `PLOT` subcommand: new `WITH` keyword allows you to fix covariate values for profile plots.

***GRAPH*****Release 13.0**

- `PANEL` subcommand introduced.
- `INTERVAL` subcommand introduced.

***HOST*****Release 13.0**

- Command introduced.

***INCLUDE*****Release 16.0**

- `ENCODING` keyword added for Unicode support.

***INSERT*****Release 13.0**

- Command introduced.

**Release 16.0**

- `ENCODING` keyword added for Unicode support.

**KEYED DATA LIST**

Release 16.0

- `ENCODING` subcommand added for Unicode support.

**KNN**

Release 17.0

- Command introduced.

**LINEAR**

Release 19

- Command introduced.

**LOGISTIC REGRESSION**

Release 13.0

- `OUTFILE` subcommand introduced.

Release 14.0

- Modification to the method of recoding string variables. [For more information, see the topic Overview on p. 1029.](#)

**MISSING VALUES**

Release 16.0

- Limitation preventing assignment of missing values to strings with a defined width greater than eight bytes removed.

**MLP**

Release 16.0

- Command introduced.

**MODEL CLOSE**

Release 13.0

- Command introduced.

**MODEL HANDLE**

Release 13.0

- Command introduced.

**MODEL LIST**

Release 13.0

- Command introduced.

**MRSETS**

Release 14.0

- LABELSOURCE keyword introduced on MDGROUP subcommand.
- CATEGORYLABELS keyword introduced on MDGROUP subcommand.

**MULTIPLE CORRESPONDENCE**

Release 13.0

- Command introduced.

**MULTIPLE IMPUTATION**

Release 17.0

- Command introduced.

**NAIVEBAYES**

Release 14.0

- Command introduced.

**NOMREG**

Release 13.0

- ENTRYMETHOD keyword introduced on STEPWISE subcommand.
- REMOVALMETHOD keyword introduced on STEPWISE subcommand.
- IC keyword introduced on PRINT subcommand.

Release 15.0

- ASSOCIATION keyword introduced on PRINT subcommand.

Release 17.0

- Added support for SET THREADS and SET MCACHE.

**NONPARAMETRIC**

Release 18

- Command introduced.

***NPAR TESTS***

Release 17.0

- Increased limits on number of variables allowed in the analysis.

***OLAP CUBES***

Release 19.0

- HIDESMALLCOUNTS subcommand introduced.

***OMS***

Release 13.0

- TREES keyword introduced on SELECT subcommand.
- IMAGES, IMAGEROOT, CHARTSIZE, and IMAGEFORMAT keywords introduced on DESTINATION subcommand.

Release 14.0

- XMLWORKSPACE keyword introduced on DESTINATION subcommand.

Release 16.0

- IMAGEFORMAT=VML introduced for FORMAT=HTML on DESTINATION subcommand.
- IMAGEMAP keyword introduced for FORMAT=HTML on DESTINATION subcommand.
- FORMAT=SPV introduced for saving output in Viewer format.
- CHARTFORMAT keyword introduced.
- TREEFORMAT keyword introduced.
- TABLES keyword introduced.
- FORMAT=SVWSOXML is no longer supported.

Release 17.0

- MODELS keyword introduced on SELECT subcommand.
- FORMAT=DOC, XLS, PDF, and SPW introduced.
- MODELFORMAT keyword introduced.

Release 19.0

- IMAGEFORMAT=VML introduced for FORMAT=OXML on DESTINATION subcommand.

***ONEWAY***

Release 19.0

- TEMPLATE subcommand introduced.

**OPTIMAL BINNING**

Release 15.0

- Command introduced.

**OUTPUT ACTIVATE**

Release 15.0

- Command introduced.

**OUTPUT CLOSE**

Release 15.0

- Command introduced.

**OUTPUT DISPLAY**

Release 15.0

- Command introduced.

**OUTPUT EXPORT**

Release 17.0

- Command introduced.

**OUTPUT NAME**

Release 15.0

- Command introduced.

**OUTPUT NEW**

Release 15.0

- Command introduced.

Release 16.0

- `TYPE` keyword is obsolete and is ignored.

**OUTPUT OPEN**

Release 15.0

- Command introduced.

Release 17.0

- `LOCK` keyword introduced.

***OUTPUT SAVE***

Release 15.0

- Command introduced.

Release 16.0

- TYPE keyword introduced.

Release 17.0

- LOCK keyword introduced.

***PER ATTRIBUTES***

Release 16.0

- Command introduced as PER ATTRIBUTES.

Release 17.0

- VERSIONLABEL keyword extended to support multiple labels.

Release 18.0

- PER ATTRIBUTES deprecated. Command name changed to REPOSITORY ATTRIBUTES.

***PER CONNECT***

Release 15.0

- Command introduced as PER CONNECT.

Release 17.0

- DOMAIN keyword deprecated on the LOGIN subcommand.
- PROVIDER keyword introduced on the LOGIN subcommand.

Release 18.0

- PER CONNECT deprecated. Command name changed to REPOSITORY CONNECT.

***PER COPY***

Release 16.0

- Command introduced as PER COPY.

Release 18.0

- PER COPY deprecated. Command name changed to REPOSITORY COPY.



**PLANCARDS**

Release 14.0

- `PAGINATE` subcommand is obsolete and no longer supported.

**PLS**

Release 16.0

- Command introduced.

**POINT**

Release 16.0

- `ENCODING` subcommand added for Unicode support.

**PLOT**

Release 19.0

- `TEMPLATE` subcommand introduced.

**PREFSCAL**

Release 14.0

- Command introduced.

**PRINT**

Release 16.0

- `ENCODING` subcommand added for Unicode support.

**PRINT EJECT**

Release 16.0

- `ENCODING` subcommand added for Unicode support.

**PRINT SPACE**

Release 16.0

- `ENCODING` subcommand added for Unicode support.

**RBF**

Release 16.0

- Command introduced.

**REGRESSION**

Release 13.0

- `PARAMETER` keyword introduced on `OUTFILE` subcommand.

Release 16.0

- Added support for `SET THREADS` and `SET MCACHE`.

Release 17.0

- Added option to specify confidence level on `CI` keyword of `STATISTICS` subcommand.

Release 19.0

- `TEMPLATE` subcommand introduced.

**RELIABILITY**

Release 17.0

- Increased limits on numbers of variables allowed on the `VARIABLES` and `SCALE` lists.

**REPEATING DATA**

Release 16.0

- `ENCODING` subcommand added for Unicode support.

**REPOSITORY ATTRIBUTES**

Release 16.0

- Command introduced as `PER ATTRIBUTES`.

Release 17.0

- `VERSIONLABEL` keyword extended to support multiple labels.

Release 18.0

- `PER ATTRIBUTES` deprecated. Command name changed to `REPOSITORY ATTRIBUTES`.

**REPOSITORY CONNECT**

Release 15.0

- Command introduced as `PER CONNECT`.

Release 17.0

- `DOMAIN` keyword deprecated on the `LOGIN` subcommand.
- `PROVIDER` keyword introduced on the `LOGIN` subcommand.

Release 18.0

- `PER CONNECT` deprecated. Command name changed to `REPOSITORY CONNECT`.

### ***REPOSITORY COPY***

Release 16.0

- Command introduced as `PER COPY`.

Release 18.0

- `PER COPY` deprecated. Command name changed to `REPOSITORY COPY`.

### ***RESPONSE RATE***

Release 18.0

- Command introduced.

### ***ROC***

Release 18.0

- `MODELQUALITY` keyword introduced.

### ***SAVE DATA COLLECTION***

Release 15.0

- Command introduced as `SAVE DIMENSIONS`.

Release 18.0

- `SAVE DIMENSIONS` deprecated. Command name changed to `SAVE DATA COLLECTION`.

### ***SAVE TRANSLATE***

Release 14.0

- Value `STATA` added to list for `TYPE` subcommand.
- `EDITION` subcommand introduced for `TYPE=STATA`.
- `SQL` subcommand introduced.
- `MISSING` subcommand introduced.
- Field/column names specified on the `RENAME` subcommand can contain characters (for example, spaces, commas, slashes, plus signs) that are not allowed in IBM® SPSS® Statistics variable names.
- Continuation lines for connection strings on the `CONNECT` subcommand do not need to begin with a plus sign.

Release 15.0

- `ENCRYPTED` subcommand introduced.

- Value `CSV` added to list for `TYPE` subcommand.
- `TEXTOPTIONS` subcommand introduced for `TYPE=CSV` and `TYPE=TAB`.

#### Release 16.0

- `VERSION=12` introduced for writing data in Excel 2007 XLSX format with `TYPE=XLS`.

#### Release 17.0

- `UNENCRYPTED` subcommand introduced.

#### Release 18.0

- `VERSION=9` introduced for writing SAS 9+ files with `TYPE=SAS`.

#### Release 19

- `ENCODING` subcommand introduced.

### ***SCRIPT***

#### Release 16.0

- Scripts run from the `SCRIPT` command now run synchronously with the command syntax stream.

#### Release 17.0

- Ability to run Python scripts introduced.

### ***SELECTPRED***

#### Release 14.0

- Command introduced.

### ***SET***

#### Release 13.0

- `RNG` and `MTINDEX` subcommands introduced.
- Default for `MXERRS` subcommand increased to 100.
- `SORT` subcommand introduced.
- `LOCALE` subcommand introduced.

#### Release 14.0

- Default for `WORKSPACE` subcommand increased to 6144.

#### Release 15.0

- `LABELS` replaces `VALUES` as the default for the `TNUMBERS` subcommand.

- JOURNAL subcommand is obsolete and no longer supported.
- Value EXTERNAL added to list for SORT subcommand, replacing the value INTERNAL as the default. Value SS is deprecated.

#### Release 16.0

- MCACHE subcommand introduced.
- THREADS subcommand introduced.
- UNICODE subcommand introduced.

#### Release 16.0.1

- BOTHLARGE keyword introduced for the TFIT subcommand.

#### Release 17.0

- FUZZBITS subcommand introduced.
- MIOUTPUT subcommand introduced.

#### Release 18.0

- ROWSBREAK, CELLSBREAK, and TOLERANCE subcommands introduced for controlling display of large pivot tables.
- ZCOMPRESSION subcommand introduced.
- COMPRESSION subcommand is obsolete and ignored.
- REPDEFER subcommand introduced.

#### Release 19.0

- XVERSION subcommand introduced.
- OATTRS subcommand introduced.
- DIGITGROUPING subcommand introduced.
- TABLERENDER subcommand introduced.
- CMPTRANS subcommand introduced.

### **SHIFT VALUES**

#### Release 17.0

- Command introduced.

### **SHOW**

#### Release 13.0

- BLKSIZE and BUFNO subcommands are obsolete and no longer supported.
- HANDLES subcommand introduced.
- SORT subcommand introduced.

**Release 15.0**

- TMSRECORDING subcommand introduced.

**Release 16.0**

- UNICODE subcommand introduced.
- MCACHE subcommand introduced.
- THREADS subcommand introduced.

**Release 17.0**

- FUZZBITS subcommand introduced.

**Release 18.0**

- EXTPATHS subcommand introduced.
- ZCOMPRESSION subcommand introduced.
- COMPRESSION subcommand removed because it is obsolete.
- REPDEFER subcommand introduced.

**Release 19.0**

- TABLERENDER subcommand introduced.
- XVERSION subcommand introduced.
- OATTRS subcommand introduced.
- DIGITGROUPING subcommand introduced.
- CMPTRANS subcommand introduced.

***SORT VARIABLES*****Release 16.0.**

- Command introduced.

**Release 18.0.**

- ROLE keyword introduced.

***SPCHART*****Release 15.0**

- (XBARONLY) keyword introduced on XR and XS subcommands.
- RULES subcommand introduced.
- ID subcommand introduced.

**Release 19.0**

- CPCHART subcommand introduced.
- NORMAL subcommand introduced.

- REFERENCE subcommand introduced.
- Following keywords introduced on STATISTICS subcommand: N, MEAN, STDDEV, CAPSIGMA, LSL, USL, TARGET, AZLOUT, AZUOUT, CZLOUT, CZUOUT, PZLOUT, PZUOUT.

### **TMS BEGIN**

Release 15.0

- Command introduced.

Release 16.0

- Added support for new string functions CHAR.CONCAT, CHAR.LENGTH, and CHAR.SUBSTR within TMS blocks.

### **TMS END**

Release 15.0

- Command introduced.

### **TMS IMPORT**

Release 18

- Command introduced.

### **TMS MERGE**

Release 15.0

- Command introduced.

### **TREE**

Release 13.0

- Command introduced.

Release 18.0

- TARGETRESPONSE subcommand introduced.

### **TSAPPLY**

Release 14.0

- Command introduced.

### **TSMODEL**

Release 14.0

- Command introduced.

***TSPLIT***

Release 14.0

- For plots with one variable, REFERENCE keyword modified to allow specification of a value.

***UNIANOVA***

Release 17.0

- POSTHOC subcommand: T2, T3, GH, and C keywords are not valid when multiple factors in the model.

***VALIDATEDATA***

Release 14.0

- Command introduced.

***VALUE LABELS***

Release 14.0

- The maximum length of a value label is extended to 120 bytes (previous limit was 60 bytes).

Release 16.0

- Limitation preventing assignment of missing values to strings with a defined width greater than eight bytes removed.

***VARIABLE ATTRIBUTE***

Release 14.0

- Command introduced.

***VARIABLE ROLE***

Release 18.0

- Command introduced.

***WRITE***

Release 16.0

- ENCODING subcommand added for Unicode support.

***XGRAPH***

Release 13.0

- Command introduced.



## **Functions**

### Release 13.0

- `APPLYMODEL` and `STRAPPLYMODEL` functions introduced.
- `DATEDIFF` and `DATESUM` functions introduced.

### Release 14.0

- `REPLACE` function introduced.
- `VALUELABEL` function introduced.

### Release 16.0

- `CHAR.INDEX` function introduced.
- `CHAR.LENGTH` function introduced.
- `CHAR.LPAD` function introduced.
- `CHAR.MBLEN` function introduced.
- `CHAR.RINDEX` function introduced.
- `CHAR.RPAD` function introduced.
- `CHAR.SUBSTR` function introduced.
- `NORMALIZE` function introduced.
- `NTRIM` function introduced.
- `STRUNC` function introduced.

### Release 17.0

- `MEDIAN` function introduced.
- *mult* and *fuzzbits* arguments introduced for the `RND` and `TRUNC` functions.
- `NEIGHBOR` and `DISTANCE` functions added to `APPLYMODEL` and `STRAPPLYMODEL`.

# Universals

This part of the *Command Syntax Reference* discusses general topics pertinent to using command syntax. The topics are divided into five sections:

- **Commands** explains command syntax, including command specification, command order, and running commands in different modes. In this section, you will learn how to read syntax charts, which summarize command syntax in diagrams and provide an easy reference. Discussions of individual commands are found in an alphabetical reference in the next part of this manual.
- **Files** discusses different types of files used by the program. Terms frequently mentioned in this manual are defined. This section provides an overview of how files are handled.
- **Variables and Variable Types and Formats** contain important information about general rules and conventions regarding variables and variable definition.
- **Transformations** describes expressions that can be used in data transformation. Functions and operators are defined and illustrated. In this section, you will find a complete list of available functions and how to use them.

## Commands

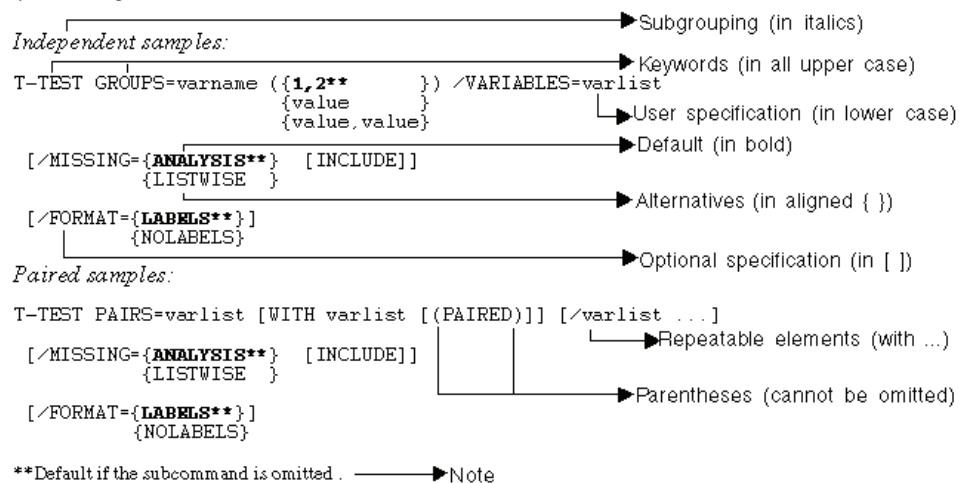
**Commands** are the instructions that you give the program to initiate an action. For the program to interpret your commands correctly, you must follow certain rules.

### *Syntax Diagrams*

Each command described in this manual includes a syntax diagram that shows all of the subcommands, keywords, and specifications allowed for that command. By recognizing symbols and different type fonts, you can use the syntax diagram as a quick reference for any command.

- Lines of text in italics indicate limitation or operation mode of the command.
- Elements shown in upper case are keywords to identify commands, subcommands, functions, operators, and other specifications. In the sample syntax diagram below, `T-TEST` is the command and `GROUPS` is a subcommand.
- Elements in lower case describe specifications that you supply. For example, `varlist` indicates that you need to supply a list of variables.
- Elements in bold are defaults. There are two types of defaults. When the default is followed by `**`, as `ANALYSIS**` is in the sample syntax diagram below, the default (`ANALYSIS`) is in effect if the subcommand (`MISSING`) is not specified. If a default is not followed by `**`, it is in effect when the subcommand (or keyword) is specified by itself.

Figure 2-1  
Syntax diagram



- Parentheses, apostrophes, and quotation marks are required where indicated.
- Unless otherwise noted, elements enclosed in square brackets ([ ]) are optional. For some commands, square brackets are part of the required syntax. The command description explains which specifications are required and which are optional.
- Braces ({ }) indicate a choice between elements. You can specify any one of the elements enclosed within the aligned braces.
- Ellipses indicate that you can repeat an element in the specification. The specification `T-TEST PAIRS=varlist [WITH varlist [(PAIRED)]] [/varlist ...]` means that you can specify multiple variable lists with optional `WITH` variables and the keyword `PAIRED` in parentheses.
- Most abbreviations are obvious; for example, `varname` stands for variable name and `varlist` stands for a variable list.
- The command terminator is not shown in the syntax diagram.

### Command Specification

The following rules apply to all commands:

- Commands begin with a keyword that is the name of the command and often have additional specifications, such as subcommands and user specifications. Refer to the discussion of each command to see which subcommands and additional specifications are required.
- Commands and any command specifications can be entered in upper and lower case. Commands, subcommands, keywords, and variable names are translated to upper case before processing. All user specifications, including variable names, labels, and data values, preserve upper and lower case.
- Spaces can be added between specifications at any point where a single blank is allowed. In addition, lines can be broken at any point where a single blank is allowed. There are two exceptions: the `END DATA` command can have only one space between words, and string specifications on commands such as `TITLE`, `SUBTITLE`, `VARIABLE LABELS`, and `VALUE`

LABELS can be broken across two lines only by specifying a plus sign (+) between string segments. For more information, see the topic [String Values in Command Specifications](#) on p. 40.

- Many command names and keywords can be abbreviated to the first three or more characters that can be resolved without ambiguity. For example, COMPUTE can be abbreviated to COMP but not COM because the latter does not adequately distinguish it from COMMENT. Some commands, however, require that all specifications be spelled out completely. This restriction is noted in the syntax chart for those commands.

## **Running Commands**

You can run commands in either batch (production) or interactive mode. In batch mode, commands are read and acted upon as a batch, so the system knows that a command is complete when it encounters a new command. In interactive mode, commands are processed immediately, and you must use a command terminator to indicate when a command is complete.

### **Interactive Mode**

The following rules apply to command specifications in interactive mode:

- Each command must start on a new line. Commands can begin in any column of a command line and continue for as many lines as needed. The exception is the END DATA command, which must begin in the first column of the first line after the end of data.
- Each command should end with a period as a command terminator. It is best to omit the terminator on BEGIN DATA, however, so that inline data are treated as one continuous specification.
- The command terminator must be the last nonblank character in a command.
- In the absence of a period as the command terminator, a blank line is interpreted as a command terminator.

*Note:* For compatibility with other modes of command execution (including command files run with INSERT or INCLUDE commands in an interactive session), each line of command syntax should not exceed 256 bytes.

### **Batch (Production) Mode**

The following rules apply to command specifications in batch mode:

- All commands in the command file must begin in column 1. You can use plus (+) or minus (–) signs in the first column if you want to indent the command specification to make the command file more readable.
- If multiple lines are used for a command, column 1 of each continuation line must be blank.
- Command terminators are optional.
- A line cannot exceed 256 bytes; any additional characters are truncated.

The following is a sample command file that will run in either interactive or batch mode:

```
GET FILE=/MYFILES/BANK.SAV'
```

```
/KEEP ID TIME SEX JOBCAT SALBEG SALNOW  
/RENAME SALNOW = SAL90.  
  
DO IF TIME LT 82.  
+ COMPUTE RATE=0.05.  
ELSE.  
+ COMPUTE RATE=0.04.  
END IF.  
  
COMPUTE SALNOW=(1+RATE)*SAL90.  
  
EXAMINE VARIABLES=SALNOW BY SEX.
```

## ***Subcommands***

Many commands include additional specifications called **subcommands**.

- Subcommands begin with a keyword that is the name of the subcommand. Most subcommands include additional specifications.
- Some subcommands are followed by an equals sign before additional specifications. The equals sign is usually optional but is required where ambiguity is possible in the specification. To avoid ambiguity, it is best to use the equals signs as shown in the syntax diagrams in this manual.
- Most subcommands can be named in any order. However, some commands require a specific subcommand order. The description of each command includes a section on subcommand order.
- Subcommands are separated from each other by a slash. To avoid ambiguity, it is best to use the slashes as shown in the syntax diagrams in this manual.

## ***Keywords***

Keywords identify commands, subcommands, functions, operators, and other specifications.

- Keywords identifying logical operators (AND, OR, and NOT); relational operators (EQ, GE, GT, LE, LT, and NE); and ALL, BY, TO, and WITH are reserved words and cannot be used as variable names.

## ***Values in Command Specifications***

The following rules apply to values specified in commands:

- A single lowercase character in the syntax diagram, such as *n*, *w*, or *d*, indicates a user-specified value.
- The value can be an integer or a real number within a restricted range, as required by the specific command or subcommand. For exact restrictions, read the individual command description.
- A number specified as an argument to a subcommand can be entered with or without leading zeros.

## String Values in Command Specifications

- Each string specified in a command should be enclosed in single or double quotes.
- To specify a single quote or apostrophe within a quoted string, either enclose the entire string in double quotes or double the single quote/apostrophe. Both of the following specifications are valid:

```
'Client's Satisfaction'
```

```
"Client's Satisfaction"
```

- To specify double quotes within a string, use single quotes to enclose the string:

```
'Categories Labeled "UNSTANDARD" in the Report'
```

- String specifications can be broken across command lines by specifying each string segment within quotes and using a plus (+) sign to join segments. For example,

```
'One, Two'
```

can be specified as

```
'One, '  
+ ' Two'
```

The plus sign can be specified on either the first or the second line of the broken string. Any blanks separating the two segments must be enclosed within one or the other string segment.

- Multiple blank spaces within quoted strings are preserved and can be significant. For example, “This string” and “This    string” are treated as different values.

## Delimiters

Delimiters are used to separate data values, keywords, arguments, and specifications.

- A blank is usually used to separate one specification from another, except when another delimiter serves the same purpose or when a comma is required.
- Commas are required to separate arguments to functions. Otherwise, blanks are generally valid substitutes for commas.
- Arithmetic operators (+, −, \*, and /) serve as delimiters in expressions.
- Blanks can be used before and after operators or equals signs to improve readability, but commas cannot.
- Special delimiters include parentheses, apostrophes, quotation marks, the slash, and the equals sign. Blanks before and after special delimiters are optional.
- The slash is used primarily to separate subcommands and lists of variables. Although slashes are sometimes optional, it is best to enter them as shown in the syntax diagrams.
- The equals sign is used between a keyword and its specifications, as in `STATISTICS=MEAN`, and to show equivalence, as in `COMPUTE target variable=expression`. Equals signs following keywords are frequently optional but are sometimes required. In general, you should follow the format of the syntax charts and examples and always include equals signs wherever they are shown.

## Command Order

Command order is more often than not a matter of common sense and follows this logical sequence: variable definition, data transformation, and statistical analysis. For example, you cannot label, transform, analyze, or use a variable in any way before it exists. The following general rules apply:

- Commands that define variables for a session (`DATA LIST`, `GET`, `GET DATA`, `MATRIX DATA`, etc.) must precede commands that assign labels or missing values to those variables; they must also precede transformation and procedure commands that use those variables.
- Transformation commands (`IF`, `COUNT`, `COMPUTE`, etc.) that are used to create and modify variables must precede commands that assign labels or missing values to those variables, and they must also precede the procedures that use those variables.
- Generally, the logical outcome of command processing determines command order. For example, a procedure that creates new variables in the active dataset must precede a procedure that uses those new variables.

In addition to observing the rules above, it is often important to distinguish between commands that cause the data to be read and those that do not, and between those that are stored pending execution with the next command that reads the data and those that take effect immediately without requiring that the data be read.

- Commands that cause the data to be read, as well as execute pending transformations, include all statistical procedures (e.g., `CROSSTABS`, `FREQUENCIES`, `REGRESSION`); some commands that save/write the contents of the active dataset (e.g., `DATASET COPY`, `SAVE TRANSLATE`, `SAVE`); `AGGREGATE`; `AUTORECODE`; `EXECUTE`; `RANK`; and `SORT CASES`.
- Commands that are stored, pending execution with the next command that reads the data, include transformation commands that modify or create new data values (e.g., `COMPUTE`, `RECODE`), commands that define conditional actions (e.g., `DO IF`, `IF`, `SELECT IF`), `PRINT`, `WRITE`, and `XSAVE`. For a comprehensive list of these commands, see [Commands That Are Stored, Pending Execution](#) on p. 44.
- Commands that take effect immediately without reading the data or executing pending commands include transformations that alter dictionary information without affecting the data values (e.g., `MISSING VALUES`, `VALUE LABELS`) and commands that don't require an active dataset (e.g., `DISPLAY`, `HOST`, `INSERT`, `OMS`, `SET`). In addition to taking effect immediately, these commands are also processed unconditionally. For example, when included within a `DO IF` structure, these commands run regardless of whether or not the condition is ever met. For a comprehensive list of these commands, see [Commands That Take Effect Immediately](#) on p. 42.

### Example

```
DO IF expense = 0.  
- COMPUTE profit=-99.  
- MISSING VALUES expense (0).  
ELSE.  
- COMPUTE profit=income-expense.  
END IF.  
LIST VARIABLES=expense profit.
```

- COMPUTE precedes MISSING VALUES and is processed first; however, execution is delayed until the data are read.
- MISSING VALUES takes effect as soon as it is encountered, *even if the condition is never met* (i.e., even if there are no cases where *expense=0*).
- LIST causes the data to be read; thus, both COMPUTE and LIST are executed during the same data pass.
- Because MISSING VALUES is already in effect by this time, the first condition in the DO IF structure will never be met, because an *expense* value of 0 is considered missing and so the condition evaluates to missing when *expense* is 0.

### **Commands That Take Effect Immediately**

These commands take effect immediately. They do not read the active dataset and do not execute pending transformations.

#### **Commands That Modify the Dictionary**

ADD DOCUMENT

ADD VALUE LABELS

APPLY DICTIONARY

DATAFILE ATTRIBUTE

DELETE VARIABLES

DOCUMENT

DROP DOCUMENTS

EXTENSION

FILE LABEL

FORMATS

MISSING VALUES

MRSETS

NUMERIC

OUTPUT EXPORT

PRINT FORMATS

RENAME VARIABLES

STRING

TMS IMPORT

TMS MERGE

VALUE LABELS

VARIABLE ALIGNMENT



VARIABLE ATTRIBUTE

VARIABLE LABELS

VARIABLE LEVEL

VARIABLE ROLE

VARIABLE WIDTH

WEIGHT

WRITE FORMATS

***Other Commands That Take Effect Immediately***

CD

CLEAR TIME PROGRAM

CLEAR TRANSFORMATIONS

CSPLAN

DATASET CLOSE

DATASET DECLARE

DATASET DISPLAY

DATASET NAME

DISPLAY

ECHO

ERASE

FILE HANDLE

FILTER

HOST

INCLUDE

INSERT

MODEL CLOSE

MODEL HANDLE

MODEL LIST

N OF CASES

NEW FILE

OMS

OMSEND

OMSINFO

OMSLOG

OUTPUT ACTIVATE  
OUTPUT CLOSE  
OUTPUT DISPLAY  
OUTPUT NAME  
OUTPUT NEW  
OUTPUT OPEN  
OUTPUT SAVE  
PERMISSIONS  
PRESERVE  
READ MODEL  
RESTORE  
SAVE MODEL  
SCRIPT  
SET  
SHOW  
SPLIT FILE  
SUBTITLE  
SYSDATA INFO  
TDISPLAY  
TITLE  
TSET  
TSHOW  
USE

***Commands That Are Stored, Pending Execution***

These commands are stored, pending execution with the next command that reads the data.

BOOTSTRAP  
BREAK  
CACHE  
COMPUTE  
COUNT  
DO IF  
DO REPEAT-END REPEAT  
IF

LEAVE  
LOOP-END LOOP  
PRINT  
PRINT EJECT  
PRINT SPACE  
RECODE  
SAMPLE  
SELECT IF  
TEMPORARY  
TIME PROGRAM  
WRITE  
XSAVE

## **Files**

IBM® SPSS® Statistics reads, creates, and writes different types of files. This section provides an overview of these types and discusses concepts and rules that apply to all files.

### **Command File**

A command file is a text file that contains syntax commands. You can type commands in a syntax window in an interactive session, use the Paste button in dialog boxes to paste generated commands into a syntax window, and/or use any text editor to create a command file. You can also edit a journal file to produce a command file. [For more information, see the topic Journal File on p. 46.](#) The following is an example of a simple command file that contains both commands and inline data:

```
DATA LIST /ID 1-3 Gender 4 (A) Age 5-6 Opinion1 TO Opinion5 7-11.  
BEGIN DATA  
001F2621221  
002M5611122  
003F3422212  
329M2121212  
END DATA.  
LIST.
```

- Case does not matter for commands but is significant for inline data. If you specified *f* for female and *m* for male in column 4 of the data line, the value of *Gender* would be *f* or *m* instead of *F* or *M* as it is now.
- Commands can be in upper or lower case. Uppercase characters are used for all commands throughout this manual only to distinguish them from other text.

## Journal File

IBM® SPSS® Statistics keeps a journal file to record all commands either run from a syntax window or generated from a dialog box during a session. You can retrieve this file with any text editor and review it to learn how the session went. You can also edit the file to build a new command file and use it in another run. An edited and tested journal file can be saved and used later for repeated tasks. The journal file also records any error or warning messages generated by commands. You can rerun these commands after making corrections and removing the messages.

The journal file is controlled by the File Locations tab of the Options dialog box, available from the Edit menu. You can turn journaling off and on, append or overwrite the journal file, and select the journal filename and location. By default, commands from subsequent sessions are appended to the journal.

The following example is a journal file for a short session with a warning message.

**Figure 2-2**

*Records from a journal file*

```
DATA LIST /ID 1-3 Gender 4 (A) Age 5-6 Opinion1 TO Opinion5 7-11.  
BEGIN DATA  
001F2621221  
002M5611122  
003F3422212  
004F45112L2  
>Warning # 1102  
>An invalid numeric field has been found. The result has been set to the  
>system-missing value.  
END DATA.  
LIST.
```

- The warning message, marked by the > symbol, tells you that an invalid numeric field has been found. Checking the last data line, you will notice that column 10 is *L*, which is probably a typographic error. You can correct the typo (for example, by changing the *L* to 1), delete the warning message, and submit the file again.

## Data Files

A wide variety of data file formats can be read and written, including raw data files created by a data entry device or a text editor, formatted data files produced by a data management program, data files generated by other software packages, and IBM® SPSS® Statistics data files.

### Raw Data Files

Raw data files contain only data, either generated by a programming language or entered with a data entry device or a text editor. Raw data arranged in almost any format can be read, including raw matrix materials and nonprintable codes. User-entered data can be embedded within a command file as inline data ([BEGIN DATA-END DATA](#)) or saved as an external file. Nonprintable machine codes are usually stored in an external file.

Commands that read raw data files include:

- [GET DATA](#)
- [DATA LIST](#)
- [MATRIX DATA](#)

Complex and hierarchical raw data files can be read using commands such as:

- [INPUT PROGRAM](#)
- [FILE TYPE](#)
- [REREAD](#)
- [REPEATING DATA](#)

### ***Data Files Created by Other Applications***

You can read files from a variety of other software applications, including:

- Excel spreadsheets ([GET DATA](#) command).
- Database tables ([GET DATA](#) command).
- IBM® SPSS® Data Collection data sources ([GET DATA](#) command).
- Delimited (including tab-delimited and CSV) and fixed-format text data files ([DATA LIST](#), [GET DATA](#)).
- dBase and Lotus files ([GET TRANSLATE](#) command).
- SAS datasets ([GET SAS](#) command).
- Stata data files ([GET STATA](#) command).

### ***IBM SPSS Statistics Data Files***

IBM® SPSS® Statistics data files are files specifically formatted for use by SPSS Statistics, containing both data and the metadata (dictionary) that define the data.

- To save the active dataset in SPSS Statistics format, use [SAVE](#) or [XSAVE](#). On most operating systems, the default extension of a saved SPSS Statistics data file is *.sav*. SPSS Statistics data files can also be matrix files created with the `MATRIX=OUT` subcommand on procedures that write matrices.
- To open SPSS Statistics data files, use [GET](#).

### ***IBM SPSS Statistics Data File Structure***

The basic structure of SPSS Statistics data files is similar to a database table:

- Rows (records) are cases. Each row represents a case or an observation. For example, each individual respondent to a questionnaire is a case.
- Columns (fields) are variables. Each column represents a variable or characteristic that is being measured. For example, each item on a questionnaire is a variable.

SPSS Statistics data files also contain metadata that describes and defines the data contained in the file. This descriptive information is called the **dictionary**. The information contained in the dictionary includes:

- Variable names and descriptive variable labels ([VARIABLE LABELS](#) command).
- Descriptive values labels ([VALUE LABELS](#) command).

- Missing values definitions ([MISSING VALUES](#) command).
- Print and write formats ([FORMATS](#) command).

Use `DISPLAY DICTIONARY` to display the dictionary for the active dataset. [For more information, see the topic DISPLAY on p. 648.](#) You can also use `SYSFILE INFO` to display dictionary information for any SPSS Statistics data file.

### **Long Variable Names**

In some instances, data files with variable names longer than eight bytes require special consideration:

- If you save a data file in portable format (see [EXPORT](#) on p. 696), variable names that exceed eight bytes are converted to unique eight-character names. For example, *mylongrootname1*, *mylongrootname2*, and *mylongrootname3* would be converted to *mylongro*, *mylong\_2*, and *mylong\_3*, respectively.
- When using data files with variable names longer than eight bytes in version 10.x or 11.x, unique, eight-byte versions of variable names are used; however, the original variable names are preserved for use in release 12.0 or later. In releases prior to 10.0, the original long variable names are lost if you save the data file.
- Matrix data files (commonly created with the `MATRIX OUT` subcommand, available in some procedures) in which the `VARNAME_` variable is longer than an eight-byte string cannot be read by releases prior to 12.0.

## **Variables**

The columns in IBM® SPSS® Statistics data files are **variables**. Variables are similar to fields in a database table.

- Variable names can be defined with numerous commands, including [DATA LIST](#), [GET DATA](#), [NUMERIC](#), [STRING](#), [VECTOR](#), [COMPUTE](#), and [RECODE](#). They can be changed with the [RENAME VARIABLES](#) command.
- Optional variable attributes can include descriptive variable labels ([VARIABLE LABELS](#) command), value labels ([VALUE LABELS](#) command), and missing value definitions ([MISSING VALUES](#) command).

The following sections provide information on variable naming rules, syntax for referring to inclusive lists of variables (keywords `ALL` and `TO`), scratch (temporary) variables, and system variables.

## Variable Names

Variable names are stored in the dictionary of the data file. Observe the following rules when establishing variable names or referring to variables by their names on commands:

- Each variable name must be unique; duplication is not allowed.
- Variable names can be up to 64 bytes long, and the first character must be a letter or one of the characters @, #, or \$. Subsequent characters can be any combination of letters, numbers, nonpunctuation characters, and a period (.). In code page mode, sixty-four bytes typically means 64 characters in single-byte languages (for example, English, French, German, Spanish, Italian, Hebrew, Russian, Greek, Arabic, and Thai) and 32 characters in double-byte languages (for example, Japanese, Chinese, and Korean). Many string characters that only take one byte in code page mode take two or more bytes in Unicode mode. For example, é is one byte in code page format but is two bytes in Unicode format; so *résumé* is six bytes in a code page file and eight bytes in Unicode mode. For information on Unicode mode, see [SET command, UNICODE subcommand](#).

*Note:* Letters include any nonpunctuation characters used in writing ordinary words in the languages supported in the platform's character set.

- Variable names cannot contain spaces.
- A # character in the first position of a variable name defines a scratch variable. You can only create scratch variables with command syntax. You cannot specify a # as the first character of a variable in dialog boxes that create new variables. [For more information, see the topic Scratch Variables on p. 51.](#)
- A \$ sign in the first position indicates that the variable is a system variable. [For more information, see the topic System Variables on p. 53.](#) The \$ sign is not allowed as the initial character of a user-defined variable.
- The period, the underscore, and the characters \$, #, and @ can be used within variable names. For example, *A\_.\$@#I* is a valid variable name.
- Variable names ending with a period should be avoided, since the period may be interpreted as a command terminator. You can only create variables that end with a period in command syntax. You cannot create variables that end with a period in dialog boxes that create new variables.
- Variable names ending in underscores should be avoided, since such names may conflict with names of variables automatically created by commands and procedures.
- Reserved keywords cannot be used as variable names. Reserved keywords are ALL, AND, BY, EQ, GE, GT, LE, LT, NE, NOT, OR, TO, and WITH.
- Variable names can be defined with any mixture of uppercase and lowercase characters, and case is preserved for display purposes.
- When long variable names need to wrap onto multiple lines in output, lines are broken at underscores, periods, and points where content changes from lower case to upper case.

### Mixed Case Variable Names

Variable names can be defined with any mixture of upper- and lowercase characters, and case is preserved for display purposes.

- Variable names are stored and displayed exactly as specified on commands that read data or create new variables. For example, `compute NewVar = 1` creates a new variable that will be displayed as *NewVar* in the Data Editor and in output from any procedures that display variable names.
- Commands that refer to existing variable names are not case sensitive. For example, `FREQUENCIES VARIABLES = newvar`, `FREQUENCIES VARIABLES = NEWVAR`, and `FREQUENCIES VARIABLES = NewVar` are all functionally equivalent.
- In languages such as Japanese, where some characters exist in both narrow and wide forms, these characters are considered different and are displayed using the form in which they were entered.
- When long variable names need to wrap onto multiple lines in output, attempts are made to break lines at underscores, periods, and changes from lower to upper case.

You can use the `RENAME VARIABLES` command to change the case of any characters in a variable name.

### **Example**

```
RENAME VARIABLES (newvariable = NewVariable).
```

- For the existing variable name specification, case is ignored. Any combination of upper and lower case will work.
- For the new variable name, case will be preserved as entered for display purposes.

For more information, see the `RENAME VARIABLES` command.

### **Long Variable Names**

In some instances, data files with variable names longer than eight bytes require special consideration:

- If you save a data file in portable format (see [EXPORT](#) on p. 696), variable names that exceed eight bytes are converted to unique eight-character names. For example, *mylongrootname1*, *mylongrootname2*, and *mylongrootname3* would be converted to *mylongro*, *mylong\_2*, and *mylong\_3*, respectively.
- When using data files with variable names longer than eight bytes in version 10.x or 11.x, unique, eight-byte versions of variable names are used; however, the original variable names are preserved for use in release 12.0 or later. In releases prior to 10.0, the original long variable names are lost if you save the data file.
- Matrix data files (commonly created with the `MATRIX OUT` subcommand, available in some procedures) in which the `VARNAME_` variable is longer than an eight-byte string cannot be read by releases prior to 12.0.

### **Keyword TO**

You can establish names for a set of variables or refer to any number of consecutive variables by specifying the beginning and the ending variables joined by the keyword `TO`.



To establish names for a set of variables with the keyword `TO`, use a character prefix with a numeric suffix.

- The prefix can be any valid name. Both the beginning and ending variables must use the same prefix.
- The numeric suffix can be any integer, but the first number must be smaller than the second. For example, `ITEM1 TO ITEM5` establishes five variables named `ITEM1`, `ITEM2`, `ITEM3`, `ITEM4`, and `ITEM5`.
- Leading zeros used in numeric suffixes are included in the variable name. For example, `V001 TO V100` establishes 100 variables—`V001`, `V002`, `V003`, ..., `V100`. `V1 TO V100` establishes 100 variables—`V1`, `V2`, `V3`, ..., `V100`.

The keyword `TO` can also be used on procedures and other commands to refer to consecutive variables on the active dataset. For example, `AVAR TO VARB` refers to the variables `AVAR` and all subsequent variables up to and including `VARB`.

- In most cases, the `TO` specification uses the variable order on the active dataset. Use the `DISPLAY` command to see the order of variables on the active dataset.
- On some subcommands, the order in which variables are named on a previous subcommand, usually the `VARIABLES` subcommand, is used to determine which variables are consecutive and therefore are implied by the `TO` specification. This is noted in the description of individual commands.

## Keyword ALL

The keyword `ALL` can be used in many commands to specify all of the variables in the active dataset. For example,

```
FREQUENCIES /VARIABLES = ALL.
```

or

```
OLAP CUBES income by ALL.
```

In the second example, a separate table will be created for every variable in the data file, including a table of `income` by `income`.

## Scratch Variables

You can use **scratch variables** to facilitate operations in transformation blocks and input programs.

- To create a scratch variable, specify a variable name that begins with the `#` character—for example, `#ID`. Scratch variables can be either numeric or string.
- Scratch variables are initialized to 0 for numeric variables or blank for string variables.
- Scratch variables cannot be used in procedures and cannot be saved in a data file (but they can be written to an external text file with `PRINT` or `WRITE`).
- Scratch variables cannot be assigned missing values, variable labels, or value labels.

- Scratch variables can be created between procedures but are always discarded as the next procedure begins.
- Scratch variables are discarded once a `TEMPORARY` command is specified.
- The keyword `TO` cannot refer to scratch variables and permanent variables at the same time.
- Scratch variables cannot be specified on a `WEIGHT` command.
- Scratch variable cannot be specified on the `LEAVE` command.
- Scratch variables are not reinitialized when a new case is read. Their values are always carried across cases. (So using a scratch variable can be essentially equivalent to using the `LEAVE` command.)

Because scratch variables are discarded, they are often useful as loop index variables and as other variables that do not need to be retained at the end of a transformation block. [For more information, see the topic Indexing Clause on p. 1059.](#) Because scratch variables are not reinitialized for each case, they are also useful in loops that span cases in an input program. [For more information, see the topic Creating Data on p. 1065.](#)

### Example

```
DATA LIST LIST (" ") /Name (A15).
BEGIN DATA
Nick Lowe
Dave Edmunds
END DATA.
STRING LastName (A15).
COMPUTE #index=INDEX(Name, " ").
COMPUTE LastName=SUBSTR(Name, #index+1).
LIST.
```

Figure 2-3

*Listing of case values*

Name	LastName
Nick Lowe	Lowe
Dave Edmunds	Edmunds

- `#index` is a scratch variable that is set to the numeric position of the first occurrence of a blank space in `Name`.
- The scratch variable is then used in the second `COMPUTE` command to determine the starting position of `LastName` within `Name`.
- The default `LIST` command will list the values of all variables for all cases. It does not include `#index` because `LIST` is a procedure that reads the data, and all scratch variables are discarded at that point.

In this example, you could have obtained the same end result without the scratch variable, using:

```
COMPUTE LastName=SUBSTR(Name, INDEX(Name, " ")+1).
```

The use of a scratch variable here simply makes the code easier to read.

**Example: Scratch variable initialization**

```

DATA LIST FREE /Var1.
BEGIN DATA
2 2 2
END DATA.
COMPUTE Var2=Var1+Var2.
COMPUTE Var3=0.
COMPUTE Var3=Var1+Var3.
COMPUTE #ScratchVar=Var1+#ScratchVar.
COMPUTE Var4=#ScratchVar.
LIST.

```

**Figure 2-4**

Listing of case values

Var1	Var2	Var3	Var4
2.00	.	2.00	2.00
2.00	.	2.00	4.00
2.00	.	2.00	6.00

- The new variable *Var2* is reinitialized to system-missing for each case, therefore *Var1+Var2* always results in system-missing.
- The new variable *Var3* is reset to 0 for each case (`COMPUTE Var3=0`), therefore *Var1+Var3* is always equivalent to *Var1+0*.
- *#ScratchVar* is initialized to 0 for the first case and is not reinitialized for subsequent cases; so *Var1+#ScratchVar* is equivalent to *Var1+0* for the first case, *Var1+2* for the second case, and *Var1+4* for the third case.
- *Var4* is set to the value of *#ScratchVar* in this example so that the value can be displayed in the case listing.

In this example, the commands:

```

COMPUTE #ScratchVar=Var1+#ScratchVar.
COMPUTE Var4=#ScratchVar.

```

are equivalent to:

```

COMPUTE Var4=Var1+Var4.
LEAVE Var4.

```

**System Variables**

**System variables** are special variables created during a working session to keep system-required information, such as the number of cases read by the system, the system-missing value, and the current date. System variables can be used in data transformations.

- The names of system variables begin with a dollar sign (\$).

- You cannot modify a system variable or alter its print or write format. Except for these restrictions, you can use system variables anywhere that a normal variable is used in the transformation language.
- System variables are not available for procedures.

<b>\$CASENUM</b>	<i>Current case sequence number.</i> For each case, <i>\$CASENUM</i> is the number of cases read up to and including that case. The format is <code>F8.0</code> . The value of <i>\$CASENUM</i> is not necessarily the row number in a Data Editor window (available in windowed environments), and the value changes if the file is sorted or new cases are inserted before the end of the file.
<b>\$SYSMIS</b>	<i>System-missing value.</i> The system-missing value displays as a period (.) or whatever is used as the decimal point.
<b>\$JDATE</b>	<i>Current date in number of days from October 14, 1582</i> (day 1 of the Gregorian calendar). The format is <code>F6.0</code> .
<b>\$DATE</b>	<i>Current date in international date format with two-digit year.</i> The format is <code>A9</code> in the form <code>dd-mmm-yy</code> .
<b>\$DATE11</b>	<i>Current date in international date format with four-digit year.</i> The format is <code>A11</code> in the form <code>dd-mmm-yyyy</code> .
<b>\$TIME</b>	<i>Current date and time.</i> <i>\$TIME</i> represents the number of seconds from midnight, October 14, 1582, to the date and time when the transformation command is executed. The format is <code>F20</code> . You can display this as a date in a number of different <a href="#">date formats</a> . You can also use it in <a href="#">date and time functions</a> .
<b>\$LENGTH</b>	<i>The current page length.</i> The format is <code>F11.0</code> . For more information, see <code>SET</code> .
<b>\$WIDTH</b>	<i>The current page width.</i> The format is <code>F3.0</code> . For more information, see <code>SET</code> .

## Variable Types and Formats

There are two basic variable types:

- **String.** Also referred to alphanumeric. String values are stored as codes listed in the IBM® SPSS® Statistics character set. [For more information, see the topic IMPORT/EXPORT Character Sets on p. 2189.](#)
- **Numeric.** Numeric values are stored internally as double-precision floating-point numbers.

Variable formats determine how raw data is read into storage and how values are displayed and written. For example, all dates and times are stored internally as numeric values, but you can use date and time format specifications to both read and display date and time values in standard date and time formats. The following sections provide details on how formats are specified and how those formats affect how data are read, displayed, and written.

## Input and Output Formats

Values are read according to their **input** format and displayed according to their **output** format. The input and output formats differ in several ways.

- The input format is either specified or implied on the `DATA LIST`, `GET DATA`, or other data definition commands. It is in effect only when cases are built in an active dataset.

- Output formats are automatically generated from input formats, with output formats expanded to include punctuation characters, such as decimal indicators, grouping symbols, and dollar signs. For example, an input format of `DOLLAR7.2` will generate an output format of `DOLLAR10.2` to accommodate the dollar sign, grouping symbol (comma), and decimal indicator (period).
- The formats (specified or default) on `NUMERIC`, `STRING`, `COMPUTE`, or other commands that create new variables are output formats. You must specify adequate widths to accommodate all punctuation characters.
- The output format is in effect during the entire working session (unless explicitly changed) and is saved in the dictionary of IBM® SPSS® Statistics data files.
- Output formats for numeric variables can be changed with `FORMATS`, `PRINT FORMATS`, and `WRITE FORMATS`.

### **String Variable Formats**

- The values of string variables can contain numbers, letters, and special characters and can be up to 32,767 bytes.
- System-missing values cannot be generated for string variables, since any character is a legal string value.
- When a transformation command that creates or modifies a string variable yields a missing or undefined result, a null string is assigned. The variable displays as blanks and is not treated as missing.
- String formats are used to read and write string variables. The input values can be alphanumeric characters (`A` format) or the hexadecimal representation of alphanumeric characters (`AHEX` format).
- For fixed-format raw data, the width can be explicitly specified on commands such as `DATA LIST` and `GET DATA` or implied if column-style specifications are used. For freefield data, the default width is 1; if the input string may be longer, `w` must be explicitly specified. Input strings shorter than the specified width are right-padded with blanks.
- The output format for a string variable is always `A`. The width is determined by the input format or the format assigned on the `STRING` command. Once defined, the width of a string variable can only be changed with the `ALTER TYPE` command.

#### ***A Format (Standard Characters)***

The `A` format is used to read standard characters. Characters can include letters, numbers, punctuation marks, blanks, and most other characters on your keyboard. Numbers entered as values for string variables cannot be used in calculations unless you convert them to numeric format with the `NUMBER` function. [For more information, see the topic String/numeric conversion functions on p. 111.](#)

*Fixed data:*

With fixed-format input data, any punctuation—including leading, trailing, and embedded blanks—within the column specifications is included in the string value. For example, a string value of

```
Mr. Ed
```

(with one embedded blank) is distinguished from a value of

```
Mr.  Ed
```

(with two embedded blanks). It is also distinguished from a string value of

```
MR. ED
```

(all upper case), and all three are treated as separate values. These can be important considerations for any procedures, transformations, or data selection commands involving string variables. Consider the following example:

```
DATA LIST FIXED /ALPHAVAR 1-10 (A) .
BEGIN DATA
Mr. Ed
Mr. Ed
MR. ED
Mr.  Ed
  Mr. Ed
END DATA.
AUTORECODE ALPHAVAR /INTO NUMVAR.
LIST.
```

AUTORECODE recodes the values into consecutive integers. The following figure shows the recoded values.

**Figure 2-5**

*Different string values illustrated*

ALPHAVAR	NUMVAR
Mr. Ed	4
Mr. Ed	4
MR. ED	2
Mr.  Ed	3
Mr. Ed	1

### **AHEX Format (Hexadecimal Characters)**

The AHEX format is used to read the hexadecimal representation of standard characters. Each set of two hexadecimal characters represents one standard character. For codes used on different operating systems, see *IMPORT/EXPORT Character Sets* on p. 2189.

- The *w* specification refers to columns of the hexadecimal representation and must be an even number. Leading, trailing, and embedded blanks are not allowed, and only valid hexadecimal characters can be used in input values.
- For some operating systems (e.g., IBM CMS), letters in hexadecimal values must be upper case.

- The default output format for variables read with the `AHEX` input format is the `A` format. The default width is half the specified input width. For example, an input format of `AHEX14` generates an output format of `A7`.
- Used as an output format, the `AHEX` format displays the printable characters in the hexadecimal characters specific to your system. The following commands run on a UNIX system—where `A=41` (decimal 65), `a=61` (decimal 97), and so on—produce the output shown below:

```
DATA LIST FIXED
  /A,B,C,D,E,F,G,H,I,J,K,L,M,N,O,P,Q,R,S,T,U,V,W,X,Y,Z 1-26 (A) .
FORMATS ALL (AHEX2) .
BEGIN DATA
ABCDEFGHIJKLMNPOQRSTUVWXYZ
abcdefghijklmnopqrstuvwxyz
END DATA.
LIST.
```

**Figure 2-6**

*Display of hexadecimal representation of the character set with `AHEX` format*

A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P	Q	R	S	T	U	V	W	X	Y	Z
41	42	43	44	45	46	47	48	49	4A	4B	4C	4D	4E	4F	50	51	52	53	54	55	56	57	58	59	5A
61	62	63	64	65	66	67	68	69	6A	6B	6C	6D	6E	6F	70	71	72	73	74	75	76	77	78	79	7A

## **Numeric Variable Formats**

- By default, if no format is explicitly specified, commands that read raw data—such as `DATA LIST` and `GET DATA`—assume that variables are numeric with an `F` format type. The default width depends on whether the data are in fixed or freefield format. For a discussion of fixed data and freefield data, see [DATA LIST](#) on p. 550.
- Numeric variables created by `COMPUTE`, `COUNT`, or other commands that create numeric variables are assigned a format type of `F8.2` (or the default numeric format defined on [SET FORMAT](#)).
- If a data value exceeds its width specification, an attempt is made to display some value nevertheless. First, the decimals are rounded, then punctuation characters are taken out, then scientific notation is tried, and if there is still not enough space, an ellipsis (...) is displayed, indicating that a value is present but cannot be displayed in the assigned width.
- The output format does not affect the value stored in the file. A numeric value is always stored in double precision.
- For all numeric formats, the maximum width is 40.
- For numeric formats where decimals are allowed, the maximum number of decimals is 16.
- For default numeric (`F`) format and scientific notation (`E`) format, the decimal indicator of the input data from text data sources (read by commands such as `DATA LIST` and `GET DATA`) must match the IBM® SPSS® Statistics locale decimal indicator (period or comma). Use [SET DECIMAL](#) to set the decimal indicator. Use [SHOW DECIMAL](#) to display the current decimal indicator.

## F, N, and E Formats

The following table lists the formats most commonly used to read in and write out numeric data. Format names are followed by total width (*w*) and an optional number of decimal positions (*d*). For example, a format of F5.2 represents a numeric value with a total width of 5, including two decimal positions and a decimal indicator.

Table 2-1  
Common numeric formats

Format type	Description	Sample format	Sample input	Output for fixed input		Output for freefield input	
				Format	Value	Format	Value
Fw.d	Standard numeric	F5.0	1234	F5.0	1234	F5.0	1234
			1.234		1*		1*
		F5.2	1234	F6.2	1234.0	F6.2	1234.0
			1.234		1.23*		1.23*
Nw.d	Restricted numeric	N5.0	00123	F5.0	123	F5.0	123
			123		. †		123
		N5.2	12345	F6.2	12345	F6.2	12345
			12.34		. †		. †
Ew.d	Scientific notation	E8.0	1234E3	E10.3	1.234E+06	E10.3	1.234E+06‡
			1234		1.234E+03		1.234E+03

\* Only the display is truncated. The value is stored in full precision.

† System-missing value.

‡ Scientific notation is accepted in input data with F, COMMA, DOLLAR, DOT, and PCT formats. The same rules apply as specified below.

### For fixed data:

- With the N format, only unsigned integers are allowed as input values. Values not padded with leading zeros to the specified width or those containing decimal points are assigned the system-missing value. This input format is useful for reading and checking values that should be integers containing leading zeros.
- The E format reads all forms of scientific notation. If the sign is omitted, + is assumed. If the sign (+ or -) is specified before the exponent, the E or D can be omitted. A single space is permitted after the E or D and/or after the sign. If both the sign and the letter E or D are omitted, implied decimal places are assumed. For example, 1.234E3, 1.234+3, 1.234E+3, 1.234D3, 1.234D+3, 1.234E 3, and 1234 are all legitimate values. Only the last value can imply decimal places.
- E format input values can be up to 40 characters wide and include up to 15 decimal positions.
- The default output width (*w*) for the E format is either the specified input width or the number of specified decimal positions plus 7 (*d*+7), whichever is greater. The minimum width is 10 and the minimum decimal places are 3.
- The DATA LIST command can read fixed-format numeric data with **implied** decimal positions. [For more information, see the topic Implied Decimal Positions on p. 565.](#)



*For freefield data:*

- F format *w* and *d* specifications do not affect how data are read. They only determine the output formats (expanded, if necessary). 1234 is always read as 1234 in freefield data, but a specified F5.2 format will be expanded to F6.2 and the value will be displayed as 1234.0 (the last decimal place is rounded because of lack of space).
- When the N format is used for freefield data, input values with embedded decimal indicators are assigned the system-missing value, but integer input values without leading zeroes are treated as valid. For example, with an input format of N5.0, a value of 123 is treated the same as a value of 00123, but a value of 12.34 is assigned the system-missing value.
- The E format for freefield data follows the same rules as for fixed data except that no blank space is permitted in the value. Thus, 1.234E3 and 1.234+3 are allowed, but the value 1.234 3 will cause mistakes when the data are read.
- The default output E format and the width and decimal place limitations are the same as with fixed data.

### ***N (Restricted Numeric) Output Format***

N format input values are assigned an F output format. To display, print, and write N format values with leading zeroes, use the `FORMATS` command to specify N as the output format. [For more information, see the topic FORMATS on p. 750.](#)

### ***COMMA, DOT, DOLLAR, and PCT Formats***

The numeric formats listed below read and write data with embedded punctuation characters and symbols, such as commas, dots, and dollar and percent signs. The input data may or may not contain such characters. The data values read in are stored as numbers but displayed using the appropriate formats.

- **DOLLAR.** Numeric values with a leading dollar sign, a comma used as the grouping separator, and a period used as the decimal indicator. For example, \$1,234.56.
- **COMMA.** Numeric values with a comma used as the grouping separator and a period used as decimal indicator. For example, 1,234.56.
- **DOT.** Numeric values with a period used as the grouping separator and a comma used as the decimal indicator. For example, 1.234,56.
- **PCT.** Numeric values with a trailing percent sign. For example, 123.45%.

The input data values may or may not contain the punctuation characters allowed by the specified format, but the data values may not contain characters *not* allowed by the format. For example, with a DOLLAR input format, input values of 1234.56, 1,234.56, and \$1,234.56 are all valid and stored internally as the same value—but with a COMMA input format, the input value with a leading dollar sign would be assigned the system-missing value.

```
DATA LIST LIST (" ") /dollarVar (DOLLAR9.2) commaVar (COMMA9.2)
    dotVar (DOT9.2) pctVar (PCT9.2).
BEGIN DATA
1234 1234 1234 1234
$1,234.00 1,234.00 1.234,00 1234.00%
END DATA.
```

LIST.

**Figure 2-7**

*Output illustrating DOLLAR, COMMA, DOT, and PCT formats*

dollarVar	commaVar	dotVar	pctVar
\$1,234.00	1,234.00	1.234,00	1234.00%
\$1,234.00	1,234.00	1.234,00	1234.00%

Other formats that use punctuation characters and symbols are date and time formats and custom currency formats. For more information on date and time formats, see [Date and Time Formats](#) on p. 63. Custom currency formats are output formats only, and are defined with the [SET](#) command.

### **Binary and Hexadecimal Formats**

Data can be read and written in formats used by a number of programming languages such as PL/I, COBOL, FORTRAN, and Assembler. The data can be binary, hexadecimal, or zoned decimal. Formats described in this section can be used both as input formats and output formats, but with fixed data only.

The default output format for all formats described in this section is an equivalent F format, allowing the maximum number of columns for values with symbols and punctuation. To change the default, use `FORMATS` or `WRITE FORMATS`.

*IBw.d (integer binary):*

The `IB` format reads fields that contain fixed-point binary (integer) data. The data might be generated by COBOL using `COMPUTATIONAL` data items, by FORTRAN using `INTEGER*2` or `INTEGER*4`, or by Assembler using fullword and halfword items. The general format is a signed binary number that is 16 or 32 bits in length.

The general syntax for the `IB` format is `IBw.d`, where `w` is the field width in bytes (omitted for column-style specifications) and `d` is the number of digits to the right of the decimal point. Since the width is expressed in bytes and the number of decimal positions is expressed in digits, `d` can be greater than `w`. For example, both of the following commands are valid:

```
DATA LIST FIXED /VAR1 (IB4.8).
```

```
DATA LIST FIXED /VAR1 1-4 (IB,8).
```

Widths of 2 and 4 represent standard 16-bit and 32-bit integers, respectively. Fields read with the `IB` format are treated as signed. For example, the one-byte binary value 11111111 would be read as -1.

*PIBw.d (positive integer binary):*

The `PIB` format is essentially the same as `IB` except that negative numbers are not allowed. This restriction allows one additional bit of magnitude. The same one-byte value 11111111 would be read as 255.

*PIBHEX<sub>w</sub> (hexadecimal of PIB):*

The PIBHEX format reads hexadecimal numbers as unsigned integers and writes positive integers as hexadecimal numbers. The general syntax for the PIBHEX format is PIBHEX<sub>w</sub>, where *w* indicates the total number of hexadecimal characters. The *w* specification must be an even number with a maximum of 16.

For input data, each hexadecimal number must consist of the exact number of characters. No signs, decimal points, or leading and trailing blanks are allowed. For some operating systems (such as IBM CMS), hexadecimal characters must be upper case. The following example illustrates the kind of data that the PIBHEX format can read:

```
DATA LIST FIXED
  /VAR1 1-4 (PIBHEX) VAR2 6-9 (PIBHEX) VAR3 11-14 (PIBHEX) .
BEGIN DATA
0001 0002 0003
0004 0005 0006
0007 0008 0009
000A 000B 000C
000D 000E 000F
00F0 0B2C FFFF
END DATA.
LIST.
```

The values for *VAR1*, *VAR2*, and *VAR3* are listed in the figure below. The PIBHEX format can also be used to write decimal values as hexadecimal numbers, which may be useful for programmers.

**Figure 2-8**

*Output displaying values read in PIBHEX format*

VAR1	VAR2	VAR3
1	2	3
4	5	6
7	8	9
10	11	12
13	14	15
240	2860	65535

*Zw.d (zoned decimal):*

The z format reads data values that contain zoned decimal data. Such numbers may be generated by COBOL systems using DISPLAY data items, by PL/I systems using PICTURE data items, or by Assembler using zoned decimal data items.

In zoned decimal format, one digit is represented by one byte, generally hexadecimal F1 representing 1, F2 representing 2, and so on. The last byte, however, combines the sign for the number with the last digit. In the last byte, hexadecimal A, F, or C assigns +, and B, D, or E assigns -. For example, hexadecimal D1 represents 1 for the last digit and assigns the minus sign (-) to the number.

The general syntax of the z format is Zw.d, where *w* is the total number of bytes (which is the same as columns) and *d* is the number of decimals. For input data, values can appear anywhere within the column specifications. Both leading and trailing blanks are allowed. Decimals can be implied by the input format specification or explicitly coded in the data. Explicitly coded decimals override the input format specifications.

The following example illustrates how the Z format reads zoned decimals in their printed forms on IBM mainframe and PC systems. The printed form for the sign zone (A to I for +1 to +9, and so on) may vary from system to system.

```
DATA LIST FIXED /VAR1 1-5 (Z) VAR2 7-11 (Z,2) VAR3 13-17 (Z)
  VAR4 19-23 (Z,2) VAR5 25-29 (Z) VAR6 31-35 (Z,2).
BEGIN DATA
1234A 1234A 1234B 1234B 1234C 1234C
1234D 1234D 1234E 1234E 1234F 1234F
1234G 1234G 1234H 1234H 1234I 1234I
1234J 1234J 1234K 1234K 1234L 1234L
1234M 1234M 1234N 1234N 1234O 1234O
1234P 1234P 1234Q 1234Q 1234R 1234R
1234{ 1234{ 1234} 1234} 1.23M 1.23M
END DATA.
LIST.
```

The values for *VAR1* to *VAR6* are listed in the following figure.

**Figure 2-9**

*Output displaying values read in Z format*

VAR1	VAR2	VAR3	VAR4	VAR5	VAR6
12341	123.41	12342	123.42	12343	123.43
12344	123.44	12345	123.45	12346	123.46
12347	123.47	12348	123.48	12349	123.49
-12341	-123.41	-12342	-123.42	-12343	-123.43
-12344	-123.44	-12345	-123.45	-12346	-123.46
-12347	-123.47	-12348	-123.48	-12349	-123.49
12340	123.40	-12340	-123.40	-1	-1.23

The default output format for the Z format is the equivalent F format, as shown in the figure. The default output width is based on the input width specification plus one column for the sign and one column for the implied decimal point (if specified). For example, an input format of Z4.0 generates an output format of F5.0, and an input format of Z4.2 generates an output format of F6.2.

*Pw.d (packed decimal):*

The P format is used to read fields with packed decimal numbers. Such numbers are generated by COBOL using COMPUTATIONAL-3 data items and by Assembler using packed decimal data items. The general format of a packed decimal field is two four-bit digits in each byte of the field except the last. The last byte contains a single digit in its four leftmost bits and a four-bit sign in its rightmost bits. If the last four bits are 1111 (hexadecimal F), the value is positive; if they are 1101 (hexadecimal D), the value is negative. One byte under the P format can represent numbers from -9 to 9.

The general syntax of the P format is Pw.d, where w is the number of bytes (not digits) and d is the number of digits to the right of the implied decimal point. The number of digits in a field is  $(2^w - 1)$ .

*PKw.d (unsigned packed decimal):*

The PK format is essentially the same as P except that there is no sign. That is, even the rightmost byte contains two digits, and negative data cannot be represented. One byte under the PK format can represent numbers from 0 to 99. The number of digits in a field is  $2^w$ .

*RBw (real binary):*

The RB format is used to read data values that contain internal format floating-point numbers. Such numbers are generated by COBOL using COMPUTATIONAL-1 or COMPUTATIONAL-2 data items, by PL/I using FLOATING DECIMAL data items, by FORTRAN using REAL or REAL\*8 data items, or by Assembler using floating-point data items.

The general syntax of the RB format is RBw, where w is the total number of bytes. The width specification must be an even number between 2 and 8. Normally, a width specification of 8 is used to read double-precision values, and a width of 4 is used to read single-precision values.

*RBHEXw (hexadecimal of RB):*

The RBHEX format interprets a series of hexadecimal characters as a number that represents a floating-point number. This representation is system-specific. If the field width is less than twice the width of a floating-point number, the value is right-padded with binary zeros. For some operating systems (for example, IBM CMS), letters in hexadecimal values must be upper case.

The general syntax of the RBHEX format is RBHEXw, where w indicates the total number of columns. The width must be an even number. The values are real (floating-point) numbers. Leading and trailing blanks are not allowed. Any data values shorter than the specified input width must be padded with leading zeros.

## **Date and Time Formats**

Date and time formats are both input and output formats. Like numeric formats, each input format generates a default output format, automatically expanded (if necessary) to accommodate display width. Internally, all date and time format values are stored as a number of seconds: date formats (e.g., DATE, ADATE, SDATE, DATETIME) are stored as the number of seconds since October 14, 1582; time formats (TIME, DTIME) are stored as a number of seconds that represents a time interval (e.g., 10:00:00 is stored internally as 36000, which is 60 seconds x 60 minutes x 10 hours).

- All date and time formats have a minimum input width, and some have a different minimum output. Wherever the input minimum width is less than the output minimum, the width is expanded automatically when displaying or printing values. However, when you specify output formats, you must allow enough space for displaying the date and time in the format you choose.
- Input data shorter than the specified width are correctly evaluated as long as all the necessary elements are present. For example, with the TIME format, 1:2, 01 2, and 01:02 are all correctly evaluated even though the minimum width is 5. However, if only one element (hours or minutes) is present, you must use a time function to aggregate or convert the data. [For more information, see the topic Date and time functions on p. 98.](#)
- If a date or time value cannot be completely displayed in the specified width, values are truncated in the output. For example, an input time value of 1:20:59 (1 hour, 20 minutes, 59 seconds) displayed with a width of 5 will generate an output value of 01:20, not 01:21. The truncation of output does not affect the numeric value stored in the working file.

The following table shows all available date and time formats, where w indicates the total number of columns and d (if present) indicates the number of decimal places for fractional seconds. The example shows the output format with the minimum width and default decimal positions (if

applicable). The format allowed in the input data is much less restrictive. [For more information, see the topic Input Data Specification on p. 65.](#)

Table 2-2  
Date and time formats

Format type	Description	Min w		Max w	Max d	General form	Example
		In	Out				
DATEw	International date	9	9	40		dd-mmm-yy	28-OCT-90
		10	11			dd-mmm-yyyy	28-OCT-1990
ADATEw	American date	8	8	40		mm/dd/yy	10/28/90
		10	10			mm/dd/yyyy	10/28/1990
EDATEw	European date	8	8	40		dd.mm.yy	28.10.90
		10	10			dd.mm.yyyy	28.10.1990
JDATEw	Julian date	5	5	40		yyddd	90301
		7	7			yyyyddd	1990301
SDATEw	Sortable date*	8	8	40		yy/mm/dd	90/10/28
		10	10			yyyy/mm/dd	1990/10/28
QYRw	Quarter and year	4	6	40		q Q yy	4 Q 90
		6	8			q Q yyyy	4 Q 1990
MOYRw	Month and year	6	6	40		mmm yy	OCT 90
		8	8			mmm yyyy	OCT 1990
WKYRw	Week and year	6	8	40		ww WK yy	43 WK 90
		8	10			ww WK yyyy	43 WK 1990
WKDAYw	Day of the week	2	2	40		(name of the day)	SU
MONTHw	Month	3	3	40		(name of the month)	JAN
TIMEw	Time	5	5	40		hh:mm	01:02
TIMEw.d		10	10	40	16	hh:mm:ss.s	01:02:34.75
DTIMEw	Days and time	1	1	40		dd hh:mm	20 08:03
DTIMEw.d		13	13	40	16	dd hh:mm:ss.s	20 08:03:00
DATETIMEw	Date and time	17	17	40		dd-mmm-yyyy hh:mm	20-JUN-1990 08:03
DATETIMEw.d		22	22	40	16	dd-mmm-yyyy hh:mm:ss.s	20-JUN-1990 08:03:00

\* All date and time formats produce sortable data. *SDATE*, a date format used in a number of Asian countries, can be sorted in its character form and is used as a sortable format by many programmers.

## ***Input Data Specification***

The following general rules apply to date and time input formats:

- The century value for two-digit years is defined by the *SET EPOCH* value. By default, the century range begins 69 years prior to the current year and ends 30 years after the current year. Whether all four digits or only two digits are displayed in output depends on the width specification on the format.

- Dashes, periods, commas, slashes, or blanks can be used as delimiters in the input values. For example, with the *DATE* format, the following input forms are all acceptable:

28-OCT-90 28/10/1990 28.OCT.90 28 October, 1990

The displayed values, however, will be the same: *28-OCT-90* or *28-OCT-1990*, depending on whether the specified width allows 11 characters in output.

- The *JDATE* format does not allow internal delimiters and requires leading zeros for day values of less than 100 and two-digit-year values of less than 10. For example, for January 1, 1990, the following two specifications are acceptable:

90001 1990001

However, neither of the following is acceptable:

90 1 90/1

- Months can be represented in digits, Roman numerals, or three-character abbreviations, and they can be fully spelled out. For example, all of the following specifications are acceptable for October:

10 X OCT October

- The quarter in *QYR* format is expressed as 1, 2, 3, or 4. It must be separated from the year by the letter *Q*. Blanks can be used as additional delimiters. For example, for the fourth quarter of 1990, all of the following specifications are acceptable:

4Q90 4Q1990 4 Q 90 4 Q 1990

On some operating systems, such as IBM CMS, *Q* must be upper case. The displayed output is *4 Q 90* or *4 Q 1990*, depending on whether the width specified allows all four digits of the year.

- The week in the *WKYR* format is expressed as a number from 1 to 53. Week 1 begins on January 1, week 2 on January 8, and so on. The value may be different from the number of the calendar week. The week and year must be separated by the string *WK*. Blanks can be used as additional delimiters. For example, for the 43rd week of 1990, all of the following specifications are acceptable:

43WK90 43WK1990 43 WK 90 43 WK 1990

On some operating systems, such as IBM CMS, *WK* must be upper case. The displayed output is *43 WK 90* or *43 WK 1990*, depending on whether the specified width allows enough space for all four digits of the year.

- In time specifications, colons can be used as delimiters between hours, minutes, and seconds. Hours and minutes are required, but seconds are optional. A period is required to separate seconds from fractional seconds. Hours can be of unlimited magnitude, but the maximum value for minutes is 59 and for seconds 59.999. . . .

- Data values can contain a sign (+ or -) in TIME and DTIME formats to represent time intervals before or after a point in time.

**Example: DATE, ADATE, and JDATE**

```
DATA LIST FIXED
  /VAR1 1-17 (DATE) VAR2 21-37 (ADATE) VAR3 41-47 (JDATE) .
BEGIN DATA
28-10-90          10/28/90          90301
28.OCT.1990      X 28 1990        1990301
28 October, 2001 Oct. 28, 2001    2001301
END DATA.
LIST.
```

- Internally, all date format variables are stored as the number of seconds from 0 hours, 0 minutes, and 0 seconds of Oct. 14, 1582.

The LIST output from these commands is shown in the following figure.

**Figure 2-10**  
Output illustrating DATE, ADATE, and JDATE formats

VAR1	VAR2	VAR3
28-OCT-1990	10/28/1990	1990301
28-OCT-1990	10/28/1990	1990301
28-OCT-2001	10/28/2001	2001301

**Example: QYR, MOYR, and WKYR**

```
DATA LIST FIXED /VAR1 1-10 (QYR) VAR2 12-25 (MOYR) VAR3 28-37 (WKYR) .
BEGIN DATA
4Q90             10/90             43WK90
4 Q 90          Oct-1990         43 WK 1990
4 Q 2001        October, 2001    43 WK 2001
END DATA.
LIST.
```

- Internally, the value of a QYR variable is stored as midnight of the first day of the first month of the specified quarter, the value of a MOYR variable is stored as midnight of the first day of the specified month, and the value of a WKYR format variable is stored as midnight of the first day of the specified week. Thus, *4Q90* and *10/90* are both equivalent to October 1, 1990, and *43WK90* is equivalent to October 22, 1990.

The LIST output from these commands is shown in the following figure.

**Figure 2-11**  
Output illustrating QYR, MOYR, and WKYR formats

VAR1	VAR2	VAR3
4 Q 1990	OCT 1990	43 WK 1990
4 Q 1990	OCT 1990	43 WK 1990
4 Q 2001	OCT 2001	43 WK 2001

**Example: TIME**

```
DATA LIST FIXED
  /VAR1 1-11 (TIME,2) VAR2 13-21 (TIME) VAR3 23-28 (TIME) .
BEGIN DATA
```



```
1:2:34.75  1:2:34.75  1:2:34
END DATA.
LIST.
```

- TIME reads and writes time of the day or a time interval.
- Internally, the TIME values are stored as the number of seconds from midnight of the day or of the time interval.

The LIST output from these commands is shown in the following figure.

**Figure 2-12**

*Output illustrating TIME format*

```
VAR1      VAR2      VAR3
1:02:34.75  1:02:34  1:02
```

**Example: WKDAY and MONTH**

```
DATA LIST FIXED
  /VAR1 1-9 (WKDAY) VAR2 10-18 (WKDAY)
  VAR3 20-29 (MONTH) VAR4 30-32 (MONTH) VAR5 35-37 (MONTH).
BEGIN DATA
Sunday  Sunday  January  1      Jan
Monday  Monday  February 2      Feb
Tues    Tues    March     3      Mar
Wed     Wed     April     4      Apr
Th      Th      Oct       10     Oct
Fr      Fr      Nov       11     Nov
Sa      Sa      Dec       12     Dec
END DATA.
FORMATS VAR2 VAR5 (F2).
LIST.
```

- WKDAY reads and writes the day of the week; MONTH reads and writes the month of the year.
- Values for WKDAY are entered as strings but stored as numbers. They can be used in arithmetic operations but not in string functions.
- Values for MONTH can be entered either as strings or as numbers but are stored as numbers. They can be used in arithmetic operations but not in string functions.
- To display the values as numbers, assign an F format to the variable, as was done for VAR2 and VAR5 in the above example.

The LIST output from these commands is shown in the following figure.

**Figure 2-13**

*Output illustrating WKDAY and MONTH formats*

```
VAR1  VAR2      VAR3  VAR4  VAR5
SUNDAY  1  JANUARY  JAN  1
MONDAY  2  FEBRUARY FEB  2
TUESDAY 3  MARCH    MAR  3
WEDNESDAY 4  APRIL    APR  4
THURSDAY 5  OCTOBER  OCT 10
FRIDAY  6  NOVEMBER NOV 11
SATURDAY 7  DECEMBER DEC 12
```

**Example: DTIME and DATETIME**

```
DATA LIST FIXED /VAR1 1-14 (DTIME) VAR2 18-42 (DATETIME) .
BEGIN DATA
20 8:3                20-6-90 8:3
20:8:03:46          20/JUN/1990 8:03:46
20 08 03 46.75      20 June, 2001 08 03 46.75
END DATA.
LIST.
```

- DTIME and DATETIME read and write time intervals.
- The decimal point explicitly coded in the input data for fractional seconds.
- The DTIME format allows a – or + sign in the data value to indicate a time interval before or after a point in time.
- Internally, values for a DTIME variable are stored as the number of seconds of the time interval, while those for a DATETIME variable are stored as the number of seconds from 0 hours, 0 minutes, and 0 seconds of Oct. 14, 1582.

The LIST output from these commands is shown in the following figure.

Figure 2-14  
Output illustrating DTIME and DATETIME formats

VAR1	VAR2
20 08:03:00	20-JUN-1990 08:03:00
20 08:03:46	20-JUN-1990 08:03:46
20 08:03:46	20-JUN-2001 08:03:46

**FORTRAN-like Input Format Specifications**

You can use FORTRAN-like input format specifications to define formats for a set of variables, as in the following example:

```
DATA LIST FILE=HUBDATA RECORDS=3
      /MOHIRED, YRHIRED, DEPT1 TO DEPT4 (T12, 2F2.0, 4(1X,F1.0)).
```

- The specification T12 in parentheses tabs to the 12th column. The first variable (*MOHIRED*) will be read beginning from column 12.
- The specification 2F2.0 assigns the format F2.0 to two adjacent variables (*MOHIRED* and *YRHIRED*).
- The next four variables (*DEPT1* to *DEPT4*) are each assigned the format F1.0. The 4 in 4(1X,F1.0) distributes the same format to four consecutive variables. 1X skips one column before each variable. (The column-skipping specification placed within the parentheses is distributed to each variable.)

**Transformation Expressions**

Transformation expressions are used in commands such as COMPUTE, IF, DO IF, LOOP IF, and SELECT IF.

### **Release history**

#### Release 13.0

- `APPLYMODEL` and `STRAPPLYMODEL` functions introduced.
- `DATEDIFF` and `DATESUM` functions introduced.

#### Release 14.0

- `REPLACE` function introduced.
- `VALUELABEL` function introduced.

#### Release 16.0

- `CHAR.INDEX` function introduced.
- `CHAR.LENGTH` function introduced.
- `CHAR.LPAD` function introduced.
- `CHAR.MBLEN` function introduced.
- `CHAR.RINDEX` function introduced.
- `CHAR.RPAD` function introduced.
- `CHAR.SUBSTR` function introduced.
- `NORMALIZE` function introduced.
- `NTRIM` function introduced.
- `STRUNC` function introduced.

#### Release 17.0

- `MEDIAN` function introduced.
- *mult* and *fuzzbits* arguments introduced for the `RND` and `TRUNC` functions.
- `NEIGHBOR` and `DISTANCE` functions added to `APPLYMODEL` and `STRAPPLYMODEL`.

## **Numeric expressions**

Numeric expressions can be used with the `COMPUTE` and `IF` commands and as part of a logical expression for commands such as `IF`, `DO IF`, `LOOP IF`, and `SELECT IF`. Arithmetic expressions can also appear in the index portion of a `LOOP` command, on the `REPEATING DATA` command, and on the `PRINT SPACES` command.

New numeric variables created with transformation expressions have an unknown measurement level until after the next command that reads the data (such as a statistical or charting procedure or the `EXECUTE` command). For information on default measurement level assignment, see [SET SCALEMIN](#).

### **Arithmetic operations**

The following arithmetic operators are available:

+	<i>Addition</i>
-	<i>Subtraction</i>
*	<i>Multiplication</i>
/	<i>Division</i>
**	<i>Exponentiation</i>

- No two operators can appear consecutively.
- Arithmetic operators cannot be implied. For example, (VAR1) (VAR2) is not a legal specification; you must specify VAR1\*VAR2.
- Arithmetic operators and parentheses serve as delimiters. To improve readability, blanks (not commas) can be inserted before and after an operator.
- To form complex expressions, you can use variables, constants, and functions with arithmetic operators.
- The order of execution is as follows: functions; exponentiation; multiplication, division, and unary -; and addition and subtraction.
- Operators at the same level are executed from left to right.
- To override the order of operation, use parentheses. Execution begins with the innermost set of parentheses and progresses out.

### **Numeric constants**

- Constants used in numeric expressions or as arguments to functions can be integer or noninteger, depending on the application or function.
- You can specify as many digits in a constant as needed as long as you understand the precision restrictions of your computer.
- Numeric constants can be signed (+ or -) but cannot contain any other special characters, such as the comma or dollar sign.
- Numeric constants can be expressed with scientific notation. Thus, the exponent for a constant in scientific notation is limited to two digits. The range of values allowed for exponents in scientific notation is from -99 to +99.

### **Complex numeric arguments**

- Except where explicitly restricted, complex expressions can be formed by nesting functions and arithmetic operators as arguments to functions.
- The order of execution for complex numeric arguments is as follows: functions; exponentiation; multiplication, division, and unary -; and addition and subtraction.
- To control the order of execution in complex numeric arguments, use parentheses.

### **Arithmetic operations with date and time variables**

Most date and time variables are stored internally as the number of seconds from a particular date or as a time interval and therefore can be used in arithmetic operations. Many operations involving dates and time can be accomplished with the extensive collection of [date and time functions](#).

- A *date* is a floating-point number representing the number of seconds from midnight, October 14, 1582. Dates, which represent a particular point in time, are stored as the number of seconds to that date. For example, October 28, 2007, is stored as 13,412,908,800.
- A date includes the time of day, which is the time interval past midnight. When time of day is not given, it is taken as 00:00 and the date is an even multiple of 86,400 (the number of seconds in a day).
- A *time interval* is a floating-point number representing the number of seconds in a time period, for example, an hour, minute, or day. For example, the value representing 5.5 days is 475,200; the value representing the time interval 14:08:17 is 50,897.
- QYR, MOYR, and WKYR variables are stored as midnight of the first day of the respective quarter, month, and week of the year. Therefore, *1 Q 90*, *1/90*, and *1 WK 90* are all equivalents of January 1, 1990, 0:0:00.
- WKDAY variables are stored as 1 to 7 and MONTH variables as 1 to 12.

You can perform virtually any arithmetic operation with both date format and time format variables. Of course, not all of these operations are particularly useful. You can calculate the number of days between two dates by subtracting one date from the other—but adding two dates does not produce a very meaningful result.

By default, any new numeric variables that you compute are displayed in F format. In the case of calculations involving time and date variables, this means that the default output is expressed as a number of seconds. Use the `FORMATS` (or `PRINT FORMATS`) command to specify an appropriate format for the computed variable.

#### **Example**

```
DATA LIST FREE /Date1 Date2 (2ADATE10).
BEGIN DATA
6/20/2006 10/28/2006
END DATA.
COMPUTE DateDiff1=(Date2-Date1)/60/60/24.
COMPUTE DateDiff2=DATEDIFF(Date2,Date1, "days").
COMPUTE FutureDate1=Date2+(10*60*60*24).
COMPUTE FutureDate2=DATESUM(Date2, 10, "days").
FORMATS FutureDate1 FutureDate2 (ADATE10).
```

- The first two `COMPUTE` commands both calculate the number of days between two dates. In the first one, `Date2-Date1` yields the number of seconds between the two dates, which is then converted to the number of days by dividing by number of seconds in a minute, number of minutes in an hour, and number of hours in a day. In the second one, the `DATEDIFF` function is used to obtain the equivalent result, but instead of an arithmetic formula to produce a result expressed in days, it simply includes the argument `"days"`.

- The second pair of COMPUTE commands both calculate a date 10 days from *Date2*. In the first one, 10 days needs to be converted to the number of seconds in ten days before it can be added to *Date2*. In the second one, the "days" argument in the DATESUM function handles that conversion.
- The FORMATS command is used to display the results of the second two COMPUTE commands as dates, since the default format is F, which would display the results as the number of seconds since October 14, 1582.

For more information on date and time functions, see [Date and time functions](#) on p. 98.

### **Conditional statements and case selection based on dates**

To specify a date as a value in a conditional statement, use one of the data aggregation functions to express the date value. For example,

```
***this works***.
SELECT IF datevar >= date.mdy(3,1,2006).
***the following do not work***.
SELECT IF datevar >= 3/1/2006. /*this will select dates >= 0.0015.
SELECT IF datevar >= "3/1/2006" /*this will generate an error.
```

For more information, see the topic [Aggregation functions](#) on p. 98.

### **Domain errors**

Domain errors occur when numeric expressions are mathematically undefined or cannot be represented numerically on the computer for reasons other than missing data. Two common examples are division by 0 and the square root of a negative number. When there is a domain error, a warning is issued, and the system-missing value is assigned to the expression. For example, the command COMPUTE TESTVAR = TRUNC(SQRT(X/Y) \* .5) returns system-missing if X/Y is negative or if Y is 0.

The following are domain errors in numeric expressions:

**	<i>A negative number to a noninteger power.</i>
/	<i>A divisor of 0.</i>
MOD	<i>A divisor of 0.</i>
SQRT	<i>A negative argument.</i>
EXP	<i>An argument that produces a result too large to be represented on the computer.</i>
LG10	<i>A negative or 0 argument.</i>
LN	<i>A negative or 0 argument.</i>
ARSIN	<i>An argument whose absolute value exceeds 1.</i>
NORMAL	<i>A negative or 0 argument.</i>
PROBIT	<i>A negative or 0 argument, or an argument 1 or greater.</i>

## Numeric functions

Numeric functions can be used in any numeric expression on `IF`, `SELECT IF`, `DO IF`, `ELSE IF`, `LOOP IF`, `END LOOP IF`, and `COMPUTE` commands. Numeric functions always return numbers (or the system-missing value whenever the result is indeterminate). The expression to be transformed by a function is called the *argument*. Most functions have a variable or a list of variables as arguments.

- In numeric functions with two or more arguments, each argument must be separated by a comma. Blanks alone cannot be used to separate variable names, expressions, or constants in transformation expressions.
- Arguments should be enclosed in parentheses, as in `TRUNC(INCOME)`, where the `TRUNC` function returns the integer portion of the variable *INCOME*.
- Multiple arguments should be separated by commas, as in `MEAN(Q1, Q2, Q3)`, where the `MEAN` function returns the mean of variables *Q1*, *Q2*, and *Q3*.

### Example

```
COMPUTE Square_Root = Sqrt(var4).
COMPUTE Remainder = MOD(var4, 3).
COMPUTE Average = MEAN.3(var1, var2, var3, var4).
COMPUTE Trunc_Mean = TRUNC(MEAN(var1 TO var4)).
```

- `Sqrt(var4)` returns the square root of the value of *var4* for each case.
- `MOD(var4, 3)` returns the remainder (modulus) from dividing the value of *var4* by 3.
- `MEAN.3(var1, var2, var3, var4)` returns the mean of the four specified variables, provided that at least three of them have nonmissing values. The divisor for the calculation of the mean is the number of nonmissing values.
- `TRUNC(MEAN(var1 TO var4))` computes the mean of the values for the inclusive range of variables and then truncates the result. Since no minimum number of nonmissing values is specified for the function, a mean will be calculated (and truncated) as long as at least one of the variables has a nonmissing value for that case.

## Arithmetic functions

- All arithmetic functions except `MOD`, `RND` and `TRUNC` have single arguments; `MOD` has two while `RND` and `TRUNC` have from one to three. Multiple arguments must be separated by a comma.
- Arguments can be numeric expressions, as in `RND(A**2/B)`.

**ABS.** `ABS(numexpr)`. Numeric. Returns the absolute value of `numexpr`, which must be numeric.

**RND.** `RND(numexpr[,mult,fuzzbits])`. Numeric. With a single argument, returns the integer nearest to that argument. Numbers ending in .5 exactly are rounded away from 0. For example, `RND(-4.5)` rounds to -5. The optional second argument, `mult`, specifies that the result is an integer multiple of this value—for example, `RND(-4.57,0.1) = -4.6`. The value must be numeric but cannot be 0. The default is 1.

The optional third argument, `fuzzbits`, is the number of least-significant bits by which the internal representation of `numexpr` (expressed as a 64-bit floating point binary) may fall short of the threshold for rounding up (e.g., 0.5 when rounding to an integer) but still be rounded up. For example, the sum  $9.62 - 5.82 - 9.21 + 6.91$  has an internal representation of 1.499999999999998 (on an Intel processor). With `fuzzbits` set to 0 and `mult` set to 1, this expression will round to 1.0, although the exact sum is 1.50 which would round to 2.0. Allowing the rounding threshold to have a small fuzziness compensates for the minute differences between calculations with floating point numbers and exact results. In this case, adding a fuzziness of 4 bits is sufficient to produce the expected result of 2.0.

If the argument `fuzzbits` is omitted, the value specified by `SET FUZZBITS` is used. The installed setting of `FUZZBITS` is 6, which should be sufficient for most applications. Setting `fuzzbits` to 0 produces the same results as in release 10. Setting `fuzzbits` to 10 produces the same results as in releases 11 and 12.

To produce the same results as in release 13, use the following expression in place of the `RND` function:

```
TRUNC(numexpr, 1, 0) + ((.5+TRUNC(numexpr, 1, 0)-numexpr) <max(1e-13, min(.5, numexpr*1e-13)) )
```

To produce the same results as in releases 14, 15, and 16 use:

```
RND(numexpr, 1, 12.5-ln(max(1e-50, abs(numexpr))) / ln(2))
```

**TRUNC.** `TRUNC(numexpr[,mult,fuzzbits])`. Numeric. Returns the value of `numexpr` truncated toward 0. The optional second argument, `mult`, specifies that the result is an integer multiple of this value—for example, `TRUNC(4.579,0.1) = 4.5`. The value must be numeric but cannot be 0. The default is 1.

The optional third argument, `fuzzbits`, is the number of least-significant bits by which the internal representation of `numexpr` (expressed as a 64-bit floating point binary) may fall short of the nearest rounding boundary and be rounded up before truncating. For example, the sum  $9.62 - 5.82 - 9.21 + 6.91$  has an internal representation of 1.499999999999998 (on an Intel processor). With `fuzzbits` set to 0 and `mult` set to 0.1, this expression will truncate to 1.4, although the exact sum is 1.50 which would truncate to 1.5. Adding a small fuzziness to the nearest rounding boundary (in this case, 1.5) compensates for the minute differences between calculations with floating point numbers and exact results. In this case, adding a fuzziness of 5 bits is sufficient to produce the expected result of 1.5.

If the argument `fuzzbits` is omitted, the value specified by `SET FUZZBITS` is used. The installed setting of `FUZZBITS` is 6, which should be sufficient for most applications. Setting `fuzzbits` to 0 produces the same results as in release 10. Setting `fuzzbits` to 10 produces the same results as in releases 11 and 12.

To produce the same results as in release 13 use:

```
TRUNC(numexpr, 1, 0) + (TRUNC(numexpr, 1, 0) + 1 - numexpr <= 1e-13)
```

To produce the same results as in releases 14, 15, and 16 use:

```
TRUNC(numexpr, 1, 12.5-ln(max(1e-50, abs(numexpr))) / ln(2))
```



**MOD.** MOD(numexpr,modulus). Numeric. Returns the remainder when numexpr is divided by modulus. Both arguments must be numeric, and modulus must not be 0.

**SQRT.** SQRT(numexpr). Numeric. Returns the positive square root of numexpr, which must be numeric and not negative.

**EXP.** EXP(numexpr). Numeric. Returns e raised to the power numexpr, where e is the base of the natural logarithms and numexpr is numeric. Large values of numexpr may produce results that exceed the capacity of the machine.

**LG10.** LG10(numexpr). Numeric. Returns the base-10 logarithm of numexpr, which must be numeric and greater than 0.

**LN.** LN(numexpr). Numeric. Returns the base-e logarithm of numexpr, which must be numeric and greater than 0.

**LNGAMMA.** LNGAMMA(numexpr). Numeric. Returns the logarithm of the complete Gamma function of numexpr, which must be numeric and greater than 0.

**ARSIN.** ARSIN(numexpr). Numeric. Returns the inverse sine (arcsine), in radians, of numexpr, which must evaluate to a numeric value between -1 and +1.

**ARTAN.** ARTAN(numexpr). Numeric. Returns the inverse tangent (arctangent), in radians, of numexpr, which must be numeric.

**SIN.** SIN(radians). Numeric. Returns the sine of radians, which must be a numeric value, measured in radians.

**COS.** COS(radians). Numeric. Returns the cosine of radians, which must be a numeric value, measured in radians.

## **Statistical functions**

- Each argument to a statistical function (expression, variable name, or constant) must be separated by a comma.
- The *.n* suffix can be used with all statistical functions to specify the number of valid arguments. For example, MEAN . 2 ( A , B , C , D ) returns the mean of the valid values for variables *A*, *B*, *C*, and *D* only if at least two of the variables have valid values. The default for *n* is 2 for SD, VARIANCE, and CFVAR and 1 for other statistical functions. If the number specified exceeds the number of arguments in the function, the result is system-missing.
- The keyword TO can be used to refer to a set of variables in the argument list.

**SUM.** SUM(numexpr,numexpr[,..]). Numeric. Returns the sum of its arguments that have valid, nonmissing values. This function requires two or more arguments, which must be numeric. You can specify a minimum number of valid arguments for this function to be evaluated.

**MEAN.** MEAN(numexpr,numexpr[,..]). Numeric. Returns the arithmetic mean of its arguments that have valid, nonmissing values. This function requires two or more arguments, which must be numeric. You can specify a minimum number of valid arguments for this function to be evaluated.

**MEDIAN.** MEDIAN(numexpr,numexpr[,..]). Numeric. Returns the median (50th percentile) of its arguments that have valid, nonmissing values. This function requires two or more arguments, which must be numeric. You can specify a minimum number of valid arguments for this function to be evaluated.

**SD.** SD(numexpr,numexpr[,...]). Numeric. Returns the standard deviation of its arguments that have valid, nonmissing values. This function requires two or more arguments, which must be numeric. You can specify a minimum number of valid arguments for this function to be evaluated.

**VARIANCE.** VARIANCE(numexpr,numexpr[,...]). Numeric. Returns the variance of its arguments that have valid values. This function requires two or more arguments, which must be numeric. You can specify a minimum number of valid arguments for this function to be evaluated.

**CFVAR.** CFVAR(numexpr,numexpr[,...]). Numeric. Returns the coefficient of variation (the standard deviation divided by the mean) of its arguments that have valid values. This function requires two or more arguments, which must be numeric. You can specify a minimum number of valid arguments for this function to be evaluated.

**MIN.** MIN(value,value[,...]). Numeric or string. Returns the minimum value of its arguments that have valid, nonmissing values. This function requires two or more arguments. For numeric values, you can specify a minimum number of valid arguments for this function to be evaluated.

**MAX.** MAX(value,value[,...]). Numeric or string. Returns the maximum value of its arguments that have valid values. This function requires two or more arguments. For numeric values, you can specify a minimum number of valid arguments for this function to be evaluated.

### **Example**

```
COMPUTE maxsum=MAX.2(SUM(var1 TO var3), SUM(var4 TO var6)).
```

- MAX.2 will return the maximum of the two sums provided that both sums are nonmissing.
- The .2 refers to the number of nonmissing arguments for the MAX function, which has only two arguments because each SUM function is considered a single argument.
- The new variable *maxsum* will be nonmissing if at least one variable specified for each SUM function is nonmissing.

## **Random variable and distribution functions**

Random variable and distribution function keywords are all of the form `prefix.suffix`, where the prefix specifies the function to be applied to the distribution and the suffix specifies the distribution.

- Random variable and distribution functions take both constants and variables for arguments.
- A function argument, if required, must come first and is denoted by  $x$  (quantile, which must fall in the range of values for the distribution) for cumulative distribution and probability density functions and  $p$  (probability) for inverse distribution functions.
- All random variable and distribution functions must specify distribution parameters as noted in their definitions.
- All arguments are real numbers.
- Restrictions to distribution parameters apply to all functions for that distribution. Restrictions for the function parameter  $x$  apply to that particular distribution function. The program issues a warning and returns system-missing when it encounters an out-of-range value for an argument.

The following are possible prefixes:

<b>CDF</b>	<i>Cumulative distribution function.</i> A cumulative distribution function <code>CDF.d_spec(x, a, ...)</code> returns a probability $p$ that a variate with the specified distribution ( <code>d_spec</code> ) falls below $x$ for continuous functions and at or below $x$ for discrete functions.
<b>IDF</b>	<i>Inverse distribution function.</i> Inverse distribution functions are not available for discrete distributions. An inverse distribution function <code>IDF.d_spec(p, a, ...)</code> returns a value $x$ such that <code>CDF.d_spec(x, a, ...)</code> = $p$ with the specified distribution ( <code>d_spec</code> ).
<b>PDF</b>	<i>Probability density function.</i> A probability density function <code>PDF.d_spec(x, a, ...)</code> returns the density of the specified distribution ( <code>d_spec</code> ) at $x$ for continuous functions and the probability that a random variable with the specified distribution equals $x$ for discrete functions.
<b>RV</b>	<i>Random number generation function.</i> A random number generation function <code>RV.d_spec(a, ...)</code> generates an independent observation with the specified distribution ( <code>d_spec</code> ).
<b>NCDF</b>	<i>Noncentral cumulative distribution function.</i> A noncentral distribution function <code>NCDF.d_spec(x, a, b, ...)</code> returns a probability $p$ that a variate with the specified noncentral distribution falls below $x$ . It is available only for beta, chi-square, $F$ , and Student's $t$ .
<b>NPDF</b>	<i>Noncentral probability density function.</i> A noncentral probability density function <code>NCDF.d_spec(x, a, ...)</code> returns the density of the specified distribution ( <code>d_spec</code> ) at $x$ . It is available only for beta, chi-square, $F$ , and Student's $t$ .
<b>SIG</b>	<i>Tail probability function.</i> A tail probability function <code>SIG.d_spec(x, a, ...)</code> returns a probability $p$ that a variate with the specified distribution ( <code>d_spec</code> ) is larger than $x$ . The tail probability function is equal to 1 minus the cumulative distribution function.

The following are suffixes for continuous distributions:

<b>BETA</b>	<i>Beta distribution.</i> The beta distribution takes values in the range $0 < x < 1$ and has two shape parameters, $\alpha$ and $\beta$ . Both $\alpha$ and $\beta$ must be positive, and they have the property that the mean of the distribution is $\alpha/(\alpha+\beta)$ . <b>Common uses.</b> The beta distribution is used in Bayesian analyses as a conjugate to the binomial distribution.
-------------	---

**Functions.** The CDF, IDF, PDF, NCDF, NPDF, and RV functions are available.

The beta distribution has PDF, CDF, and IDF

$$f(x; \alpha, \beta) = \frac{1}{B(\alpha, \beta)} x^{\alpha-1} (1-x)^{\beta-1}$$

$$F(x; \alpha, \beta) = \text{IB}(x; \alpha, \beta)$$

$$F^{-1}(p; \alpha, \beta) = \text{IB}^{-1}(p; \alpha, \beta)$$

where  $B(a, b) = \int_0^1 x^{a-1} (1-x)^{b-1} dx$  is the beta function and

$\text{IB}(x; a, b) = \int_0^x \frac{1}{B(a, b)} t^{a-1} (1-t)^{b-1} dt$  is the incomplete beta function.

**Relationship to other distributions.**

- When  $\alpha=\beta=1$ , the beta( $\alpha,\beta$ ) distribution is equivalent to the uniform(0,1) distribution.
- The beta( $\alpha,\beta$ ) distribution is the distribution of  $X/(X+Y)$  where  $X$  and  $Y$  are variables that have chi-square distributions with degrees of freedom parameters  $2\alpha$  and  $2\beta$ , respectively.

*Noncentral beta distribution.* The noncentral beta distribution is a generalization of the beta distribution that takes values in the range  $0 < x < 1$  and has an extra noncentrality parameter,  $\lambda$ , which must be greater than or equal to 0.

**Functions.**

The noncentral beta distribution has PDF, CDF, and IDF

$$f(x; \alpha, \beta, \lambda) = \sum_{j=0}^{\infty} \frac{1}{j!} \left(\frac{\lambda}{2}\right)^j e^{-\lambda/2} \frac{x^{\alpha+j-1} (1-x)^{\beta-1}}{B(\alpha+j; \beta)}$$

$$F(x; \alpha, \beta, \lambda) = \sum_{j=0}^{\infty} \frac{1}{j!} \left(\frac{\lambda}{2}\right)^j e^{-\lambda/2} \text{IB}(x; \alpha = j, \beta)$$

where  $B(a, b) = \int_0^1 x^{a-1} (1-x)^{b-1} dx$  is the beta function and

$\text{IB}(x; a, b) = \int_0^x \frac{1}{B(a, b)} t^{a-1} (1-t)^{b-1} dt$  is the incomplete beta function.

**Relationship to other distributions.**

- When  $\lambda$  equals 0, this distribution reduces to the beta distribution.
- The noncentral beta( $\alpha,\beta,\lambda$ ) distribution is the distribution of  $X/(X+Y)$  where  $X$  is a variable that has a noncentral chi-square( $2\alpha,\lambda$ ) distribution, and  $Y$  is a variable that has a central chi-square( $2\beta$ ) distribution.

**BVNOR**

*Bivariate normal distribution.* The bivariate normal distribution takes real values and has one correlation parameter,  $\rho$ , which must be between  $-1$  and  $1$ , inclusive.

**Functions.** The CDF and PDF functions are available and require two quantiles,  $x_1$  and  $x_2$ .

The bivariate normal distribution has PDF

$$f(x_1, x_2; \rho) = \frac{1}{2\pi(1-\rho^2)^{1/2}} \exp\left(\frac{-1}{2(1-\rho^2)}(x_1^2 - 2\rho x_1 x_2 + x_2^2)\right)$$

The CDF does not have a closed form and is computed by approximation.

**Relationship to other distributions.**

- Two variables with a bivariate normal( $\rho$ ) distribution with correlation  $\rho$  have marginal normal distributions with a mean of 0 and a standard deviation of 1.

**CAUCHY**

*Cauchy distribution.* The Cauchy distribution takes real values and has a location parameter,  $\theta$ , and a scale parameter,  $\zeta$ ;  $\zeta$  must be positive. The Cauchy distribution is symmetric about the location parameter, but has such slowly decaying tails that the distribution does not have a computable mean.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The Cauchy distribution has PDF, CDF, and IDF

$$f(x; \theta, \varsigma) = \frac{1}{\pi\varsigma} \left(1 + \left(\frac{x-\theta}{\varsigma}\right)^2\right)^{-1}$$

$$F(x; \theta, \varsigma) = \frac{1}{2} + \frac{1}{\pi} \tan^{-1}\left(\frac{x-\theta}{\varsigma}\right)$$

$$F^{-1}(p; \theta, \varsigma) = \theta + \varsigma \tan(\pi(p - 1/2))$$

**Relationship to other distributions.**

- A “standardized” Cauchy variate,  $(x-\theta)/\varsigma$ , has a  $t$  distribution with 1 degree of freedom.

## CHISQ

*Chi-square distribution.* The chi-square( $\nu$ ) distribution takes values in the range  $x \geq 0$  and has one degrees of freedom parameter,  $\nu$ ; it must be positive and has the property that the mean of the distribution is  $\nu$ .

**Functions.** The CDF, IDF, PDF, RV, NCDF, NPDF, and SIG functions are available.

The chi-square distribution has PDF, CDF, and IDF

$$f(x; \nu) = \frac{1}{2^{\nu/2} \Gamma(\nu/2)} x^{(\nu/2)-1} e^{-x/2}$$

$$F(x; \nu) = IG\left(\frac{x}{2}; \frac{\nu}{2}\right)$$

$$F^{-1}(p; \nu) = 2IG^{-1}\left(p; \frac{\nu}{2}\right)$$

where  $\Gamma(a) = \int_0^{\infty} x^{a-1} e^{-x} dx$  is the gamma function and

$IG(x; a) = \int_0^x \frac{1}{\Gamma(a)} t^{a-1} e^{-t} dt$  is the incomplete gamma function.

**Relationship to other distributions.**

- The chi-square( $\nu$ ) distribution is the distribution of the sum of squares of  $\nu$  independent normal(0,1) random variates.
- The chi-square( $\nu$ ) distribution is equivalent to the gamma( $\nu/2$ ,  $1/2$ ) distribution.

*Noncentral chi-square distribution.* The noncentral chi-square distribution is a generalization of the chi-square distribution that takes values in the range  $x \geq 0$  and has an extra noncentrality parameter,  $\lambda$ , which must be greater than or equal to 0.

**Functions.**

The noncentral chi-square distribution has PDF and CDF

$$f(x; \nu, \lambda) = \sum_{j=0}^{\infty} \frac{1}{j!} \left(\frac{\lambda}{2}\right)^j e^{-\lambda/2} \frac{x^{\nu/2+j-1} e^{-x/2}}{2^{\nu/2+j} \Gamma(\nu/2+j)}$$

$$F(x; \nu, \lambda) = \sum_{j=0}^{\infty} \frac{1}{j!} \left(\frac{\lambda}{2}\right)^j e^{-\lambda/2} \text{IG}\left(\frac{x}{2}; \frac{\nu}{2} + j\right)$$

where  $\Gamma(a) = \int_0^{\infty} x^{a-1} e^{-x} dx$  is the gamma function and

$\text{IG}(x; a) = \int_0^x \frac{1}{\Gamma(a)} t^{a-1} e^{-t} dt$  is the incomplete gamma function.

**Relationship to other distributions.**

- When  $\lambda$  equals 0, this distribution reduces to the chi-square distribution.
- The noncentral chi-square( $\nu, \lambda$ ) distribution is the distribution of the sum of squares of  $\nu$  independent normal( $\mu_i, 1$ ) random variates. Then  $\lambda = \sum \mu_i^2$ .

**EXP**

*Exponential distribution.* The exponential distribution takes values in the range  $x \geq 0$  and has one scale parameter,  $\beta$ , which must be greater than 0 and has the property that the mean of the distribution is  $1/\beta$ .

**Common uses.** In life testing, the scale parameter  $a$  represents the rate of decay.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The exponential distribution has PDF, CDF, and IDF

$$f(x; \beta) = \beta e^{-\beta x}$$

$$F(x; \beta) = 1 - e^{-\beta x}$$

$$F^{-1}(p; \beta) = -\frac{1}{\beta} \ln(1-p)$$

**Relationship to other distributions.**

- The exponential( $\beta$ ) distribution is equivalent to the gamma( $1, \beta$ ) distribution.

**F**

*F distribution.* The F distribution takes values in the range  $x \geq 0$  and has two degrees of freedom parameters,  $\nu_1$  and  $\nu_2$ , which are the “numerator” and “denominator” degrees of freedom, respectively. Both  $\nu_1$  and  $\nu_2$  must be positive.

**Common uses.** The  $F$  distribution is commonly used to test hypotheses under the Gaussian assumption.

**Functions.** The CDF, IDF, IDDF, RV, NCDF, NPDF, and SIG functions are available.

The  $F$  distribution has PDF, CDF, and IDF

$$f(x; \nu_1, \nu_2) = \frac{1}{B(\nu_1/2, \nu_2/2)} \left(\frac{\nu_1}{\nu_2}\right)^{\nu_1/2} x^{(\nu_1/2)-1} \left(1 + \frac{\nu_1}{\nu_2}x\right)^{-(\nu_1+\nu_2)/2}$$

$$F(x; \nu_1, \nu_2) = \text{IB}\left(\frac{\nu_1 x}{\nu_2 + \nu_1 x}; \frac{\nu_1}{2}, \frac{\nu_2}{2}\right)$$

$$F^{-1}(p; \nu_1, \nu_2) = \frac{\nu_2}{\nu_1} \left(\frac{\text{IB}^{-1}(p; \nu_1/2, \nu_2/2)}{1 - \text{IB}^{-1}(p; \nu_1/2, \nu_2/2)}\right)$$

where  $B(a, b) = \int_0^1 x^{a-1} (1-x)^{b-1} dx$  is the beta function and

$\text{IB}(x; a, b) = \int_0^x \frac{1}{B(a, b)} t^{a-1} (1-t)^{b-1} dt$  is the incomplete beta function.

#### Relationship to other distributions.

- The  $F(\nu_1, \nu_2)$  distribution is the distribution of  $(X/\nu_1)/(Y/\nu_2)$ , where  $X$  and  $Y$  are independent chi-square random variates with  $\nu_1$  and  $\nu_2$  degrees of freedom, respectively.

*Noncentral F distribution.* The noncentral  $F$  distribution is a generalization of the  $F$  distribution that takes values in the range  $x \geq 0$  and has an extra noncentrality parameter,  $\lambda$ , which must be greater than or equal to 0.

#### Functions.

The noncentral  $F$  distribution has PDF and CDF

$$f(x; \nu_1, b, \lambda) = \sum_{j=0}^{\infty} \frac{1}{j!} \left(\frac{\lambda}{2}\right)^j e^{-\lambda/2} \frac{(\nu_1/b)^{\nu_1/2+j}}{B(\nu_1/2+j, b/2)} x^{\nu_1/2+j-1} \left(1 + \frac{\nu_1}{b}x\right)^{-((\nu_1+b)/2+j)}$$

$$F(x; \nu_1, b, \lambda) = \sum_{j=0}^{\infty} \frac{1}{j!} \left(\frac{\lambda}{2}\right)^j e^{-\lambda/2} \text{IB}\left(\frac{\nu_1 x}{b + \nu_1 x}; \frac{\nu_1}{2} + j, \frac{b}{2}\right)$$

where  $B(a, b) = \int_0^1 x^{a-1} (1-x)^{b-1} dx$  is the beta function and

$\text{IB}(x; a, b) = \int_0^x \frac{1}{B(a, b)} t^{a-1} (1-t)^{b-1} dt$  is the incomplete beta function.

#### Relationship to other distributions.

- When  $\lambda$  equals 0, this distribution reduces to the  $F$  distribution.
- The noncentral  $F$  distribution is the distribution of  $(X/\nu_1)/(Y/\nu_2)$ , where  $X$  and  $Y$  are independent variates with noncentral chi-square( $\nu_1, \lambda$ ) and central chi-square( $\nu_2$ ) distributions, respectively.

## GAMMA

*Gamma distribution.* The gamma distribution takes values in the range  $x \geq 0$  and has one shape parameter,  $\alpha$ , and one scale parameter,  $\beta$ . Both parameters must be positive and have the property that the mean of the distribution is  $\alpha/\beta$ .

**Common uses.** The gamma distribution is commonly used in queuing theory, inventory control, and precipitation processes.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The gamma distribution has PDF, CDF, and IDF

$$f(x; \alpha, \beta) = \frac{\beta^\alpha}{\Gamma(\alpha)} x^{\alpha-1} e^{-\beta x}$$

$$F(x; \alpha, \beta) = \text{IG}(\beta x; \alpha)$$

$$F^{-1}(p; \alpha, \beta) = \frac{1}{\beta} \text{IG}^{-1}(p; \alpha)$$

where  $\Gamma(a) = \int_0^\infty x^{a-1} e^{-x} dx$  is the gamma function and

$\text{IG}(x; a) = \int_0^x \frac{1}{\Gamma(a)} t^{a-1} e^{-t} dt$  is the incomplete gamma function.

**Relationship to other distributions.**

- When  $\alpha=1$ , the gamma( $\alpha, \beta$ ) distribution reduces to the exponential( $\beta$ ) distribution.
- When  $\beta=1/2$ , the gamma( $\alpha, \beta$ ) distribution reduces to the chi-square( $2\alpha$ ) distribution.
- When  $\alpha$  is an integer, the gamma distribution is also known as the Erlang distribution.

## HALFNRM

*Half-normal distribution.* The half-normal distribution takes values in the range  $x \geq \mu$  and has one location parameter,  $\mu$ , and one scale parameter,  $\sigma$ . Parameter  $\sigma$  must be positive.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The half-normal distribution has PDF, CDF, and IDF

$$f(x; \mu, \sigma) = \phi\left(\frac{x-\mu}{\sigma}\right) / \sigma$$

$$F(x; \mu, \sigma) = 2\Phi\left(\frac{x-\mu}{\sigma}\right) - 1$$

$$F^{-1}(p; \mu, \sigma) = \mu + \sigma \Phi^{-1}\left(\frac{1+p}{2}\right)$$

**Relationship to other distributions.**

- If  $X$  has a normal( $\mu, \sigma$ ) distribution, then  $|X-\mu|$  has a half-normal( $\mu, \sigma$ ) distribution.

## IGAUSS

*Inverse Gaussian distribution.* The inverse Gaussian, or Wald, distribution takes values in the range  $x > 0$  and has two parameters,  $\mu$  and  $\lambda$ , both of which must be positive. The distribution has mean  $\mu$ .



**Common uses.** The inverse Gaussian distribution is commonly used to test hypotheses for model parameter estimates.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The inverse Gaussian distribution has PDF and CDF

$$f(x; \mu, \lambda) = \left(\frac{\lambda}{2\pi x^3}\right)^{1/2} \exp\left(-\frac{\lambda(x-\mu)^2}{2\mu^2 x}\right)$$

$$F(x; \mu, \lambda) = \Phi\left(\sqrt{\frac{\lambda}{x}}\left(-1 + \frac{x}{\mu}\right)\right) + e^{(2\lambda/\mu)} \Phi\left(-\sqrt{\frac{\lambda}{x}}\left(1 + \frac{x}{\mu}\right)\right)$$

The IDF is computed by approximation.

## LAPLACE

*Laplace or double exponential distribution.* The Laplace distribution takes real values and has one location parameter,  $\mu$ , and one scale parameter,  $\beta$ . Parameter  $\beta$  must be positive. The distribution is symmetric about  $\mu$  and has exponentially decaying tails.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The Laplace distribution has PDF, CDF, and IDF

$$f(x; \mu, \beta) = \frac{1}{2\beta} e^{-|x-\mu|/\beta}$$

$$F(x; \mu, \beta) = \begin{cases} \frac{1}{2} e^{(x-\mu)/\beta} & x \leq \mu \\ 1 - \frac{1}{2} e^{(\mu-x)/\beta} & x > \mu \end{cases}$$

$$F^{-1}(p; \mu, \beta) = \begin{cases} \mu + \beta \ln(2p) & 0 \leq p \leq \frac{1}{2} \\ \mu - \beta \ln(2(1-p)) & \frac{1}{2} < p \leq 1 \end{cases}$$

## LOGISTIC

*Logistic distribution.* The logistic distribution takes real values and has one location parameter,  $\mu$ , and one scale parameter,  $\varsigma$ . Parameter  $\varsigma$  must be positive. The distribution is symmetric about  $\mu$  and has longer tails than the normal distribution.

**Common uses.** The logistic distribution is used to model growth curves.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The logistic distribution has PDF, CDF, and IDF

$$f(x; \mu, \varsigma) = \frac{1}{\varsigma} e^{-(x-\mu)/\varsigma} \left(1 + e^{-(x-\mu)/\varsigma}\right)^{-2}$$

$$F(x; \mu, \varsigma) = \frac{1}{1 + e^{-(x-\mu)/\varsigma}}$$

$$F^{-1}(p; \mu, \varsigma) = \mu + \varsigma \ln\left(\frac{p}{1-p}\right)$$

## LNORMAL

*Lognormal distribution.* The lognormal distribution takes values in the range  $x \geq 0$  and has two parameters,  $\eta$  and  $\sigma$ , both of which must be positive.

**Common uses.** Lognormal is used in the distribution of particle sizes in aggregates, flood flows, concentrations of air contaminants, and failure time.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The lognormal distribution has PDF, CDF, and IDF

$$f(x; \eta, \sigma) = \frac{1}{x\sigma\sqrt{2\pi}} e^{-(\ln(x/\eta))^2/(2\sigma^2)}$$

$$F(x; \eta, \sigma) = \Phi\left(\frac{1}{\sigma} \ln\left(\frac{x}{\eta}\right)\right)$$

$$F^{-1}(p; \eta, \sigma) = \eta e^{\sigma\Phi^{-1}(p)}$$

**Relationship to other distributions.**

- If  $X$  has a lognormal( $\eta, \sigma$ ) distribution, then  $\ln(X)$  has a normal( $\ln(\eta), \sigma$ ) distribution.

## NORMAL

*Normal distribution.* The normal, or Gaussian, distribution takes real values and has one location parameter,  $\mu$ , and one scale parameter,  $\sigma$ . Parameter  $\sigma$  must be positive. The distribution has mean  $\mu$  and standard deviation  $\sigma$ .

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The normal distribution has PDF, CDF, and IDF

$$f(x; \mu, \sigma) = \frac{1}{\sigma\sqrt{2\pi}} e^{-(x-\mu)^2/(2\sigma^2)}$$

$$F(x; \mu, \sigma) = \Phi\left(\frac{x-\mu}{\sigma}\right)$$

$$F^{-1}(p; \mu, \sigma) = \mu + \sigma\Phi^{-1}(p)$$

**Relationship to other distributions.**

- If  $X$  has a normal( $\mu, \sigma$ ) distribution, then  $\exp(X)$  has a normal( $\exp(\mu), \sigma$ ) distribution.

Three functions in releases earlier than 6.0 are special cases of the normal distribution functions: `CDFNORM(arg)=CDF.NORMAL(x, 0, 1)`, where `arg` is  $x$ ; `PROBIT(arg)=IDF.NORMAL(p, 0, 1)`, where `arg` is  $p$ ; and `NORMAL(arg)=RV.NORMAL(0, sigma)`, where `arg` is  $\sigma$ .

## PARETO

*Pareto distribution.* The Pareto distribution takes values in the range  $x_{\min} < x$  and has a threshold parameter,  $x_{\min}$ , and a shape parameter,  $\alpha$ . Both parameters must be positive.

**Common uses.** Pareto is commonly used in economics as a model for a density function with a slowly decaying tail.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The Pareto distribution has PDF, CDF, and IDF

$$f(x; x_{\min}, \alpha) = \frac{\alpha}{x_{\min}} \left( \frac{x_{\min}}{x} \right)^{\alpha+1}$$

$$F(x; x_{\min}, \alpha) = 1 - \left( \frac{x_{\min}}{x} \right)^{\alpha}$$

$$F^{-1}(p; x_{\min}, \alpha) = x_{\min} (1 - p)^{-1/\alpha}$$

### SMOD

*Studentized maximum modulus distribution.* The Studentized maximum modulus distribution takes values in the range  $x > 0$  and has a number of comparisons parameter,  $k^*$ , and degrees of freedom parameter,  $\nu$ , both of which must be greater than or equal to 1.

**Common uses.** The Studentized maximum modulus is commonly used in post hoc multiple comparisons for GLM and ANOVA.

**Functions.** The CDF and IDF functions are available, and are computed by approximation.

### SRANGE

*Studentized range distribution.* The Studentized range distribution takes values in the range  $x > 0$  and has a number of samples parameter,  $k$ , and degrees of freedom parameter,  $\nu$ , both of which must be greater than or equal to 1.

**Common uses.** The Studentized range is commonly used in post hoc multiple comparisons for GLM and ANOVA.

**Functions.** The CDF and IDF functions are available, and are computed by approximation.

### T

*Student t distribution.* The Student t distribution takes real values and has one degrees of freedom parameter,  $\nu$ , which must be positive. The Student t distribution is symmetric about 0.

**Common uses.** The major uses of the Student  $t$  distribution are to test hypotheses and construct confidence intervals for means of data.

**Functions.** The CDF, IDF, PDF, RV, NCDF, and NPDF functions are available.

The  $t$  distribution has PDF, CDF, and IDF

$$f(x; \nu) = \frac{1}{\sqrt{\nu} B(\nu/2, 1/2)} \left( 1 + \frac{x^2}{\nu} \right)^{-(\nu+1)/2}$$

$$F(x; \nu) = \begin{cases} \frac{1}{2} \text{IB} \left( \frac{\nu}{\nu+x^2}; \frac{\nu}{2}, \frac{1}{2} \right) & x \leq 0 \\ 1 - \frac{1}{2} \text{IB} \left( \frac{\nu}{\nu+x^2}; \frac{\nu}{2}, \frac{1}{2} \right) & x > 0 \end{cases}$$

$$F^{-1}(p; \nu) = \begin{cases} -\sqrt{\nu(1/(\text{IB}^{-1}(2p; \frac{\nu}{2}, \frac{1}{2})) - 1)} & 0 \leq p \leq \frac{1}{2} \\ \sqrt{\nu(1/(\text{IB}^{-1}(2(1-p); \frac{\nu}{2}, \frac{1}{2})) - 1)} & \frac{1}{2} < p \leq 1 \end{cases}$$

where  $B(a, b) = \int_0^1 x^{a-1}(1-x)^{b-1} dx$  is the beta function and  
 $IB(x; a, b) = \int_0^x \frac{1}{B(a, b)} t^{a-1}(1-t)^{b-1} dt$  is the incomplete beta function.

**Relationship to other distributions.**

- The  $t(\nu)$  distribution is the distribution of  $X/Y$ , where  $X$  is a normal(0,1) variate and  $Y$  is a chi-square( $\nu$ ) variate divided by  $\nu$ .
- The square of a  $t(\nu)$  variate has an  $F(1, \nu)$  distribution.
- The  $t(\nu)$  distribution approaches the normal(0,1) distribution as  $\nu$  approaches infinity.

*Noncentral t distribution.* The noncentral t distribution is a generalization of the t distribution that takes real values and has an extra noncentrality parameter,  $\lambda$ , which must be greater than or equal to 0. When  $\lambda$  equals 0, this distribution reduces to the t distribution.

**Functions.**

The noncentral t distribution has PDF and CDF

$$f(x; \nu, \lambda) = \sum_{j=0}^{\infty} \frac{1}{j!} (\lambda\sqrt{2})^j e^{-\lambda^2/2} \frac{\Gamma((\nu+j+1)/2)}{\Gamma(\nu/2)\Gamma(1/2)} \frac{x^j}{\nu^{(j+1)/2}} \left(1 + \frac{x^2}{\nu}\right)^{-(\nu+j+1)/2}$$

$$F(x; \nu, \lambda) = \begin{cases} \frac{1}{2} \sum_{j=0}^{\infty} \frac{1}{j!} (-\lambda\sqrt{2})^j e^{-\lambda^2/2} \frac{\Gamma((j+1)/2)}{\Gamma(1/2)} IB\left(\frac{\nu}{\nu+x^2}; \frac{\nu}{2}, \frac{j+1}{2}\right) & x \leq 0 \\ 1 - \frac{1}{2} \sum_{j=0}^{\infty} \frac{1}{j!} (\lambda\sqrt{2})^j e^{-\lambda^2/2} \frac{\Gamma((j+1)/2)}{\Gamma(1/2)} IB\left(\frac{\nu}{\nu+x^2}; \frac{\nu}{2}, \frac{j+1}{2}\right) & x > 0 \end{cases}$$

where  $B(a, b) = \int_0^1 x^{a-1}(1-x)^{b-1} dx$  is the beta function and  
 $IB(x; a, b) = \int_0^x \frac{1}{B(a, b)} t^{a-1}(1-t)^{b-1} dt$  is the incomplete beta function.

**Relationship to other distributions.**

- The noncentral  $t(\nu, \lambda)$  distribution is the distribution of  $X/Y$ , where  $X$  is a normal( $\lambda, 1$ ) variate and  $Y$  is a central chi-square( $\nu$ ) variate divided by  $\nu$ .

**UNIFORM**

*Uniform distribution.* The uniform distribution takes values in the range  $a < x < b$  and has a minimum value parameter,  $a$ , and a maximum value parameter,  $b$ .

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The uniform distribution has PDF, CDF, and IDF

$$f(x; a, b) = \frac{1}{b-a}$$

$$F(x; a, b) = \frac{x-a}{b-a}$$

$$F^{-1}(p; a, b) = a + (b-a)p$$

The uniform random number function in releases earlier than 6.0 is a special case:  $\text{UNIFORM}(\text{arg}) = \text{RV} . \text{UNIFORM}(0, b)$ , where  $\text{arg}$  is parameter  $b$ . Among other uses, the uniform distribution commonly models the round-off error.

## WEIBULL

*Weibull distribution.* The Weibull distribution takes values in the range  $x \geq 0$  and has one scale parameter,  $\beta$ , and one shape parameter,  $\alpha$ , both of which must be positive.

**Common uses.** The Weibull distribution is commonly used in survival analysis.

**Functions.** The CDF, IDF, PDF, and RV functions are available.

The Weibull distribution has PDF, CDF, and IDF

$$f(x; \beta, \alpha) = \frac{\alpha}{\beta} \left(\frac{x}{\beta}\right)^{\alpha-1} e^{-(x/\beta)^\alpha}$$

$$F(x; \beta, \alpha) = 1 - e^{-(x/\beta)^\alpha}$$

$$F^{-1}(p; \beta, \alpha) = \beta(-\ln(1-p))^{1/\alpha}$$

**Relationship to other distributions.**

- A Weibull( $\beta, 1$ ) distribution is equivalent to an exponential( $\beta$ ) distribution.

The following are suffixes for discrete distributions:

## BERNOULLI

*Bernoulli distribution.* The Bernoulli distribution takes values 0 or 1 and has one success probability parameter,  $\theta$ , which must be between 0 and 1, inclusive.

**Functions.** The CDF, PDF, and RV functions are available.

The Bernoulli distribution has PDF and CDF

$$f(x; \theta) = \theta^x (1 - \theta)^{1-x}$$

$$F(x; \theta) = \begin{cases} 1 - \theta & x = 0 \\ 1 & x = 1 \end{cases}$$

**Relationship to other distributions.**

- The Bernoulli distribution is a special case of the binomial distribution and is used in simple success-failure experiments.

## BINOM

*Binomial distribution.* The binomial distribution takes integer values  $0 \leq x \leq n$ , representing the number of successes in  $n$  trials, and has one number of trials parameter,  $n$ , and one success probability parameter,  $\theta$ . Parameter  $n$  must be a positive integer and parameter  $\theta$  must be between 0 and 1, inclusive.

**Common uses.** The binomial distribution is used in independently replicated success-failure experiments.

**Functions.** The CDF, PDF, and RV functions are available.

The binomial distribution has PDF and CDF

$$f(x; n, \theta) = \binom{n}{x} \theta^x (1 - \theta)^{n-x}$$

$$F(x; n, \theta) = \begin{cases} 1 - \text{IB}(\theta; x + 1, n - x) & x = 0, 1, \dots, n - 1 \\ 1 & x = n \end{cases}$$

where  $\text{IB}(x; a, b) = \int_0^x \frac{1}{\text{B}(a, b)} t^{a-1} (1-t)^{b-1} dt$  is the incomplete beta function.

## GEOM

*Geometric distribution.* The geometric distribution takes integer values  $x \geq 1$ , representing the number of trials needed (including the last trial) before a success is observed, and has one success probability parameter,  $\theta$ , which must be between 0 and 1, inclusive.

**Functions.** The CDF, PDF, and RV functions are available.

The geometric distribution has PDF and CDF

$$f(x; \theta) = \theta(1 - \theta)^{x-1}$$

$$F(x; \theta) = 1 - (1 - \theta)^x$$

### Relationship to other distributions.

- The geometric( $\theta$ ) distribution is equivalent to the negative binomial ( $1, \theta$ ) distribution.

## HYPER

*Hypergeometric distribution.* The hypergeometric distribution takes integer values in the range  $\max(0, N_p + n - N) \leq x \leq \min(N_p, n)$ , and has three parameters,  $N$ ,  $n$ , and  $N_p$ , where  $N$  is the total number of objects in an urn model,  $n$  is the number of objects randomly drawn without replacement from the urn,  $N_p$  is the number of objects with a given characteristic, and  $x$  is the number of objects with the given characteristic observed out of the withdrawn objects. All three parameters are positive integers, and both  $n$  and  $N_p$  must be less than or equal to  $N$ .

**Functions.** The CDF, PDF, and RV functions are available.

The hypergeometric distribution has PDF and CDF

$$f(x; N, n, Np) = \frac{\binom{Np}{x} \binom{N-Np}{n-x}}{\binom{N}{n}}$$

$$F(x; N, n, Np) = \sum_{k=\max(0, n+Np-N)}^x \text{Prob}(X=k)$$

## NEGBIN

*Negative binomial distribution.* The negative binomial distribution takes integer values in the range  $x \geq r$ , where  $x$  is the number of trials needed (including the last trial) before  $r$  successes are observed, and has one threshold parameter,  $r$ , and one success probability parameter,  $\theta$ . Parameter  $r$  must be a positive integer and parameter  $\theta$  must be greater than 0 and less than or equal to 1.

**Functions.** The CDF, PDF, and RV functions are available.

The negative binomial distribution has PDF and CDF

$$f(x; r, \theta) = \binom{x-1}{r-1} \theta^r (1-\theta)^{x-r}$$

$$F(x; r, \theta) = \text{IB}(\theta; r, x - r + 1)$$

where  $\text{IB}(x; a, b) = \int_0^x \frac{1}{\text{B}(a, b)} t^{a-1} (1-t)^{b-1} dt$  is the incomplete beta function.

### Relationship to other distributions.

- The negative binomial(1,  $\theta$ ) distribution is equivalent to the geometric( $\theta$ ) distribution.

## POISSON

*Poisson distribution.* The Poisson distribution takes integer values in the range  $x \geq 0$  and has one rate or mean parameter,  $\lambda$ . Parameter  $\lambda$  must be positive.

**Common uses.** The Poisson distribution is used in modeling the distribution of counts, such as traffic counts and insect counts.

**Functions.** The CDF, PDF, and RV functions are available.

The Poisson distribution has PDF and CDF

$$f(x; \lambda) = \frac{\lambda^x}{x!} e^{-\lambda}$$

$$F(x; \lambda) = 1 - \text{IG}(\lambda; x+1)$$

where  $IG(x; a) = \int_0^x \frac{1}{\Gamma(a)} t^{a-1} e^{-t} dt$  is the incomplete gamma function.

### **Probability Density Functions**

The following functions give the value of the density function with the specified distribution at the value *quant*, the first argument. Subsequent arguments are the parameters of the distribution. Note the period in each function name.

**PDF.BERNOULLI.** PDF.BERNOULLI(quant, prob). Numeric. Returns the probability that a value from the Bernoulli distribution, with the given probability parameter, will be equal to quant.

**PDF.BETA.** PDF.BETA(quant, shape1, shape2). Numeric. Returns the probability density of the beta distribution, with the given shape parameters, at quant.

**PDF.BINOM.** PDF.BINOM(quant, n, prob). Numeric. Returns the probability that the number of successes in n trials, with probability prob of success in each, will be equal to quant. When n is 1, this is the same as PDF.BERNOULLI.

**PDF.BVNOR.** PDF.BVNOR(quant1, quant2, corr). Numeric. Returns the probability density of the standard bivariate normal distribution, with the given correlation parameter, at quant1, quant2.

**PDF.CAUCHY.** PDF.CAUCHY(quant, loc, scale). Numeric. Returns the probability density of the Cauchy distribution, with the given location and scale parameters, at quant.

**PDF.CHISQ.** PDF.CHISQ(quant, df). Numeric. Returns the probability density of the chi-square distribution, with df degrees of freedom, at quant.

**PDF.EXP.** PDF.EXP(quant, shape). Numeric. Returns the probability density of the exponential distribution, with the given shape parameter, at quant.

**PDF.F.** PDF.F(quant, df1, df2). Numeric. Returns the probability density of the F distribution, with degrees of freedom df1 and df2, at quant.

**PDF.GAMMA.** PDF.GAMMA(quant, shape, scale). Numeric. Returns the probability density of the gamma distribution, with the given shape and scale parameters, at quant.

**PDF.GEOM.** PDF.GEOM(quant, prob). Numeric. Returns the probability that the number of trials to obtain a success, when the probability of success is given by prob, will be equal to quant.

**PDF.HALFNRM.** PDF.HALFNRM(quant, mean, stddev). Numeric. Returns the probability density of the half normal distribution, with specified mean and standard deviation, at quant.

**PDF.HYPER.** PDF.HYPER(quant, total, sample, hits). Numeric. Returns the probability that the number of objects with a specified characteristic, when sample objects are randomly selected from a universe of size total in which hits have the specified characteristic, will be equal to quant.



**PDF.IGAUSS.** PDF.IGAUSS(quant, loc, scale). Numeric. Returns the probability density of the inverse Gaussian distribution, with the given location and scale parameters, at quant.

**PDF.LAPLACE.** PDF.LAPLACE(quant, mean, scale). Numeric. Returns the probability density of the Laplace distribution, with the specified mean and scale parameters, at quant.

**PDF.LOGISTIC.** PDF.LOGISTIC(quant, mean, scale). Numeric. Returns the probability density of the logistic distribution, with the specified mean and scale parameters, at quant.

**PDF.LNORMAL.** PDF.LNORMAL(quant, a, b). Numeric. Returns the probability density of the log-normal distribution, with the specified parameters, at quant.

**PDF.NEGBIN.** PDF.NEGBIN(quant, thresh, prob). Numeric. Returns the probability that the number of trials to obtain a success, when the threshold parameter is thresh and the probability of success is given by prob, will be equal to quant.

**PDF.NORMAL.** PDF.NORMAL(quant, mean, stddev). Numeric. Returns the probability density of the normal distribution, with specified mean and standard deviation, at quant.

**PDF.PARETO.** PDF.PARETO(quant, threshold, shape). Numeric. Returns the probability density of the Pareto distribution, with the specified threshold and shape parameters, at quant.

**PDF.POISSON.** PDF.POISSON(quant, mean). Numeric. Returns the probability that a value from the Poisson distribution, with the specified mean or rate parameter, will be equal to quant.

**PDF.T.** PDF.T(quant, df). Numeric. Returns the probability density of Student's t distribution, with the specified degrees of freedom df, at quant.

**PDF.UNIFORM.** PDF.UNIFORM(quant, min, max). Numeric. Returns the probability density of the uniform distribution, with the specified minimum and maximum, at quant.

**PDF.WEIBULL.** PDF.WEIBULL(quant, a, b). Numeric. Returns the probability density of the Weibull distribution, with the specified parameters, at quant.

**NPDF.BETA.** NPDF.BETA(quant, shape1, shape2, nc). Numeric. Returns the probability density of the noncentral beta distribution, with the given shape and noncentrality parameters, at quant.

**NPDF.CHISQ.** NPDF.CHISQ(quant, df, nc). Numeric. Returns the probability density of the noncentral chi-square distribution, with df degrees of freedom and the specified noncentrality parameter, at quant.

**NPDF.F.** NPDF.F(quant, df1, df2, nc). Numeric. Returns the probability density of the noncentral F distribution, with degrees of freedom df1 and df2 and noncentrality nc, at quant.

**NPDF.T.** NPDF.T(quant, df, nc). Numeric. Returns the probability density of the noncentral Student's t distribution, with the specified degrees of freedom df and noncentrality nc, at quant.

**Tail probability functions**

The following functions give the probability that a random variable with the specified distribution will be greater than *quant*, the first argument. Subsequent arguments are the parameters of the distribution. Note the period in each function name.

**SIG.CHISQ.** SIG.CHISQ(quant, df). Numeric. Returns the cumulative probability that a value from the chi-square distribution, with df degrees of freedom, will be greater than quant

**SIG.F.** These significance values should not be used to test hypotheses about the F values in this table. Cluster analysis specifically attempts to maximize between-group variance, and the significance values reported here do not reflect this.

**Cumulative distribution functions**

The following functions give the probability that a random variable with the specified distribution will be less than *quant*, the first argument. Subsequent arguments are the parameters of the distribution. Note the period in each function name.

**CDF.BERNOULLI.** CDF.BERNOULLI(quant, prob). Numeric. Returns the cumulative probability that a value from the Bernoulli distribution, with the given probability parameter, will be less than or equal to quant.

**CDF.BETA.** CDF.BETA(quant, shape1, shape2). Numeric. Returns the cumulative probability that a value from the Beta distribution, with the given shape parameters, will be less than quant.

**CDF.BINOM.** CDF.BINOM(quant, n, prob). Numeric. Returns the cumulative probability that the number of successes in n trials, with probability prob of success in each, will be less than or equal to quant. When n is 1, this is the same as CDF.BERNOULLI.

**CDF.BVNOR.** CDF.BVNOR(quant1, quant2, corr). Numeric. Returns the cumulative probability that a value from the standard bivariate normal distribution, with the given correlation parameter, will be less than quant1 and quant2.

**CDF.CAUCHY.** CDF.CAUCHY(quant, loc, scale). Numeric. Returns the cumulative probability that a value from the Cauchy distribution, with the given location and scale parameters, will be less than quant.

**CDF.CHISQ.** CDF.CHISQ(quant, df). Numeric. Returns the cumulative probability that a value from the chi-square distribution, with df degrees of freedom, will be less than quant.

**CDF.EXP.** CDF.EXP(quant, scale). Numeric. Returns the cumulative probability that a value from the exponential distribution, with the given scale parameter, will be less than quant.

**CDF.F.** CDF.F(quant, df1, df2). Numeric. Returns the cumulative probability that a value from the F distribution, with degrees of freedom df1 and df2, will be less than quant.

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**CDF.GAMMA.** CDF.GAMMA(quant, shape, scale). Numeric. Returns the cumulative probability that a value from the Gamma distribution, with the given shape and scale parameters, will be less than quant.

**CDF.GEOM.** CDF.GEOM(quant, prob). Numeric. Returns the cumulative probability that the number of trials to obtain a success, when the probability of success is given by prob, will be less than or equal to quant.

**CDF.HALFNRM.** CDF.HALFNRM(quant, mean, stddev). Numeric. Returns the cumulative probability that a value from the half normal distribution, with specified mean and standard deviation, will be less than quant.

**CDF.HYPER.** CDF.HYPER(quant, total, sample, hits). Numeric. Returns the cumulative probability that the number of objects with a specified characteristic, when sample objects are randomly selected from a universe of size total in which hits have the specified characteristic, will be less than or equal to quant.

**CDF.IGAUSS.** CDF.IGAUSS(quant, loc, scale). Numeric. Returns the cumulative probability that a value from the inverse Gaussian distribution, with the given location and scale parameters, will be less than quant.

**CDF.LAPLACE.** CDF.LAPLACE(quant, mean, scale). Numeric. Returns the cumulative probability that a value from the Laplace distribution, with the specified mean and scale parameters, will be less than quant.

**CDF.LOGISTIC.** CDF.LOGISTIC(quant, mean, scale). Numeric. Returns the cumulative probability that a value from the logistic distribution, with the specified mean and scale parameters, will be less than quant.

**CDF.LNORMAL.** CDF.LNORMAL(quant, a, b). Numeric. Returns the cumulative probability that a value from the log-normal distribution, with the specified parameters, will be less than quant.

**CDF.NEGBIN.** CDF.NEGBIN(quant, thresh, prob). Numeric. Returns the cumulative probability that the number of trials to obtain a success, when the threshold parameter is thresh and the probability of success is given by prob, will be less than or equal to quant.

**CDF.NORM.** CDF.NORM(zvalue). Numeric. Returns the probability that a random variable with mean 0 and standard deviation 1 would be less than zvalue, which must be numeric.

**CDF.NORMAL.** CDF.NORMAL(quant, mean, stddev). Numeric. Returns the cumulative probability that a value from the normal distribution, with specified mean and standard deviation, will be less than quant.

**CDF.PARETO.** CDF.PARETO(quant, threshold, shape). Numeric. Returns the cumulative probability that a value from the Pareto distribution, with the specified threshold and shape parameters, will be less than quant.

**CDF.POISSON.** CDF.POISSON(quant, mean). Numeric. Returns the cumulative probability that a value from the Poisson distribution, with the specified mean or rate parameter, will be less than or equal to quant.

**CDF.SMOD.** CDF.SMOD(quant, a, b). Numeric. Returns the cumulative probability that a value from the Studentized maximum modulus, with the specified parameters, will be less than quant.

**CDF.SRANGE.** CDF.SRANGE(quant, a, b). Numeric. Returns the cumulative probability that a value from the Studentized range statistic, with the specified parameters, will be less than quant.

**CDF.T.** CDF.T(quant, df). Numeric. Returns the cumulative probability that a value from Student's t distribution, with the specified degrees of freedom df, will be less than quant.

**CDF.UNIFORM.** CDF.UNIFORM(quant, min, max). Numeric. Returns the cumulative probability that a value from the uniform distribution, with the specified minimum and maximum, will be less than quant.

**CDF.WEIBULL.** CDF.WEIBULL(quant, a, b). Numeric. Returns the cumulative probability that a value from the Weibull distribution, with the specified parameters, will be less than quant.

**NCDF.BETA.** NCDF.BETA(quant, shape1, shape2, nc). Numeric. Returns the cumulative probability that a value from the noncentral Beta distribution, with the given shape and noncentrality parameters, will be less than quant.

**NCDF.CHISQ.** NCDF.CHISQ(quant, df, nc). Numeric. Returns the cumulative probability that a value from the noncentral chi-square distribution, with df degrees of freedom and the specified noncentrality parameter, will be less than quant.

**NCDF.F.** NCDF.F(quant, df1, df2, nc). Numeric. Returns the cumulative probability that a value from the noncentral F distribution, with degrees of freedom df1 and df2, and noncentrality nc, will be less than quant.

**NCDF.T.** NCDF.T(quant, df, nc). Numeric. Returns the cumulative probability that a value from the noncentral Student's t distribution, with the specified degrees of freedom df and noncentrality nc, will be less than quant.

### ***Inverse distribution functions***

The following functions give the value in a specified distribution having a cumulative probability equal to *prob*, the first argument. Subsequent arguments are the parameters of the distribution. Note the period in each function name.

**IDF.BETA.** IDF.BETA(prob, shape1, shape2). Numeric. Returns the value from the Beta distribution, with the given shape parameters, for which the cumulative probability is prob.

**IDF.CAUCHY.** IDF.CAUCHY(prob, loc, scale). Numeric. Returns the value from the Cauchy distribution, with the given location and scale parameters, for which the cumulative probability is prob.

**IDF.CHISQ.** IDF.CHISQ(prob, df). Numeric. Returns the value from the chi-square distribution, with the specified degrees of freedom df, for which the cumulative probability is prob. For example, the chi-square value that is significant at the 0.05 level with 3 degrees of freedom is IDF.CHISQ(0.95,3).

**IDF.EXP.** IDF.EXP(p, scale). Numeric. Returns the value of an exponentially decaying variable, with rate of decay scale, for which the cumulative probability is p.

**IDF.F.** IDF.F(prob, df1, df2). Numeric. Returns the value from the F distribution, with the specified degrees of freedom, for which the cumulative probability is prob. For example, the F value that is significant at the 0.05 level with 3 and 100 degrees of freedom is IDF.F(0.95,3,100).

**IDF.GAMMA.** IDF.GAMMA(prob, shape, scale). Numeric. Returns the value from the Gamma distribution, with the specified shape and scale parameters, for which the cumulative probability is prob.

**IDF.HALFNRM.** IDF.HALFNRM(prob, mean, stddev). Numeric. Returns the value from the half normal distribution, with the specified mean and standard deviation, for which the cumulative probability is prob.

**IDF.IGAUSS.** IDF.IGAUSS(prob, loc, scale). Numeric. Returns the value from the inverse Gaussian distribution, with the given location and scale parameters, for which the cumulative probability is prob.

**IDF.LAPLACE.** IDF.LAPLACE(prob, mean, scale). Numeric. Returns the value from the Laplace distribution, with the specified mean and scale parameters, for which the cumulative probability is prob.

**IDF.LOGISTIC.** IDF.LOGISTIC(prob, mean, scale). Numeric. Returns the value from the logistic distribution, with specified mean and scale parameters, for which the cumulative probability is prob.

**IDF.LNORMAL.** IDF.LNORMAL(prob, a, b). Numeric. Returns the value from the log-normal distribution, with specified parameters, for which the cumulative probability is prob.

**IDF.NORMAL.** IDF.NORMAL(prob, mean, stddev). Numeric. Returns the value from the normal distribution, with specified mean and standard deviation, for which the cumulative probability is prob.

**IDF.PARETO.** IDF.PARETO(prob, threshold, shape). Numeric. Returns the value from the Pareto distribution, with specified threshold and scale parameters, for which the cumulative probability is prob.

**IDF.SMOD.** IDF.SMOD(prob, a, b). Numeric. Returns the value from the Studentized maximum modulus, with the specified parameters, for which the cumulative probability is prob.

**IDF.SRANGE.** IDF.SRANGE(prob, a, b). Numeric. Returns the value from the Studentized range statistic, with the specified parameters, for which the cumulative probability is prob.

**IDF.T.** IDF.T(prob, df). Numeric. Returns the value from Student's t distribution, with specified degrees of freedom df, for which the cumulative probability is prob.

**IDF.UNIFORM.** IDF.UNIFORM(prob, min, max). Numeric. Returns the value from the uniform distribution between min and max for which the cumulative probability is prob.

**IDF.WEIBULL.** IDF.WEIBULL(prob, a, b). Numeric. Returns the value from the Weibull distribution, with specified parameters, for which the cumulative probability is prob.

**PROBIT.** PROBIT(prob). Numeric. Returns the value in a standard normal distribution having a cumulative probability equal to prob. The argument prob is a probability greater than 0 and less than 1.

### ***Random variable functions***

The following functions give a random variate from a specified distribution. The arguments are the parameters of the distribution. You can repeat the sequence of pseudorandom numbers by setting a seed in the Preferences dialog box before each sequence. Note the period in each function name.

**NORMAL.** NORMAL(stddev). Numeric. Returns a normally distributed pseudorandom number from a distribution with mean 0 and standard deviation stddev, which must be a positive number. You can repeat the sequence of pseudorandom numbers by setting a seed in the Random Number Seed dialog box before each sequence.

**RV.BERNOULLI.** RV.BERNOULLI(prob). Numeric. Returns a random value from a Bernoulli distribution with the specified probability parameter prob.

**RV.BETA.** RV.BETA(shape1, shape2). Numeric. Returns a random value from a Beta distribution with specified shape parameters.

**RV.BINOM.** RV.BINOM(n, prob). Numeric. Returns a random value from a binomial distribution with specified number of trials and probability parameter.

**RV.CAUCHY.** RV.CAUCHY(loc, scale). Numeric. Returns a random value from a Cauchy distribution with specified location and scale parameters.

**RV.CHISQ.** RV.CHISQ(df). Numeric. Returns a random value from a chi-square distribution with specified degrees of freedom df.

**RV.EXP.** RV.EXP(scale). Numeric. Returns a random value from an exponential distribution with specified scale parameter.

**RV.F.** RV.F(df1, df2). Numeric. Returns a random value from an F distribution with specified degrees of freedom, df1 and df2.

**RV.GAMMA.** RV.GAMMA(shape, scale). Numeric. Returns a random value from a Gamma distribution with specified shape and scale parameters.

**RV.GEOM.** RV.GEOM(prob). Numeric. Returns a random value from a geometric distribution with specified probability parameter.

**RV.HALFNRM.** RV.HALFNRM(mean, stddev). Numeric. Returns a random value from a half normal distribution with the specified mean and standard deviation.

**RV.HYPER.** RV.HYPER(total, sample, hits). Numeric. Returns a random value from a hypergeometric distribution with specified parameters.

**RV.IGAUSS.** RV.IGAUSS(loc, scale). Numeric. Returns a random value from an inverse Gaussian distribution with the specified location and scale parameters.

**RV.LAPLACE.** RV.LAPLACE(mean, scale). Numeric. Returns a random value from a Laplace distribution with specified mean and scale parameters.

**RV.LOGISTIC.** RV.LOGISTIC(mean, scale). Numeric. Returns a random value from a logistic distribution with specified mean and scale parameters.

**RV.LNORMAL.** RV.LNORMAL(a, b). Numeric. Returns a random value from a log-normal distribution with specified parameters.

**RV.NEGBIN.** RV.NEGBIN(threshold, prob). Numeric. Returns a random value from a negative binomial distribution with specified threshold and probability parameters.

**RV.NORMAL.** RV.NORMAL(mean, stddev). Numeric. Returns a random value from a normal distribution with specified mean and standard deviation.

**RV.PARETO.** RV.PARETO(threshold, shape). Numeric. Returns a random value from a Pareto distribution with specified threshold and shape parameters.

**RV.POISSON.** RV.POISSON(mean). Numeric. Returns a random value from a Poisson distribution with specified mean/rate parameter.

**RV.T.** RV.T(df). Numeric. Returns a random value from a Student's t distribution with specified degrees of freedom df.

**RV.UNIFORM.** RV.UNIFORM(min, max). Numeric. Returns a random value from a uniform distribution with specified minimum and maximum. See also the UNIFORM function.

**WEIBULL.** RV.WEIBULL(a, b). Numeric. Returns a random value from a Weibull distribution with specified parameters.

**UNIFORM.** UNIFORM(max). Numeric. Returns a uniformly distributed pseudorandom number between 0 and the argument max, which must be numeric (but can be negative). You can repeat the sequence of pseudorandom numbers by setting the same Random Number Seed (available in the Transform menu) before each sequence.

## **Date and time functions**

Date and time functions provide aggregation, conversion, and extraction routines for dates and time intervals. Each function transforms an expression consisting of one or more arguments. Arguments can be complex expressions, variable names, or constants. Date and time expressions and variables are legitimate arguments.

### **Aggregation functions**

Aggregation functions generate date and time intervals from values that were not read by date and time input formats.

- All aggregation functions begin with `DATE` or `TIME`, depending on whether a date or a time interval is requested. This is followed by a subfunction that corresponds to the type of values found in the data.
- The subfunctions are separated from the function by a period (`.`) and are followed by an argument list specified in parentheses.
- The arguments to the `DATE` and `TIME` functions must be separated by commas and must resolve to integer values.
- Functions that contain a day argument—for example, `DATE.DMY(d,m,y)`—check the validity of the argument. The value for day must be an integer between 1 and 31. If an invalid value is encountered, a warning is displayed and the value is set to system-missing. However, if the day value is invalid for a particular month—for example, 31 in September, April, June, and November or 29 through 31 for February in nonleap years—the resulting date is placed in the next month. For example `DATE.DMY(31, 9, 2006)` returns the date value for October 1, 2006.

**DATE.DMY.** `DATE.DMY(day,month,year)`. Numeric. Returns a date value corresponding to the indicated day, month, and year. The arguments must resolve to integers, with day between 1 and 31, month between 1 and 13, and year a four-digit integer greater than 1582. To display the result as a date, assign a date format to the result variable.

**DATE.MDY.** `DATE.MDY(month,day,year)`. Numeric. Returns a date value corresponding to the indicated month, day, and year. The arguments must resolve to integers, with day between 1 and 31, month between 1 and 13, and year a four-digit integer greater than 1582. To display the result as a date, assign a date format to the result variable.

**DATE.MOYR.** `DATE.MOYR(month,year)`. Numeric. Returns a date value corresponding to the indicated month and year. The arguments must resolve to integers, with month between 1 and 13, and year a four-digit integer greater than 1582. To display the result as a date, assign a date format to the result variable.

**DATE.QYR.** `DATE.QYR(quarter,year)`. Numeric. Returns a date value corresponding to the indicated quarter and year. The arguments must resolve to integers, with quarter between 1 and 4, and year a four-digit integer greater than 1582. To display the result as a date, assign a date format to the result variable.



**DATE.WKYR.** DATE.WKYR(weeknum,year). Numeric. Returns a date value corresponding to the indicated weeknum and year. The arguments must resolve to integers, with weeknum between 1 and 53, and year a four-digit integer greater than 1582. The date value returned represents the first day of the specified week for that year. The first week starts on January 1 of each year; so the date returned for any given week value will differ between years. To display the result as a date, assign a date format to the result variable.

**DATE.YRDAY.** DATE.YRDAY(year,daynum). Numeric. Returns a date value corresponding to the indicated year and daynum. The arguments must resolve to integers, with daynum between 1 and 366 and with year being a four-digit integer greater than 1582. To display the result as a date, assign a date format to the result variable.

**TIME.DAYS.** TIME.DAYS(days). Numeric. Returns a time interval corresponding to the indicated number of days. The argument must be numeric. To display the result as a time, assign a time format to the result variable.

**TIME.HMS.** TIME.HMS(hours[,minutes,seconds]). Numeric. Returns a time interval corresponding to the indicated number of hours, minutes, and seconds. The minutes and seconds arguments are optional. Minutes and seconds must resolve to numbers less than 60 if any higher-order argument is non-zero. All arguments except the last non-zero argument must resolve to integers. For example TIME.HMS(25.5) and TIME.HMS(0,90,25.5) are valid, while TIME.HMS(25.5,30) and TIME.HMS(25,90) are invalid. All arguments must resolve to either all positive or all negative values. To display the result as a time, assign a time format to the result variable.

### Example

```
DATA LIST FREE
  /Year Month Day Hour Minute Second Days.
BEGIN DATA
2006 10 28 23 54 30 1.5
END DATA.
COMPUTE Date1=DATE.DMY(Day, Month, Year).
COMPUTE Date2=DATE.MDY(Month, Day, Year).
COMPUTE MonthYear=DATE.MOYR(Month, Year).
COMPUTE Time=TIME.HMS(Hour, Minute, Second).
COMPUTE Duration=TIME.DAYS(Days).
LIST VARIABLES=Date1 to Duration.
FORMATS
  Date1 (DATE11) Date2 (ADATE10) MonthYear (MOYR8)
  Time (TIME8) Duration (Time8).
LIST VARIABLES=Date1 to Duration.

***LIST Results Before Applying Formats***
      Date1      Date2      MonthYear      Time Duration
13381372800 13381372800 13379040000 86070 129600

***LIST Results After Applying Formats***
      Date1      Date2 MonthYear      Time Duration
28-OCT-2006 10/28/2006  OCT 2006 23:54:30 36:00:00
```

- Since dates and times are stored internally as a number of seconds, prior to applying the appropriate date or time formats, all the computed values are displayed as numbers that indicate the respective number of seconds.
- The internal values for *Date1* and *Date2* are exactly the same. The only difference between DATE.DMY and DATE.MDY is the order of the arguments.

### **Date and time conversion functions**

The conversion functions convert time intervals from one unit of time to another. Time intervals are stored as the number of seconds in the interval; the conversion functions provide a means for calculating more appropriate units, for example, converting seconds to days.

Each conversion function consists of the `CTIME` function followed by a period (`.`), the target time unit, and an argument. The argument can consist of expressions, variable names, or constants. The argument must already be a time interval. [For more information, see the topic Aggregation functions on p. 98.](#) Time conversions produce noninteger results with a default format of `F8.2`.

Since time and dates are stored internally as seconds, a function that converts to seconds is not necessary.

**CTIME.DAYS.** `CTIME.DAYS(timevalue)`. Numeric. Returns the number of days, including fractional days, in `timevalue`, which is a number of seconds, a time expression, or a time format variable.

**CTIME.HOURS.** `CTIME.HOURS(timevalue)`. Numeric. Returns the number of hours, including fractional hours, in `timevalue`, which is a number of seconds, a time expression, or a time format variable.

**CTIME.MINUTES.** `CTIME.MINUTES(timevalue)`. Numeric. Returns the number of minutes, including fractional minutes, in `timevalue`, which is a number of seconds, a time expression, or a time format variable.

**CTIME.SECONDS.** `CTIME.SECONDS(timevalue)`. Numeric. Returns the number of seconds, including fractional seconds, in `timevalue`, which is a number, a time expression, or a time format variable.

#### **Example**

```
DATA LIST FREE (" ")
  /StartDate (ADATE12) EndDate (ADATE12)
  StartDateTIme (DATETIME20) EndDateTIme (DATETIME20)
  StartTime (TIME10) EndTime (TIME10).
BEGIN DATA
3/01/2003, 4/10/2003
01-MAR-2003 12:00, 02-MAR-2003 12:00
09:30, 10:15
END DATA.
COMPUTE days = CTIME.DAYS(EndDate-StartDate).
COMPUTE hours = CTIME.HOURS(EndDateTIme-StartDateTIme).
COMPUTE minutes = CTIME.MINUTES(EndTime-StartTime).
```

- `CTIME.DAYS` calculates the difference between *EndDate* and *StartDate* in days—in this example, 40 days.
- `CTIME.HOURS` calculates the difference between *EndDateTIme* and *StartDateTIme* in hours—in this example, 24 hours.
- `CTIME.MINUTES` calculates the difference between *EndTime* and *StartTime* in minutes—in this example, 45 minutes.

**YRMODA function**

**YRMODA(arg list)**      *Convert year, month, and day to a day number.* The number returned is the number of days since October 14, 1582 (day 0 of the Gregorian calendar).

- Arguments for YRMODA can be variables, constants, or any other type of numeric expression but must yield integers.
- Year, month, and day must be specified in that order.
- The first argument can be any year between 0 and 99, or between 1582 to 47516.
- If the first argument yields a number between 00 and 99, 1900 through 1999 is assumed.
- The month can range from 1 through 13. Month 13 with day 0 yields the last day of the year. For example, YRMODA(1990, 13, 0) produces the day number for December 31, 1990. Month 13 with any other day yields the day of the first month of the coming year—for example, YRMODA(1990, 13, 1) produces the day number for January 1, 1991.
- The day can range from 0 through 31. Day 0 is the last day of the previous month regardless of whether it is 28, 29, 30, or 31. For example, YRMODA(1990, 3, 0) yields 148791.00, the day number for February 28, 1990.
- The function returns the system-missing value if any of the three arguments is missing or if the arguments do not form a valid date *after* October 14, 1582.
- Since YRMODA yields the number of days instead of seconds, you can not display it in date format unless you convert it to the number of seconds.

**Extraction functions**

The extraction functions extract subfields from dates or time intervals, targeting the day or a time from a date value. This permits you to classify events by day of the week, season, shift, and so forth.

Each extraction function begins with XDATE, followed by a period, the subfunction name (what you want to extract), and an argument.

**XDATE.DATE.** XDATE.DATE(datevalue). Numeric. Returns the date portion from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date. To display the result as a date, apply a date format to the variable.

**XDATE.HOUR.** XDATE.HOUR(datetime). Numeric. Returns the hour (an integer between 0 and 23) from a value that represents a time or a datetime. The argument can be a number, a time or datetime variable or an expression that resolves to a time or datetime value.

**XDATE.JDAY.** XDATE.JDAY(datevalue). Numeric. Returns the day of the year (an integer between 1 and 366) from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date.

**XDATE.MDAY.** XDATE.MDAY(datevalue). Numeric. Returns the day of the month (an integer between 1 and 31) from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date.

**XDATE.MINUTE.** XDATE.MINUTE(datetime). Numeric. Returns the minute (an integer between 0 and 59) from a value that represents a time or a datetime. The argument can be a number, a time or datetime variable, or an expression that resolves to a time or datetime value.

**XDATE.MONTH.** XDATE.MONTH(datevalue). Numeric. Returns the month (an integer between 1 and 12) from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date.

**XDATE.QUARTER.** XDATE.QUARTER(datevalue). Numeric. Returns the quarter of the year (an integer between 1 and 4) from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date.

**XDATE.SECOND.** XDATE.SECOND(datetime). Numeric. Returns the second (a number between 0 and 60) from a value that represents a time or a datetime. The argument can be a number, a time or datetime variable or an expression that resolves to a time or datetime value.

**XDATE.TDAY.** XDATE.TDAY(timevalue). Numeric. Returns the number of whole days (as an integer) from a numeric value that represents a time interval. The argument can be a number, a time format variable, or an expression that resolves to a time interval.

**XDATE.TIME.** XDATE.TIME(datetime). Numeric. Returns the time portion from a value that represents a time or a datetime. The argument can be a number, a time or datetime variable or an expression that resolves to a time or datetime value. To display the result as a time, apply a time format to the variable.

**XDATE.WEEK.** XDATE.WEEK(datevalue). Numeric. Returns the week number (an integer between 1 and 53) from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date.

**XDATE.WKDAY.** XDATE.WKDAY(datevalue). Numeric. Returns the day-of-week number (an integer between 1, Sunday, and 7, Saturday) from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date.

**XDATE.YEAR.** XDATE.YEAR(datevalue). Numeric. Returns the year (as a four-digit integer) from a numeric value that represents a date. The argument can be a number, a date format variable, or an expression that resolves to a date.

### Example

```
DATA LIST FREE (" , ")
  /StartDateTime (datetime25) .
BEGIN DATA
29-OCT-2003 11:23:02
1 January 1998 1:45:01
21/6/2000 2:55:13
END DATA.
COMPUTE dateonly=XDATE.DATE(StartDateTime) .
FORMATS dateonly(ADATE10) .
COMPUTE hour=XDATE.HOUR(StartDateTime) .
COMPUTE DayofWeek=XDATE.WKDAY(StartDateTime) .
COMPUTE WeekofYear=XDATE.WEEK(StartDateTime) .
COMPUTE quarter=XDATE.QUARTER(StartDateTime) .
```

- The date portion extracted with XDATE.DATE returns a date expressed in seconds; so, FORMATS is used to display the date in a readable date format.

- Day of the week is an integer between 1 (Sunday) and 7 (Saturday).
- Week of the year is an integer between 1 and 53 (January 1–7 = 1).

### **Date differences**

The DATEDIFF function calculates the difference between two date values and returns an integer (with any fraction component truncated) in the specified date/time units. The general form of the expression is

```
DATEDIFF(datetime2, datetime1, "unit").
```

where *datetime2* and *datetime1* are both date or time format variables (or numeric values that represent valid date/time values), and "unit" is one of the following string literal values, enclosed in quotes:

- Years
- Quarters
- Months
- Weeks
- Days
- Hours
- Minutes
- Seconds

### **Example**

```
DATA LIST FREE /date1 date2 (2ADATE10).
BEGIN DATA
1/1/2004 2/1/2005
1/1/2004 2/15/2005
1/30/2004 1/29/2005
END DATA.
COMPUTE years=DATEDIFF(date2, date1, "years").
```

- The result will be the integer portion of the number of years between the two dates, with any fractional component truncated.
- One "year" is defined as the same month and day, one year before or after the second date argument.
- For the first two cases, the result is 1, since in both cases the number of years is greater than or equal to 1 and less than 2.
- For the third case, the result is 0, since the difference is one day short of a year based on the definition of year.

### **Example**

```
DATA LIST FREE /date1 date2 (2ADATE10).
BEGIN DATA
1/1/2004 2/1/2004
1/1/2004 2/15/2004
```

```
1/30/2004 2/1/2004
END DATA.
COMPUTE months=DATEDIFF(date2, date1, "months").
```

- The result will be the integer portion of the number of months between the two dates, with any fractional component truncated.
- One “month” is defined as the same day of the month, one calendar month before or after the second date argument.
- For the first two cases, the result will be 1, since both February 1 and February 15, 2004, are greater than or equal to one month and less than two months after January 1, 2004.
- For the third case, the result will be 0. By definition, *any* date in February 2004 will be less than one month after January 30, 2004, resulting in a value of 0.

### **Date increments**

The DATESUM function calculates a date or time value a specified number of units from a given date or time value. The general form of the function is:

```
DATESUM(datevar, value, "unit", "method").
```

- `datevar` is a date/time format variable (or a numeric value that represents a valid date/time value).
- `value` is a positive or negative number. For variable-length units (years, quarters, months), fractional values are truncated to integers.
- `"unit"` is one of the following string literal values enclosed in quotes: years, quarters, months, weeks, days, hours, minutes, seconds.
- `"method"` is an optional specification for variable-length units (years, quarters, months) enclosed in quotes. The method can be `"rollover"` or `"closest"`. The **rollover** method advances excess days into the next month. The **closest** method uses the closest legitimate date within the month. This is the default.

### **Example**

```
DATA LIST FREE /datevar1 (ADATE10).
BEGIN DATA
2/28/2004
2/29/2004
END DATA.
COMPUTE rollover_year=DATESUM(datevar1, 1, "years", "rollover").
COMPUTE closest_year=DATESUM(datevar1, 1, "years", "closest").
COMPUTE fraction_year=DATESUM(datevar1, 1.5, "years").
FORMATS rollover_year closest_year fraction_year (ADATE10).
SUMMARIZE
  /TABLES=datevar1 rollover_year closest_year fraction_year
  /FORMAT=VALIDLIST NOCASENUM
  /CELLS=NONE.
```

**Figure 2-15**  
Results of rollover and closest year calculations

	datevar1	rollover_year	closest_year	fraction_year
1	02/28/2004	02/28/2005	02/28/2005	08/28/2005
2	02/29/2004	03/01/2005	02/28/2005	08/29/2005

- The rollover and closest methods yield the same result when incrementing February 28, 2004, by one year: February 28, 2005.
- Using the rollover method, incrementing February 29, 2004, by one year returns a value of March 1, 2005. Since there is no February 29, 2005, the excess day is rolled over to the next month.
- Using the closest method, incrementing February 29, 2004, by one year returns a value of February 28, 2005, which is the closest day in the same month of the following year.
- The results for *fraction\_year* are exactly the same as for *closest\_year* because the closest method is used by default, and the value parameter of 1.5 is truncated to 1 for variable-length units.
- All three COMPUTE commands create new variables that display values in the default F format, which for a date value is a large integer. The FORMATS command specifies the ADATE format for the new variables.

### Example

```
DATA LIST FREE /datevar1 (ADATE10).
BEGIN DATA
01/31/2003
01/31/2004
03/31/2004
05/31/2004
END DATA.
COMPUTE rollover_month=DATESUM(datevar1, 1, "months", "rollover").
COMPUTE closest_month=DATESUM(datevar1, 1, "months", "closest").
COMPUTE previous_month_rollover =
  DATESUM(datevar1, -1, "months", "rollover").
COMPUTE previous_month_closest =
  DATESUM(datevar1, -1, "months", "closest").
FORMATS rollover_month closest_month
previous_month_rollover previous_month_closest (ADATE10).
SUMMARIZE
  /TABLES=datevar1 rollover_month closest_month
  previous_month_rollover previous_month_closest
  /FORMAT=VALIDLIST NOCASENUM
  /CELLS=NONE.
```

**Figure 2-16**  
Results of month calculations

	datevar1	rollover_month	closest_month	previous_ month_rollover	previous_ month_closest
1	01/31/2003	03/03/2003	02/28/2003	12/31/2002	12/31/2002
2	01/31/2004	03/02/2004	02/29/2004	12/31/2003	12/31/2003
3	03/31/2004	05/01/2004	04/30/2004	03/02/2004	02/29/2004
4	05/31/2004	07/01/2004	06/30/2004	05/01/2004	04/30/2004

- Using the rollover method, incrementing by one month from January 31 yields a date in March, since February has a maximum of 29 days; and incrementing one month from March 31 and May 31 yields May 1 and July 1, respectively, since April and June each have only 30 days.
- Using the closest method, incrementing by one month from the last day of any month will always yield the closest valid date within the next month. For example, in a nonleap year, one month after January 31 is February 28, and one month after February 28 is March 28.
- Using the rollover method, decrementing by one month (by specifying a negative value parameter) from the last day of a month may sometimes yield unexpected results, since the excess days are rolled back to the original month. For example, one month prior to March 31 yields March 3 for nonleap years and March 2 for leap years.
- Using the closest method, decrementing by one month from the last day of the month will always yield the closest valid date within the previous month. For example, one month prior to April 30, is March 30 (not March 31), and one month prior to March 31 is February 28 in nonleap years and February 29 in leap years.

## **String expressions**

Expressions involving string variables can be used on `COMPUTE` and `IF` commands and in logical expressions on commands such as `IF`, `DO IF`, `LOOP IF`, and `SELECT IF`.

- A string expression can be a constant enclosed in quotes (for example, `'IL'`), a string function, or a string variable. [For more information, see the topic String functions on p. 106.](#)
- An expression must return a string if the target variable is a string.
- The string returned by a string expression does not have to be the same length as the target variable; no warning messages are issued if the lengths are not the same. If the target variable produced by a `COMPUTE` command is shorter, the result is right-trimmed. If the target variable is longer, the result is right-padded.

## **String functions**

- The target variable for each string function must be a string and *must have already been declared* (see `STRING`).
- Multiple arguments in a list must be separated by commas.
- When two strings are compared, the case in which they are entered is significant. The `LOWER` and `UPCASE` functions are useful for making comparisons of strings regardless of case.
- String functions that include a byte position or count argument or return a byte position or count may return different results in Unicode mode than in code page mode. For example, `é` is one byte in code page mode but is two bytes in Unicode mode; so `résumé` is six bytes in code page mode and eight bytes in Unicode mode.
- In Unicode mode, trailing blanks are always removed from the values of string variables in string functions unless explicitly preserved with the `NTRIM` function.
- In code page mode, trailing blanks are always preserved in the values of string variables unless explicitly removed with the `RTRIM` function.

For more information on Unicode mode, see [UNICODE Subcommand](#).



**CHAR.INDEX.** CHAR.INDEX(haystack, needle[, divisor]). Numeric. Returns a number indicating the character position of the first occurrence of needle in haystack. The optional third argument, divisor, is a number of characters used to divide needle into separate strings. Each substring is used for searching and the function returns the first occurrence of any of the substrings. For example, CHAR.INDEX(var1, 'abcd') will return the value of the starting position of the complete string "abcd" in the string variable var1; CHAR.INDEX(var1, 'abcd', 1) will return the value of the position of the first occurrence of any of the values in the string; and CHAR.INDEX(var1, 'abcd', 2) will return the value of the first occurrence of either "ab" or "cd". Divisor must be a positive integer and must divide evenly into the length of needle. Returns 0 if needle does not occur within haystack.

**CHAR.LENGTH.** CHAR.LENGTH(strexpr). Numeric. Returns the length of strexpr in characters, with any trailing blanks removed.

**CHAR.LPAD.** CHAR.LPAD(strexpr1,length[,strexpr2]). String. Left-pads strexpr1 to make its length the value specified by length using as many complete copies as will fit of strexpr2 as the padding string. The value of length represents the number of characters and must be a positive integer. If the optional argument strexpr2 is omitted, the value is padded with blank spaces.

**CHAR.MBLEN.** CHAR.MBLEN(strexpr,pos). Numeric. Returns the number of bytes in the character at character position pos of strexpr.

**CHAR.RINDEX.** CHAR.RINDEX(haystack,needle[,divisor]). Numeric. Returns an integer that indicates the starting character position of the last occurrence of the string needle in the string haystack. The optional third argument, divisor, is the number of characters used to divide needle into separate strings. For example, CHAR.RINDEX(var1, 'abcd') will return the starting position of the last occurrence of the entire string "abcd" in the variable var1; CHAR.RINDEX(var1, 'abcd', 1) will return the value of the position of the last occurrence of any of the values in the string; and CHAR.RINDEX(var1, 'abcd', 2) will return the value of the starting position of the last occurrence of either "ab" or "cd". Divisor must be a positive integer and must divide evenly into the length of needle. If needle is not found, the value 0 is returned.

**CHAR.RPAD.** CHAR.RPAD(strexpr1,length[,strexpr2]). String. Right-pads strexpr1 with strexpr2 to extend it to the length given by length using as many complete copies as will fit of strexpr2 as the padding string. The value of length represents the number of characters and must be a positive integer. The optional third argument strexpr2 is a quoted string or an expression that resolves to a string. If strexpr2 is omitted, the value is padded with blanks.

**CHAR.SUBSTR.** CHAR.SUBSTR(strexpr,pos[,length]). String. Returns the substring beginning at character position pos of strexpr. The optional third argument represents the number of characters in the substring. If the optional argument length is omitted, returns the substring beginning at character position pos of strexpr and running to the end of strexpr. For example CHAR.SUBSTR('abcd', 2) returns 'bcd' and CHAR.SUBSTR('abcd', 2, 2) returns 'bc'. (Note: Use the SUBSTR function instead of CHAR.SUBSTR if you want to use the function on the left side of an equals sign to replace a substring.)

**CONCAT.** CONCAT(strexpr,strexpr[,...]). String. Returns a string that is the concatenation of all its arguments, which must evaluate to strings. This function requires two or more arguments. In code page mode, if strexpr is a string variable, use RTRIM if you only want the actual string value without the right-padding to the defined variable width. For example, CONCAT(RTRIM(stringvar1), RTRIM(stringvar2)).

**LENGTH.** LENGTH(strexpr). Numeric. Returns the length of strexpr in bytes, which must be a string expression. For string variables, in Unicode mode this is the number of bytes in each value, excluding trailing blanks, but in code page mode this is the defined variable length, including trailing blanks. To get the length (in bytes) without trailing blanks in code page mode, use LENGTH(RTRIM(strexpr)).

**LOWER.** LOWER(strexpr). String. Returns strexpr with uppercase letters changed to lowercase and other characters unchanged. The argument can be a string variable or a value. For example, LOWER(name1) returns charles if the value of name1 is Charles.

**LTRIM.** LTRIM(strexpr[,char]). String. Returns strexpr with any leading instances of char removed. If char is not specified, leading blanks are removed. Char must resolve to a single character.

**MAX.** MAX(value,value[,...]). Numeric or string. Returns the maximum value of its arguments that have valid values. This function requires two or more arguments. For numeric values, you can specify a minimum number of valid arguments for this function to be evaluated.

**MIN.** MIN(value,value[,...]). Numeric or string. Returns the minimum value of its arguments that have valid, nonmissing values. This function requires two or more arguments. For numeric values, you can specify a minimum number of valid arguments for this function to be evaluated.

**MBLEN.BYTE.** MBLEN.BYTE(strexpr,pos). Numeric. Returns the number of bytes in the character at byte position pos of strexpr.

**NORMALIZE.** NORMALIZE(strexpr). String. Returns the normalized version of strexp. In Unicode mode, it returns Unicode NFC. In code page mode, it has no effect and returns strexp unmodified. The length of the result may be different from the length of the input.

**NTRIM.** NTRIM(varname). Returns the value of varname, without removing trailing blanks. The value of varname must be a variable name; it cannot be an expression.

**REPLACE.** REPLACE(a1, a2, a3[, a4]). String. In a1, instances of a2 are replaced with a3. The optional argument a4 specifies the number of occurrences to replace; if a4 is omitted, all occurrences are replaced. Arguments a1, a2, and a3 must resolve to string values (literal strings enclosed in quotes or string variables), and the optional argument a4 must resolve to a non-negative integer. For example, REPLACE("abcabc", "a", "x") returns a value of "xbcxbc" and REPLACE("abcabc", "a", "x", 1) returns a value of "xbcab".

**RTRIM.** RTRIM(strexpr[,char]). String. Trims trailing instances of char within strexpr. The optional second argument char is a single quoted character or an expression that yields a single character. If char is omitted, trailing blanks are trimmed.

**STRUNC.** STRUNC(strexpr, length). String. Returns strexp truncated to length (in bytes) and then trimmed of any trailing blanks. Truncation removes any fragment of a character that would be truncated.

**UPCASE.** UPCASE(strexpr). String. Returns strexpr with lowercase letters changed to uppercase and other characters unchanged.

### Deprecated string functions

The following functions provide functionality similar to the newer CHAR functions, but they operate at the byte level rather than the character level. For example, the INDEX function returns the byte position of *needle* within *haystack*, whereas CHAR.INDEX returns the character position. These functions are supported primarily for compatibility with previous releases.

**INDEX.** INDEX(haystack,needle[,divisor]). Numeric. Returns a number that indicates the byte position of the first occurrence of *needle* in *haystack*. The optional third argument, *divisor*, is a number of bytes used to divide *needle* into separate strings. Each substring is used for searching and the function returns the first occurrence of any of the substrings. *Divisor* must be a positive integer and must divide evenly into the length of *needle*. Returns 0 if *needle* does not occur within *haystack*.

**LPAD.** LPAD(strexpr1,length[,strexpr2]). String. Left-pads *strexpr1* to make its length the value specified by *length* using as many complete copies as will fit of *strexpr2* as the padding string. The value of *length* represents the number of bytes and must be a positive integer. If the optional argument *strexpr2* is omitted, the value is padded with blank spaces.

**RINDEX.** RINDEX(haystack,needle[,divisor]). Numeric. Returns an integer that indicates the starting byte position of the last occurrence of the string *needle* in the string *haystack*. The optional third argument, *divisor*, is the number of bytes used to divide *needle* into separate strings. *Divisor* must be a positive integer and must divide evenly into the length of *needle*. If *needle* is not found, the value 0 is returned.

**RPAD.** RPAD(strexpr1,length[,strexpr2]). String. Right-pads *strexpr1* with *strexpr2* to extend it to the length given by *length* using as many complete copies as will fit of *strexpr2* as the padding string. The value of *length* represents the number of bytes and must be a positive integer. The optional third argument *strexpr2* is a quoted string or an expression that resolves to a string. If *strexpr2* is omitted, the value is padded with blanks.

**SUBSTR.** SUBSTR(strexpr,pos[,length]). String. Returns the substring beginning at byte position *pos* of *strexpr*. The optional third argument represents the number of bytes in the substring. If the optional argument *length* is omitted, returns the substring beginning at byte position *pos* of *strexpr* and running to the end of *strexpr*. When used on the left side of an equals sign, the substring is replaced by the string specified on the right side of the equals sign. The rest of the original string remains intact. For example, SUBSTR(ALPHA6,3,1)='\*' changes the third character of all values for *ALPHA6* to \*. If the replacement string is longer or shorter than the substring, the replacement is truncated or padded with blanks on the right to an equal length.

### Example

```
STRING stringVar1 stringVar2 stringVar3 (A22).
COMPUTE stringVar1=' Does this'.
COMPUTE stringVar2='ting work?'.
COMPUTE stringVar3=
    CONCAT(RTRIM(LTRIM(stringVar1)), " ",
    REPLACE(stringVar2, "ting", "thing")).
```

- The CONCAT function concatenates the values of *stringVar1* and *stringVar2*, inserting a space as a literal string (" ") between them.

- The `RTRIM` function strips off trailing blanks from *stringVar1*. In code page mode, this is necessary to eliminate excessive space between the two concatenated string values because in code page mode all string variable values are automatically right-padded to the defined width of the string variables. In Unicode mode, this has no effect because trailing blanks are automatically removed from string variable values in Unicode mode.
- The `LTRIM` function removes the leading spaces from the beginning of the value of *stringVar1*.
- The `REPLACE` function replaces the misspelled "ting" with "thing" in *stringVar2*.

The final result is a string value of “*Does this thing work?*”

### Example

This example extracts the numeric components from a string telephone number into three numeric variables.

```
DATA LIST FREE (" ") /telephone (A16).
BEGIN DATA
111-222-3333
222 - 333 - 4444
333-444-5555
444 - 555-6666
555-666-0707
END DATA.
STRING #telstr(A16).
COMPUTE #telstr = telephone.
VECTOR tel(3,f4).
LOOP #i = 1 to 2.
- COMPUTE #dash = CHAR.INDEX(#telstr,"-").
- COMPUTE tel(#i) = NUMBER(CHAR.SUBSTR(#telstr,1,#dash-1),f10).
- COMPUTE #telstr = CHAR.SUBSTR(#telstr,#dash+1).
END LOOP.
COMPUTE tel(3) = NUMBER(#telstr,f10).
EXECUTE.
FORMATS tel1 tel2 (N3) tel3 (N4).
```

- A temporary (scratch) string variable, *#telstr*, is declared and set to the value of the original string telephone number.
- The `VECTOR` command creates three numeric variables—*tel1*, *tel2*, and *tel3*—and creates a vector containing those variables.
- The `LOOP` structure iterates twice to produce the values for *tel1* and *tel2*.
- `COMPUTE #dash = CHAR.INDEX(#telstr,"-")` creates another temporary variable, *#dash*, that contains the position of the first dash in the string value.
- On the first iteration, `COMPUTE tel(#i) = NUMBER(CHAR.SUBSTR(#telstr,1,#dash-1),f10)` extracts everything prior to the first dash, converts it to a number, and sets *tel1* to that value.
- `COMPUTE #telstr = CHAR.SUBSTR(#telstr,#dash+1)` then sets *#telstr* to the remaining portion of the string value after the first dash.
- On the second iteration, `COMPUTE #dash...` sets *#dash* to the position of the “first” dash in the modified value of *#telstr*. Since the area code and the original first dash have been removed from *#telstr*, this is the position of the dash between the exchange and the number.

- `COMPUTE tel(#)`... sets `tel2` to the numeric value of everything up to the “first” dash in the modified version of `#telstr`, which is everything after the first dash and before the second dash in the original string value.
- `COMPUTE #telstr`... then sets `#telstr` to the remaining segment of the string value—everything after the “first” dash in the modified value, which is everything after the second dash in the original value.
- After the two loop iterations are complete, `COMPUTE tel(3) = NUMBER(#telstr,f10)` sets `tel3` to the numeric value of the final segment of the original string value.

### ***String/numeric conversion functions***

**NUMBER.** `NUMBER(strexpr,format)`. Numeric. Returns the value of the string expression `strexpr` as a number. The second argument, `format`, is the numeric format used to read `strexpr`. For example, `NUMBER(stringDate,DATE11)` converts strings containing dates of the general format `dd-mmm-yyyy` to a numeric number of seconds that represent that date. (To display the value as a date, use the `FORMATS` or `PRINT FORMATS` command.) If the string cannot be read using the format, this function returns `system-missing`.

**STRING.** `STRING(numexpr,format)`. String. Returns the string that results when `numexpr` is converted to a string according to `format`. `STRING(-1.5,F5.2)` returns the string value `'-1.50'`. The second argument `format` must be a format for writing a numeric value.

#### ***Example***

```
DATA LIST FREE /tel1 tel2 tel3.
BEGIN DATA
123 456 0708
END DATA.
STRING telephone (A12).
COMPUTE telephone=
  CONCAT(STRING(tel1,N3), "-", STRING(tel2, N3), "-", STRING(tel3, N4)).
```

- A new string variable, `telephone`, is declared to contain the computed string value.
- The three numeric variables are converted to strings and concatenated with dashes between the values.
- The numeric values are converted using `N` format to preserve any leading zeros.

### ***LAG function***

**LAG.** `LAG(variable[, n])`. Numeric or string. The value of `variable` in the previous case or `n` cases before. The optional second argument, `n`, must be a positive integer; the default is 1. For example, `prev4=LAG(gnp,4)` returns the value of `gnp` for the fourth case before the current one. The first four cases have `system-missing` values for `prev4`.

- The result is of the same type (numeric or string) as the variable specified as the first argument.

- The first  $n$  cases for string variables are set to blanks. For example, if `PREV2=LAG(LNAME, 2)` is specified, blanks will be assigned to the first two cases for `PREV2`.
- When `LAG` is used with commands that select cases (for example, `SELECT IF` and `SAMPLE`), `LAG` counts cases *after* case selection, even if specified before these commands. [For more information, see the topic Command Order on p. 41.](#)

*Note:* In a series of transformation commands without any intervening `EXECUTE` commands or other commands that read the data, lag functions are calculated after all other transformations, regardless of command order. For example,

```
COMPUTE lagvar=LAG(var1).
COMPUTE var1=var1*2.
```

and

```
COMPUTE lagvar=LAG(var1).
EXECUTE.
COMPUTE var1=var1*2.
```

yield very different results for the value of `lagvar`, since the former uses the transformed value of `var1` while the latter uses the original value.

## VALUELABEL function

**VALUELABEL.** VALUELABEL(varname). String. Returns the value label for the value of variable or an empty string if there is no label for the value. The value of `varname` must be a variable name; it cannot be an expression.

### Example

```
STRING labelvar (A120).
COMPUTE labelvar=VALUELABEL(var1).
DO REPEAT varlist=var2, var3, var4
    /newvars=labelvar2, labelvar3, labelvar4.
- STRING newvars (A120).
- COMPUTE newvars=VALUELABEL(varlist).
END REPEAT.
```

## Logical expressions

Logical expressions can appear on the `COMPUTE`, `IF`, `SELECT IF`, `DO IF`, `ELSE IF`, `LOOP IF`, and `END LOOP IF` commands. A logical expression is evaluated as true or false, or as missing if it is indeterminate. A logical expression returns 1 if the expression is true, 0 if it is false, or system-missing if it is missing. Thus, logical expressions can be any expressions that yield this three-value logic.

- The simplest logical expression is a logical variable. A logical variable is any numeric variable that has the values 1, 0, or system-missing. Logical variables cannot be strings.
- Logical expressions can be simple logical variables or relations, or they can be complex logical tests involving variables, constants, functions, relational operators, logical operators, and parentheses to control the order of evaluation.

- On an `IF` command, a logical expression that is true causes the assignment expression to be executed. A logical expression that returns missing has the same effect as one that is false—that is, the assignment expression is not executed and the value of the target variable is not altered.
- On a `DO IF` command, a logical expression that is true causes the execution of the commands immediately following the `DO IF`, up to the next `ELSE IF`, `ELSE`, or `END IF`. If it is false, the next `ELSE IF` or `ELSE` command is evaluated. If the logical expression returns missing for each of these, the entire structure is skipped.
- On a `SELECT IF` command, a logical expression that is true causes the case to be selected. A logical expression that returns missing has the same effect as one that is false—that is, the case is not selected.
- On a `LOOP IF` command, a logical expression that is true causes looping to begin (or continue). A logical expression that returns missing has the same effect as one that is false—that is, the structure is skipped.
- On an `END LOOP IF` command, a logical expression that is false returns control to the `LOOP` command for that structure, and looping continues. If it is true, looping stops and the structure is terminated. A logical expression that returns a missing value has the same effect as one that is true—that is, the structure is terminated.

### Example

```
DATA LIST FREE (" ") /a.
BEGIN DATA
1, , 1, ,
END DATA.
COMPUTE b=a.
* The following does NOT work since the second condition
  is never evaluated.
DO IF a=1.
COMPUTE a1=1.
ELSE IF MISSING(a).
COMPUTE a1=2.
END IF.
* On the other hand the following works.
DO IF MISSING(b).
COMPUTE b1=2.
ELSE IF b=1.
COMPUTE b1=1.
END IF.
```

- The first `DO IF` will never yield a value of 2 for `a1` because if `a` is missing, then `DO IF a=1` evaluates as missing and control passes immediately to `END IF`. So `a1` will either be 1 or missing.
- In the second `DO IF`, however, we take care of the missing condition first; so if the value of `b` is missing, `DO IF MISSING(b)` evaluates as true and `b1` is set to 2; otherwise, `b1` is set to 1.

### String variables in logical expressions

String variables, like numeric variables, can be tested in logical expressions.

- String variables must be declared before they can be used in a string expression.

- String variables cannot be compared to numeric variables.
- If strings of different lengths are compared, the shorter string is right-padded with blanks to equal the length of the longer string.
- The magnitude of strings can be compared using `LT`, `GT`, and so on, but the outcome depends on the sorting sequence of the computer. Use with caution.
- User-missing string values are treated the same as nonmissing string values when evaluating string variables in logical expressions. In other words, all string variable values are treated as valid, nonmissing values in logical expressions.

### **Relational operators**

A relation is a logical expression that compares two values using a *relational operator*. In the command

```
IF (X EQ 0) Y=1
```

the variable `X` and `0` are expressions that yield the values to be compared by the `EQ` relational operator. The following are the relational operators:

<b>Symbol</b>	<b>Definition</b>
EQ or =	Equal to
NE or ~= or != or <>	Not equal to
LT or <	Less than
LE or <=	Less than or equal to
GT or >	Greater than
GE or >=	Greater than or equal to

- The expressions in a relation can be variables, constants, or more complicated arithmetic expressions.
- Blanks (not commas) must be used to separate the relational operator from the expressions. To make the command more readable, use extra blanks or parentheses.
- For string values, “less than” and “greater than” results can vary by locale even for the same set of characters, since the national collating sequence is used. Language order, not ASCII order, determines where certain characters fall in the sequence.

### **NOT logical operator**

The `NOT` logical operator reverses the true/false outcome of the expression that immediately follows.

- The `NOT` operator affects only the expression that immediately follows, unless a more complex logical expression is enclosed in parentheses.
- You can substitute `~` or `¬` for `NOT` as a logical operator.
- `NOT` can be used to check whether a numeric variable has the value 0, 1, or any other value. For example, all scratch variables are initialized to 0. Therefore, `NOT (#ID)` returns false or missing when `#ID` has been assigned a value other than 0.



## AND and OR logical operators

Two or more relations can be logically joined using the logical operators AND and OR. Logical operators combine relations according to the following rules:

- The ampersand (&) symbol is a valid substitute for the logical operator AND. The vertical bar ( | ) is a valid substitute for the logical operator OR.
- Only one logical operator can be used to combine two relations. However, multiple relations can be combined into a complex logical expression.
- Regardless of the number of relations and logical operators used to build a logical expression, the result is either true, false, or indeterminate because of missing values.
- Operators or expressions cannot be implied. For example, `X EQ 1 OR 2` is illegal; you must specify `X EQ 1 OR X EQ 2`.
- The ANY and RANGE functions can be used to simplify complex expressions.

**AND**            *Both relations must be true for the complex expression to be true.*

**OR**             *If either relation is true, the complex expression is true.*

The following table lists the outcomes for AND and OR combinations.

Table 2-3

*Logical outcomes*

Expression	Outcome	Expression	Outcome
true AND true	= true	true OR true	= true
true AND false	= false	true OR false	= true
false AND false	= false	false OR false	= false
true AND missing	= missing	true OR missing	= true*
missing AND missing	= missing	missing OR missing	= missing
false AND missing	= false*	false OR missing	= missing

\* Expressions where the outcome can be evaluated with incomplete information. [For more information, see the topic Missing values in logical expressions on p. 121.](#)

### Example

```
DATA LIST FREE /var1 var2 var3.
BEGIN DATA
1 1 1
1 2 1
1 2 3
4 2 4
END DATA.
SELECT IF var1 = 4 OR ((var2 > var1) AND (var1 <> var3)).
```

- Any case that meets the first condition—`var1 = 4`—will be selected, which in this example is only the last case.
- Any case that meets the second condition will also be selected. In this example, only the third case meets this condition, which contains two criteria: `var2` is greater than `var1` and `var1` is not equal to `var3`.

**Order of evaluation**

- When arithmetic operators and functions are used in a logical expression, the order of operations is functions and arithmetic operations first, then relational operators, and then logical operators.
- When more than one logical operator is used, NOT is evaluated first, then AND, and then OR.
- To change the order of evaluation, use parentheses.

**Logical functions**

- Each argument to a logical function (expression, variable name, or constant) must be separated by a comma.
- The target variable for a logical function must be numeric.
- The functions RANGE and ANY can be useful shortcuts to more complicated specifications on the IF, DO IF, and other conditional commands. For example, for non-missing values, the command

```
SELECT IF ANY (REGION, "NW" , "NE" , "SE" ) .
```

is equivalent to

```
SELECT IF ( REGION EQ "NW" OR REGION EQ "NE" OR REGION EQ "SE" ) .
```

**RANGE.** RANGE(test,lo,hi[,lo,hi,..]). Logical. Returns 1 or true if test is within any of the inclusive range(s) defined by the pairs lo, hi. Arguments must be all numeric or all strings of the same length, and each of the lo, hi pairs must be ordered with lo <= hi. Note: For string values, results can vary by locale even for the same set of characters, since the national collating sequence is used. Language order, not ASCII order, determines where certain characters fall in the sequence.

**ANY.** ANY(test,value[,value,...]). Logical. Returns 1 or true if the value of test matches any of the subsequent values; returns 0 or false otherwise. This function requires two or more arguments. For example, ANY(var1, 1, 3, 5) returns 1 if the value of var1 is 1, 3, or 5 and 0 for other values. ANY can also be used to scan a list of variables or expressions for a value. For example, ANY(1, var1, var2, var3) returns 1 if any of the three specified variables has a value of 1 and 0 if all three variables have values other than 1.

See [Treatment of missing values in arguments](#) for information on how missing values are handled by the ANY and RANGE functions.

**Scoring expressions**

Scoring expressions apply model XML from an external file to the active dataset and generate predicted values, predicted probabilities, and other values based on that model.

- Scoring expressions must be preceded by a MODEL HANDLE command that identifies the external XML model file and optionally does variable mapping.

- Scoring expressions require two arguments: the first identifies the model, and the second identifies the scoring function. An optional third argument allows users to obtain the probability (for each case) associated with a selected category, in the case of a categorical target variable. It is also used in nearest neighbor models to specify a particular neighbor.
- Prior to applying scoring functions to a set of data, a data validation analysis is performed. The analysis includes checking that data are of the correct type as well as checking that the data values are in the set of allowed values defined in the model. For example, for categorical variables, a value that is neither a valid category nor defined as user-missing would be treated as an invalid value. Values that are found to be invalid are treated as system-missing.

The following scoring expressions are available:

**ApplyModel.** ApplyModel(handle, "function", value). Numeric. Applies a particular scoring function to the input case data using the model specified by handle and where "function" is one of the following string literal values enclosed in quotes: predict, stddev, probability, confidence, nodeid, cumhazard, neighbor, distance. The model handle is the name associated with the external XML file, as defined on the MODEL HANDLE command. The optional third argument applies when the function is "probability", "neighbor", or "distance". For "probability", it specifies a category for which the probability is calculated. For "neighbor" and "distance", it specifies a particular neighbor (as an integer) for nearest neighbor models. ApplyModel returns system-missing if a value can not be computed. String values must be enclosed in quotes. For example, ApplyModel(name1, 'probability', 'reject'), where name1 is the model's handle name and 'reject' is a valid category for a target variable that is a string.

**StrApplyModel.** StrApplyModel(handle, "function", value). String. Applies a particular scoring function to the input case data using the model specified by handle and where "function" is one of the following string literal values enclosed in quotes: predict, stddev, probability, confidence, nodeid, cumhazard, neighbor, distance. The model handle is the name associated with the external XML file, as defined on the MODEL HANDLE command. The optional third argument applies when the function is "probability", "neighbor", or "distance". For "probability", it specifies a category for which the probability is calculated. For "neighbor" and "distance", it specifies a particular neighbor (as an integer) for nearest neighbor models. StrApplyModel returns a blank string if a value cannot be computed.

The following scoring functions are available:

<b>PREDICT</b>	<i>Returns the predicted value of the target variable.</i>
<b>STDDEV</b>	<i>Standard deviation.</i>
<b>PROBABILITY</b>	<i>Probability associated with a particular category of a target variable. Applies only to categorical variables. In the absence of the optional third parameter, category, this is the probability that the predicted category is the correct one for the target variable. If a particular category is specified, then this is the probability that the specified category is the correct one for the target variable.</i>
<b>CONFIDENCE</b>	<i>A probability measure associated with the predicted value of a categorical target variable. Applies only to categorical variables.</i>
<b>NODEID</b>	<i>The terminal node number. Applies only to tree models.</i>
<b>CUMHAZARD</b>	<i>Cumulative hazard value. Applies only to Cox regression models.</i>

**NEIGHBOR**

*The ID of the kth nearest neighbor:* Applies only to nearest neighbor models. In the absence of the optional third parameter, k, this is the ID of the nearest neighbor. The ID is the value of the case labels variable, if supplied, and otherwise the case number.

**DISTANCE**

*The distance to the kth nearest neighbor:* Applies only to nearest neighbor models. In the absence of the optional third parameter, k, this is the distance to the nearest neighbor. Depending on the model, either Euclidean or City Block distance will be used.

The following table lists the set of scoring functions available for each type of model that supports scoring. The function type denoted as PROBABILITY (category) refers to specification of a particular category (the optional third parameter) for the PROBABILITY function.

Table 2-4  
Supported functions by model type

Model type	Supported functions
Tree (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE, NODEID
Tree (scale target)	PREDICT, NODEID, STDDEV
Boosted Tree (C5.0)	PREDICT, CONFIDENCE
Linear Regression	PREDICT, STDDEV
Automatic Linear Models	PREDICT
Binary Logistic Regression	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Conditional Logistic Regression	PREDICT
Multinomial Logistic Regression	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
General Linear Model	PREDICT, STDDEV
Discriminant	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
TwoStep Cluster	PREDICT
K-Means Cluster	PREDICT, CONFIDENCE
Kohonen	PREDICT
Neural Net (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Neural Net (scale target)	PREDICT
Naive Bayes	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Anomaly Detection	PREDICT
Ruleset	PREDICT, CONFIDENCE
Generalized Linear Model (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Generalized Linear Model (scale target)	PREDICT, STDDEV
Generalized Linear Mixed Model (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Generalized Linear Mixed Model (scale target)	PREDICT
Ordinal Multinomial Regression	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Cox Regression	PREDICT, CUMHAZARD

Model type	Supported functions
Nearest Neighbor (scale target)	PREDICT, STDDEV, NEIGHBOR, NEIGHBOR(K), DISTANCE, DISTANCE(K)
Nearest Neighbor (categorical target)	PREDICT, PROBABILITY, CONFIDENCE, NEIGHBOR, NEIGHBOR(K), DISTANCE, DISTANCE(K)

- For the Binary Logistic Regression, Multinomial Logistic Regression, and Naive Bayes models, the value returned by the CONFIDENCE function is identical to that returned by the PROBABILITY function.
- For the K-Means model, the value returned by the CONFIDENCE function is the least distance.
- For tree and ruleset models, the confidence can be interpreted as an adjusted probability of the predicted category and is always less than the value given by PROBABILITY. For these models, the confidence value is more reliable than the value given by PROBABILITY.
- For neural network models, the confidence provides a measure of whether the predicted category is much more likely than the second-best predicted category.
- For Ordinal Multinomial Regression and Generalized Linear Model, the PROBABILITY function is supported when the target variable is binary.
- For nearest neighbor models without a target variable, the available functions are NEIGHBOR and DISTANCE.

## Missing values

Functions and simple arithmetic expressions treat missing values in different ways. In the expression

```
(var1+var2+var3)/3
```

the result is missing if a case has a missing value for any of the three variables.

In the expression

```
MEAN(var1, var2, var3)
```

the result is missing only if the case has missing values for all three variables.

For statistical functions, you can specify the minimum number of arguments that must have nonmissing values. To do so, type a period and the minimum number after the function name, as in:

```
MEAN.2(var1, var2, var3)
```

The following sections contain more information on the treatment of missing values in functions and transformation expressions, including special missing value functions.

### Treatment of missing values in arguments

If the logic of an expression is indeterminate because of missing values, the expression returns a missing value and the command is not executed. The following table summarizes how missing values are handled in arguments to various functions.

**Table 2-5**  
Missing values in arguments

Function	Returns system-missing if
MOD (x1, x2)	x1 is missing, or x2 is missing and x1 is not 0.
MAX.n (x1, x2, . . . xk) MEAN.n (x1, x2, . . . xk) MIN.n (x1, x2, . . . x1) SUM.n (x1, x2, . . . xk)	Fewer than <i>n</i> arguments are valid; the default <i>n</i> is 1.
CFVAR.n (x1, x2, . . . xk) SD.n (x1, x2, . . . xk) VARIANCE.n (x1, x2, . . . xk)	Fewer than <i>n</i> arguments are valid; the default <i>n</i> is 2.
LPAD (x1, x2, x3) LTRIM (x1, x2) RTRIM (x1, x2) RPAD (x1, x2, x3)	x1 or x2 is illegal or missing.
SUBSTR (x1, x2, x3)	x2 or x3 is illegal or missing.
NUMBER (x, format) STRING (x, format)	The conversion is invalid.
INDEX (x1, x2, x3) RINDEX (x1, x2, x3)	x3 is invalid or missing.
LAG (x, n)	x is missing <i>n</i> cases previously (and always for the first <i>n</i> cases); the default <i>n</i> is 1.
ANY (x, x1, x2, . . . xk)	For numeric values, if x is missing or all the remaining arguments are missing, the result is system-missing. For string values, user-missing value are treated as valid values, and the result is never missing.
RANGE (x, x1, x2, . . . xk1, xk2)	For numeric values, the result is system-missing if: <ul style="list-style-type: none"> <li>■ x is missing, or</li> <li>■ all the ranges defined by the remaining arguments are missing, or</li> <li>■ any range has a starting value that is higher than the ending value.</li> </ul> <p>A numeric range is missing if either of the arguments that define the range is missing. This includes ranges for which one of the arguments is equal to the value of the first argument in the expression. For example: RANGE (x, x1, x2) is missing if any of the arguments is missing, even if x1 or x2 is equal to x.</p> <p>For string values, user-missing values are treated as valid values, and the result is only missing if any range has a starting value that is higher than the ending value.</p>
VALUE (x)	x is system-missing.
MISSING (x) NMISS (x1, x2, . . . xk) NVALID (x1, x2, . . . xk) SYSMIS (x)	Never.

- Any function that is not listed in this table returns the system-missing value when the argument is missing.
- The system-missing value is displayed as a period (.) for numeric variables.
- String variables do not have system-missing values. An invalid string expression nested within a complex transformation yields a null string, which is passed to the next level of operation and treated as missing. However, an invalid string expression that is not nested is displayed as a blank string and is *not* treated as missing.

### ***Missing values in numeric expressions***

- Most numeric expressions receive the system-missing value when any one of the values in the expression is missing.
- Some arithmetic operations involving 0 can be evaluated even when the variables have missing values. These operations are:

<b>Expression</b>	<b>Result</b>
0 * missing	0
0 / missing	0
MOD(0,missing)	0

- The *.n* suffix can be used with the statistical functions SUM, MEAN, MIN, MAX, SD, VARIANCE, and CFVAR to specify the number of valid arguments that you consider acceptable. The default of *n* is 2 for SD, VARIANCE, and CFVAR, and 1 for other statistical functions. For example,

```
COMPUTE FACTOR = SUM.2(SCORE1 TO SCORE3).
```

computes the variable *FACTOR* only if a case has valid information for at least two scores. *FACTOR* is assigned the system-missing value if a case has valid values for fewer than two scores. If the number specified exceeds the number of arguments in the function, the result is system-missing.

### ***Missing values in string expressions***

- If the numeric argument (which can be an expression) for the functions LPAD and RPAD is illegal or missing, the result is a null string. If the padding or trimming is the only operation, the string is then padded to its entire length with blanks. If the operation is nested, the null string is passed to the next nested level.
- If a numeric argument to SUBSTR is illegal or missing, the result is a null string. If SUBSTR is the only operation, the string is blank. If the operation is nested, the null string is passed to the next nested level.
- If a numeric argument to INDEX or RINDEX is illegal or missing, the result is system-missing.

### ***Missing values in logical expressions***

In a simple relation, the logic is indeterminate if the expression on either side of the relational operator is missing. When two or more relations are joined by logical operators AND and OR, a missing value is always returned if all of the relations in the expression are missing. However, if

any one of the relations can be determined, IBM® SPSS® Statistics tries to return true or false according to the logical outcomes. [For more information, see the topic AND and OR logical operators on p. 115.](#)

- When two relations are joined with the `AND` operator, the logical expression can never be true if one of the relations is indeterminate. The expression can, however, be false.
- When two relations are joined with the `OR` operator, the logical expression can never be false if one relation returns missing. The expression, however, can be true.

### ***Missing value functions***

- Each argument to a missing-value function (expression, variable name, or constant) must be separated by a comma.
- With the exception of the `MISSING` function, only numeric values can be used as arguments in missing-value functions.
- The keyword `TO` can be used to refer to a set of variables in the argument list for functions `NMISS` and `NVALID`.
- The functions `MISSING` and `SYSMIS` are logical functions and can be useful shortcuts to more complicated specifications on the `IF`, `DO IF`, and other conditional commands.

**VALUE.** `VALUE(variable)`. Numeric or string. Returns the value of variable, ignoring user missing-value definitions for variable, which must be a variable name or a vector reference to a variable name.

**MISSING.** `MISSING(variable)`. Logical. Returns 1 or true if variable has a system- or user-missing value. The argument should be a variable name in the active dataset.

**SYSMIS.** `SYSMIS(numvar)`. Logical. Returns 1 or true if the value of numvar is system-missing. The argument numvar must be the name of a numeric variable in the active dataset.

**NMISS.** `NMISS(variable[,...])`. Numeric. Returns a count of the arguments that have system- and user-missing values. This function requires one or more arguments, which should be variable names in the active dataset.

**NVALID.** `NVALID(variable[,...])`. Numeric. Returns a count of the arguments that have valid, nonmissing values. This function requires one or more arguments, which should be variable names in the active dataset.



# 2SLS

2SLS is available in the Regression option.

```
2SLS [EQUATION=]dependent variable WITH predictor variable
  [/[EQUATION=]dependent variable...]
  /INSTRUMENTS=varlist
  [/ENDOGENOUS=varlist]
  [/{CONSTANT**}
  {NOCONSTANT}]
  [/PRINT=COV]
  [/SAVE = [PRED] [RESID]]
  [/APPLY[='model name']]
```

**\*\***Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
2SLS VAR01 WITH VAR02 VAR03
  /INSTRUMENTS VAR03 LAGVAR01.
```

## **Overview**

2SLS performs two-stage least-squares regression to produce consistent estimates of parameters when one or more predictor variables might be correlated with the disturbance. This situation typically occurs when your model consists of a system of simultaneous equations wherein endogenous variables are specified as predictors in one or more of the equations. The two-stage least-squares technique uses instrumental variables to produce regressors that are not contemporaneously correlated with the disturbance. Parameters of a single equation or a set of simultaneous equations can be estimated.

## **Options**

**New Variables.** You can change `NEWVAR` settings on the `TSET` command prior to `2SLS` to evaluate the regression statistics without saving the values of predicted and residual variables, or you can save the new values to replace the values that were saved earlier, or you can save the new values without erasing values that were saved earlier (see the `TSET` command). You can also use the `SAVE` subcommand on `2SLS` to override the `NONE` or the default `CURRENT` settings on `NEWVAR`.

**Covariance Matrix.** You can obtain the covariance matrix of the parameter estimates in addition to all of the other output by specifying `PRINT=DETAILED` on the `TSET` command prior to `2SLS`. You can also use the `PRINT` subcommand to obtain the covariance matrix, regardless of the setting on `PRINT`.

### **Basic Specification**

The basic specification is at least one `EQUATION` subcommand and one `INSTRUMENTS` subcommand.

- For each specified equation, `2SLS` estimates and displays the regression analysis-of-variance table, regression standard error, mean of the residuals, parameter estimates, standard errors of the parameter estimates, standardized parameter estimates, *t* statistic significance tests and probability levels for the parameter estimates, tolerance of the variables, and correlation matrix of the parameter estimates.
- If the setting on `NEWVAR` is either `ALL` or the default `CURRENT`, two new variables containing the predicted and residual values are automatically created for each equation. The variables are labeled and added to the active dataset.

### **Subcommand Order**

- Subcommands can be specified in any order.

### **Syntax Rules**

- The `INSTRUMENTS` subcommand must specify at least as many variables as are specified after `WITH` on the longest `EQUATION` subcommand.
- If a subcommand is specified more than once, the effect is cumulative (except for the `APPLY` subcommand, which executes only the last occurrence).

### **Operations**

- `2SLS` cannot produce forecasts beyond the length of any regressor series.
- `2SLS` honors the `WEIGHT` command.
- `2SLS` uses listwise deletion of missing data. Whenever a variable is missing a value for a particular observation, that observation will not be used in any of the computations.

## **Examples**

```
* 2-Stage Least Squares.
TSET NEWVAR=NONE .
2SLS buyoff WITH buycd buybk offer_type1 offer_type2
  /INSTRUMENTS offer_type1 offer_type2 lndiscdd lndiscbk buycd_1 buybk_1
  /CONSTANT .
```

- The procedure fits a model for *buyoff* using *buycd*, *buyck*, *offer\_type1*, and *offer\_type2* as explanatory variables.
- The `INSTRUMENTS` subcommand specifies that *offer\_type1*, *offer\_type2*, *lndiscdd*, *lndiscbk*, *buycd\_1*, and *buybk\_1* are used as instrument variables to predict *buycd* and *buybk*.
- All other options are set to their default values.

## ***EQUATION Subcommand***

`EQUATION` specifies the structural equations for the model and is required. The actual keyword `EQUATION` is optional.

- An equation specifies a single dependent variable, followed by keyword `WITH` and one or more predictor variables.
- You can specify more than one equation. Multiple equations are separated by slashes.

### ***Example***

```
2SLS EQUATION=Y1 WITH X1 X2
/INSTRUMENTS=X1 LAGX2 X3.
```

- In this example, *Y1* is the dependent variable, and *X1* and *X2* are the predictors. The instruments that are used to predict the *X2* values are *X1*, *LAGX2*, and *X3*.

## ***INSTRUMENTS Subcommand***

`INSTRUMENTS` specifies the instrumental variables. These variables are used to compute predicted values for the endogenous variables in the first stage of 2SLS.

- At least one `INSTRUMENTS` subcommand must be specified.
- If more than one `INSTRUMENTS` subcommand is specified, the effect is cumulative. All variables that are named on `INSTRUMENTS` subcommands are used as instruments to predict all the endogenous variables.
- Any variable in the active dataset can be named as an instrument.
- Instrumental variables can be specified on the `EQUATION` subcommand, but this specification is not required.
- The `INSTRUMENTS` subcommand must name at least as many variables as are specified after `WITH` on the longest `EQUATION` subcommand.
- If all the predictor variables are listed as the only `INSTRUMENTS`, the results are the same as results from ordinary least-squares regression.

### ***Example***

```
2SLS DEMAND WITH PRICE, INCOME
/PRICE WITH DEMAND, RAINFALL, LAGPRICE
/INSTRUMENTS=INCOME, RAINFALL, LAGPRICE.
```

- The endogenous variables are `PRICE` and `DEMAND`.
- The instruments to be used to compute predicted values for the endogenous variables are `INCOME`, `RAINFALL`, and `LAGPRICE`.

## ***ENDOGENOUS Subcommand***

All variables that are not specified on the `INSTRUMENTS` subcommand are used as endogenous variables by `2SLS`. The `ENDOGENOUS` subcommand simply allows you to document what these variables are.

- Computations are not affected by specifications on the `ENDOGENOUS` subcommand.

### ***Example***

```
2SLS Y1 WITH X1 X2 X3
      /INSTRUMENTS=X2 X4 LAGY1
      /ENDOGENOUS=Y1 X1 X3 .
```

- In this example, the `ENDOGENOUS` subcommand is specified to document the endogenous variables.

## ***CONSTANT and NOCONSTANT Subcommands***

Specify `CONSTANT` or `NOCONSTANT` to indicate whether a constant term should be estimated in the regression equation. The specification of either subcommand overrides the `CONSTANT` setting on the `TSET` command for the current procedure.

- `CONSTANT` is the default and specifies that the constant term is used as an instrument.
- `NOCONSTANT` eliminates the constant term.

## ***SAVE Subcommand***

`SAVE` saves the values of predicted and residual variables that are generated during the current session to the end of the active dataset. The default names `FITn` and `ERRn` will be generated, where `n` increments each time variables are saved for an equation. `SAVE` overrides the `NONE` or the default `CURRENT` setting on `NEWVAR` for the current procedure.

- |               |  |
|---------------|--|
| <b>PRED</b>   | <i>Save the predicted value.</i> The new variable is named <code>FIT<sub>n</sub></code> , where <code>n</code> increments each time a predicted or residual variable is saved for an equation. |
| <b>RESSID</b> | <i>Save the residual value.</i> The new variable is named <code>ERR<sub>n</sub></code> , where <code>n</code> increments each time a predicted or residual variable is saved for an equation.  |

## ***PRINT Subcommand***

`PRINT` can be used to produce an additional covariance matrix for each equation. The only specification on this subcommand is keyword `COV`. The `PRINT` subcommand overrides the `PRINT` setting on the `TSET` command for the current procedure.

## ***APPLY Subcommand***

`APPLY` allows you to use a previously defined `2SLS` model without having to repeat the specifications.

- The only specification on `APPLY` is the name of a previous model. If a model name is not specified, the model that was specified on the previous `2SLS` command is used.
- To change the series that are used with the model, enter new series names before or after the `APPLY` subcommand.
- To change one or more model specifications, specify the subcommands of only those portions that you want to change after the `APPLY` subcommand.
- If no series are specified on the command, the series that were originally specified with the model that is being reapplied are used.

**Example**

```
2SLS Y1 WITH X1 X2 / X1 WITH Y1 X2
      /INSTRUMENTS=X2 X3 .
2SLS APPLY
      /INSTRUMENTS=X2 X3 LAGX1 .
```

- In this example, the first command requests `2SLS` using `X2` and `X3` as instruments.
- The second command specifies the same equations but changes the instruments to `X2`, `X3`, and `LAGX1`.

# ACF

```
ACF VARIABLES= series names

[/DIFF={1}]
      {n}

[/SDIFF={1}]
      {n}

[/PERIOD=n]

[/{NOLOG**}]
      {LN   }

[/SEASONAL]

[/MXAUTO={16**}]
      {n   }

[/ERROR={IND**}]
      {MA  }

[/PACF]

[/APPLY [= 'model name']]
```

**\*\*Default if the subcommand is omitted and there is no corresponding specification on the TSET command.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
ACF TICKETS.
```

## Overview

ACF displays and plots the sample autocorrelation function of one or more time series. You can also display and plot the autocorrelations of transformed series by requesting natural log and differencing transformations within the procedure.

### Options

**Modifying the Series.** You can request a natural log transformation of the series using the LN subcommand and seasonal and nonseasonal differencing to any degree using the SDIFF and DIFF subcommands. With seasonal differencing, you can specify the periodicity on the PERIOD subcommand.

**Statistical Output.** With the MXAUTO subcommand, you can specify the number of lags for which you want autocorrelations to be displayed and plotted, overriding the maximum specified on TSET. You can also display and plot values at periodic lags only using the SEASONAL

subcommand. In addition to autocorrelations, you can display and plot partial autocorrelations using the `PACF` subcommand.

**Method of Calculating Standard Errors.** You can specify one of two methods of calculating the standard errors for the autocorrelations on the `SERROR` subcommand.

### ***Basic Specification***

The basic specification is one or more series names.

- For each series specified, `ACF` automatically displays the autocorrelation value, standard error, Box-Ljung statistic, and probability for each lag.
- `ACF` plots the autocorrelations and marks the bounds of two standard errors on the plot. By default, `ACF` displays and plots autocorrelations for up to 16 lags or the number of lags specified on `TSET`.
- If a method has not been specified on `TSET`, the default method of calculating the standard error (`IND`) assumes that the process is white noise.

### ***Subcommand Order***

- Subcommands can be specified in any order.

### ***Syntax Rules***

- `VARIABLES` can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each one is executed.

### ***Operations***

- Subcommand specifications apply to all series named on the `ACF` command.
- If the `LN` subcommand is specified, any differencing requested on that `ACF` command is done on the log-transformed series.
- Confidence limits are displayed in the plot, marking the bounds of two standard errors at each lag.

### ***Limitations***

- A maximum of one `VARIABLES` subcommand. There is no limit on the number of series named on the list.

## ***Example***

```
ACF VARIABLES = TICKETS
  /LN
  /DIFF=1
  /SDIFF=1
  /PER=12
  /MXAUTO=50 .
```

- This example produces a plot of the autocorrelation function for the series *TICKETS* after a natural log transformation, differencing, and seasonal differencing have been applied. Along with the plot, the autocorrelation value, standard error, Box-Ljung statistic, and probability are displayed for each lag.
- LN transforms the data using the natural logarithm (base *e*) of the series.
- DIFF differences the series once.
- SDIFF and PERIOD apply one degree of seasonal differencing with a period of 12.
- MXAUTO specifies that the maximum number of lags for which output is to be produced is 50.

## **VARIABLES Subcommand**

VARIABLES specifies the series names and is the only required subcommand.

## **DIFF Subcommand**

DIFF specifies the degree of differencing used to convert a nonstationary series to a stationary one with a constant mean and variance before the autocorrelations are computed.

- You can specify 0 or any positive integer on DIFF.
- If DIFF is specified without a value, the default is 1.
- The number of values used in the calculations decreases by 1 for each degree-1 of differencing.

### **Example**

```
ACF VARIABLES = SALES
  /DIFF=1.
```

- In this example, the series *SALES* will be differenced once before the autocorrelations are computed and plotted.

## **SDIFF Subcommand**

If the series exhibits a seasonal or periodic pattern, you can use the SDIFF subcommand to seasonally difference the series before obtaining autocorrelations.

- The specification on SDIFF indicates the degree of seasonal differencing and can be 0 or any positive integer.
- If SDIFF is specified without a value, the degree of seasonal differencing defaults to 1.
- The number of seasons used in the calculations decreases by 1 for each degree of seasonal differencing.
- The length of the period used by SDIFF is specified on the PERIOD subcommand. If the PERIOD subcommand is not specified, the periodicity established on the TSET or DATE command is used (see the PERIOD subcommand).



## ***PERIOD Subcommand***

PERIOD indicates the length of the period to be used by the SDIFF or SEASONAL subcommands.

- The specification on PERIOD indicates how many observations are in one period or season and can be any positive integer greater than 1.
- The PERIOD subcommand is ignored if it is used without the SDIFF or SEASONAL subcommands.
- If PERIOD is not specified, the periodicity established on TSET PERIOD is in effect. If TSET PERIOD is not specified, the periodicity established on the DATE command is used. If periodicity was not established anywhere, the SDIFF and SEASONAL subcommands will not be executed.

### ***Example***

```
ACF VARIABLES = SALES
  /SDIFF=1M
  /PERIOD=12 .
```

- This command applies one degree of seasonal differencing with a periodicity (season) of 12 to the series *SALES* before autocorrelations are computed.

## ***LN and NOLOG Subcommands***

LN transforms the data using the natural logarithm (base  $e$ ) of the series and is used to remove varying amplitude over time. NOLOG indicates that the data should not be log transformed. NOLOG is the default.

- If you specify LN on an ACF command, any differencing requested on that command will be done on the log-transformed series.
- There are no additional specifications on LN or NOLOG.
- Only the last LN or NOLOG subcommand on an ACF command is executed.
- If a natural log transformation is requested when there are values in the series that are less than or equal to zero, the ACF will not be produced for that series because nonpositive values cannot be log transformed.
- NOLOG is generally used with an APPLY subcommand to turn off a previous LN specification.

### ***Example***

```
ACF VARIABLES = SALES
  /LN .
```

- This command transforms the series *SALES* using the natural log transformation and then computes and plots autocorrelations.

## ***SEASONAL Subcommand***

Use the SEASONAL subcommand to focus attention on the seasonal component by displaying and plotting autocorrelations at periodic lags only.

- There are no additional specifications on `SEASONAL`.
- If `SEASONAL` is specified, values are displayed and plotted at the periodic lags indicated on the `PERIOD` subcommand. If `PERIOD` is not specified, the periodicity established on the `TSET` or `DATE` command is used (see the `PERIOD` subcommand).
- If `SEASONAL` is not specified, autocorrelations for all lags up to the maximum are displayed and plotted.

### **Example**

```
ACF VARIABLES = SALES
  /SEASONAL
  /PERIOD=12 .
```

- In this example, autocorrelations are displayed only at every 12th lag.

## **MXAUTO Subcommand**

`MXAUTO` specifies the maximum number of lags for a series.

- The specification on `MXAUTO` must be a positive integer.
- If `MXAUTO` is not specified, the default number of lags is the value set on `TSET MXAUTO`. If `TSET MXAUTO` is not specified, the default is 16.
- The value on `MXAUTO` overrides the value set on `TSET MXAUTO`.

### **Example**

```
ACF VARIABLES = SALES
  /MXAUTO=14 .
```

- This command sets the maximum number of autocorrelations to be displayed for the series `SALES` to 14.

## **SERROR Subcommand**

`SERROR` specifies the method of calculating the standard errors for the autocorrelations.

- You must specify either the keyword `IND` or `MA` on `SERROR`.
- The method specified on `SERROR` overrides the method specified on the `TSET ACFSE` command.
- If `SERROR` is not specified, the method indicated on `TSET ACFSE` is used. If `TSET ACFSE` is not specified, the default is `IND`.

**IND**      *Independence model.* The method of calculating the standard errors assumes that the underlying process is white noise.

**MA**      *MA model.* The method of calculating the standard errors is based on Bartlett's approximation. With this method, appropriate where the true `MA` order of the process is  $k-1$ , standard errors grow at increased lags (Pankratz, 1983).

**Example**

```
ACF VARIABLES = SALES
  /SERROR=MA.
```

- In this example, the standard errors of the autocorrelations are computed using the MA method.

**PACF Subcommand**

Use the PACF subcommand to display and plot sample partial autocorrelations as well as autocorrelations for each series named on the ACF command.

- There are no additional specifications on PACF.
- PACF also displays the standard errors of the partial autocorrelations and indicates the bounds of two standard errors on the plot.
- With the exception of SERROR, all other subcommands specified on that ACF command apply to both the partial autocorrelations and the autocorrelations.

**Example**

```
ACF VARIABLES = SALES
  /DIFFERENCE=1
  /PACF.
```

- This command requests both autocorrelations and partial autocorrelations for the series SALES after it has been differenced once.

**APPLY Subcommand**

APPLY allows you to use a previously defined ACF model without having to repeat the specifications.

- The only specification on APPLY is the name of a previous model in quotation marks. If a model name is not specified, the model specified on the previous ACF command is used.
- To change one or more model specifications, specify the subcommands of only those portions you want to change after the APPLY subcommand.
- If no series are specified on the ACF command, the series that were originally specified with the model being reapplied are used.
- To change the series used with the model, enter new series names before or after the APPLY subcommand.

**Example**

```
ACF VARIABLES = TICKETS
  /LN
  /DIFF=1
  /SDIFF=1
  /PERIOD=12
  /MXAUTO=50.
ACF VARIABLES = ROUNDTRP
  /APPLY.
ACF APPLY
```

```
/NOLOG.  
ACF APPLY 'MOD_2 '  
/PERIOD=6 .
```

- The first command requests a maximum of 50 autocorrelations for the series *TICKETS* after a natural log transformation, differencing, and one degree of seasonal differencing with a periodicity of 12 have been applied. This model is assigned the default name *MOD\_1*.
- The second command displays and plots the autocorrelation function for the series *ROUNDTRP* using the same model that was used for the series *TICKETS*. This model is assigned the name *MOD\_2*.
- The third command requests another autocorrelation function of the series *ROUNDTRP* using the same model but without the natural log transformation. Note that when *APPLY* is the first specification after the *ACF* command, the slash (/) before it is not necessary. This model is assigned the name *MOD\_3*.
- The fourth command reapplies *MOD\_2*, autocorrelations for the series *ROUNDTRP* with the natural log and differencing specifications, but this time with a periodicity of 6. This model is assigned the name *MOD\_4*. It differs from *MOD\_2* only in the periodicity.

## **References**

- Box, G. E. P., and G. M. Jenkins. 1976. *Time series analysis: Forecasting and control*, Rev. ed. San Francisco: Holden-Day.
- Pankratz, A. 1983. *Forecasting with univariate Box-Jenkins models: Concepts and cases*. New York: John Wiley and Sons.

# ***ADD DOCUMENT***

```
ADD DOCUMENT  
  'text'  
  'text'.
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Example***

```
ADD DOCUMENT  
  "This data file is a 10% random sample from the"  
  "master data file. It's seed value is 13254689.".
```

## ***Overview***

ADD DOCUMENT saves a block of text of any length in the active dataset. The result is equivalent to the DOCUMENT command. The documentation can be displayed with the DISPLAY DOCUMENT command.

When GET retrieves a data file, or APPLY DICTIONARY is used to apply documents from another data file, or ADD FILES, MATCH FILES, or UPDATE is used to combine data files, all documents from each specified file are copied into the working file. DROP DOCUMENTS can be used to drop those documents from the working file.

## ***Basic Specification***

The basic specification is ADD DOCUMENT followed by one or more optional lines of quoted text. The text is stored in the file dictionary when the data file is saved in IBM® SPSS® Statistics format.

## ***Syntax Rules***

- Each line must be enclosed in single or double quotation marks, following the standard rules for quoted strings.
- Each line can be up to 80 bytes long (typically 80 characters in single-byte languages), including the command name but not including the quotation marks used to enclose the text. If any line exceeds 80 bytes, an error will result and the command will not be executed.
- The text can be entered on as many lines as needed.
- Multiple ADD DOCUMENT commands can be specified for the same data file.

## ***Operations***

- The text from each ADD DOCUMENT command is appended to the end of the list of documentation, followed by the date in parentheses.
- An ADD DOCUMENT command with no quoted text string appends a date in parentheses to the documentation.

---

*ADD DOCUMENT*

- `DISPLAY DOCUMENTS` will display all documentation for the data file specified on the `ADD DOCUMENT` and/or `DOCUMENT` commands. Documentation is displayed exactly as entered; each line of the `ADD DOCUMENT` command is displayed as a separate line, and there is no line wrapping.
- `DROP DOCUMENTS` deletes all documentation created by both `ADD DOCUMENT` and `DOCUMENT`.

***Example***

If the command name and the quoted text string are specified on the same line, the command name counts toward the 80-byte line limit, so it's a good idea to put the command name on a separate line, as in:

```
ADD DOCUMENT
  "This is some text that describes this file."
```

***Example***

To insert blank lines between blocks of text, enter a null string, as in:

```
ADD DOCUMENT
  "This is some text that describes this file."
  ""
  "This is some more text preceded by a blank line."
```

# ADD FILES

```
ADD FILES FILE={'savfile'|'dataset'}
           {*
           }

[/RENAME=(old varnames=new varnames)...]

[/IN=varname]

/FILE=... [/RENAME=...] [/IN=...]

[/BY varlist]

[/MAP]

[/KEEP={ALL** }] [/DROP=varlist]
           {varlist}

[/FIRST=varname] [/LAST=varname]
```

\*\*Default if the subcommand is omitted.

## Example

```
ADD FILES FILE="/data/school1.sav" /FILE="/data/school2.sav".
```

## Overview

ADD FILES combines cases from 2 up to 50 open data sets or external IBM® SPSS® Statistics data files by concatenating or interleaving cases. When cases are **concatenated**, all cases from one file are added to the end of all cases from another file. When cases are **interleaved**, cases in the resulting file are ordered according to the values of one or more key variables.

The files specified on ADD FILES can be external SPSS Statistics data files and/or currently open datasets. The combined file becomes the new active dataset.

In general, ADD FILES is used to combine files containing the same variables but different cases. To combine files containing the same cases but different variables, use MATCH FILES. To update existing SPSS Statistics data files, use UPDATE.

## Options

**Variable Selection.** You can specify which variables from each input file are included in the new active dataset using the DROP and KEEP subcommands.

**Variable Names.** You can rename variables in each input file before combining the files using the RENAME subcommand. This permits you to combine variables that are the same but whose names differ in different input files or to separate variables that are different but have the same name.

**Variable Flag.** You can create a variable that indicates whether a case came from a particular input file using IN. When interleaving cases, you can use the FIRST or LAST subcommands to create a variable that flags the first or last case of a group of cases with the same value for the key variable.

**Variable Map.** You can request a map showing all variables in the new active dataset, their order, and the input files from which they came using the MAP subcommand.

**Basic Specification**

- The basic specification is two or more `FILE` subcommands, each of which specifies a file to be combined. If cases are to be interleaved, the `BY` subcommand specifying the key variables is also required.
- All variables from all input files are included in the new active dataset unless `DROP` or `KEEP` is specified.

**Subcommand Order**

- `RENAME` and `IN` must immediately follow the `FILE` subcommand to which they apply.
- `BY`, `FIRST`, and `LAST` must follow all `FILE` subcommands and their associated `RENAME` and `IN` subcommands.

**Syntax Rules**

- `RENAME` can be repeated after each `FILE` subcommand. `RENAME` applies only to variables in the file named on the `FILE` subcommand immediately preceding it.
- `BY` can be specified only once. However, multiple key variables can be specified on `BY`. When `BY` is used, all files must be sorted in ascending order by the key variables (see `SORT CASES`).
- `FIRST` and `LAST` can be used only when files are interleaved (when `BY` is used).
- `MAP` can be repeated as often as desired.

**Operations**

- `ADD FILES` reads all input files named on `FILE` and builds a new active dataset. `ADD FILES` is executed when the data are read by one of the procedure commands or the `EXECUTE`, `SAVE`, or `SORT CASES` commands.
- If the current active dataset is included and is specified with an asterisk (`FILE=*`), the new merged dataset replaces the active dataset. If that dataset is a named dataset, the merged dataset retains that name. If the current active dataset is not included or is specified by name (for example, `FILE=Dataset1`), a new unnamed, merged dataset is created, and it becomes the active dataset. For information on naming datasets, see [DATASET NAME](#).
- The resulting file contains complete dictionary information from the input files, including variable names, labels, print and write formats, and missing-value indicators. It also contains the documents from each input file. See `DROP DOCUMENTS` for information on deleting documents.
- For each variable, dictionary information is taken from the first file containing value labels, missing values, or a variable label for the common variable. If the first file has no such information, `ADD FILES` checks the second file, and so on, seeking dictionary information.
- Variables are copied in order from the first file specified, then from the second file specified, and so on. Variables that are not contained in all files receive the system-missing value for cases that do not have values for those variables.
- If the same variable name exists in more than one file but the format type (numeric or string) does not match, the command is not executed.
- If a numeric variable has the same name but different formats (for example, `F8.0` and `F8.2`) in different input files, the format of the variable in the first-named file is used.



- If a string variable has the same name but different formats (for example, A24 and A16) in different input files, the command is not executed.
- If the active dataset is named as an input file, any `N` and `SAMPLE` commands that have been specified are applied to the active dataset before the files are combined.
- If only one of the files is weighted, the program turns weighting off when combining cases from the two files. To weight the cases, use the `WEIGHT` command again.

### Limitations

- A maximum of 50 files can be combined on one `ADD FILES` command.
- The `TEMPORARY` command cannot be in effect if the active dataset is used as an input file.

## Examples

```
ADD FILES FILE="/data/school1.sav"
        /FILE="/data/school2.sav".
```

- `ADD FILES` concatenates cases from the IBM® SPSS® Statistics data files *school1.sav* and *school2.sav*. All cases from *school1.sav* precede all cases from *school2.sav* in the resulting file.

```
SORT CASES BY LOCATN DEPT.
ADD FILES FILE="/data/source.sav" /FILE=* /BY LOCATN DEPT
        /KEEP AVGHOUR AVGRAISE LOCATN DEPT SEX HOURLY RAISE /MAP.
SAVE OUTFILE="/data/prsnnl.sav".
```

- `SORT CASES` sorts cases in the active dataset in ascending order of their values for *LOCATN* and *DEPT*.
- `ADD FILES` combines two files: the external SPSS Statistics data file *source.sav* and the sorted active dataset. The file *source.sav* must also be sorted by *LOCATN* and *DEPT*.
- `BY` indicates that the keys for interleaving cases are *LOCATN* and *DEPT*, the same variables used on `SORT CASES`.
- `KEEP` specifies the variables to be retained in the resulting file.
- `MAP` produces a list of variables in the resulting file and the two input files.
- `SAVE` saves the resulting file as a new SPSS Statistics data file named *prsnnl.sav*.

## FILE Subcommand

`FILE` identifies the files to be combined. A separate `FILE` subcommand must be used for each input file.

- An asterisk may be specified on `FILE` to indicate the active dataset.
- Dataset names instead of file names can be used to refer to currently open datasets.
- The order in which files are named determines the order of cases in the resulting file.

### Example

```
GET DATA /TYPE=XLS /FILE='/temp/excelfile1.xls'.
```

```
DATASET NAME exceldata1.
GET DATA /TYPE=XLS /FILE='/temp/excelfile2.xls'.
ADD FILES FILE='exceldata1'
        /FILE=*
        /FILE='/temp/mydata.sav'.
```

## RENAME Subcommand

RENAME renames variables in input files *before* they are processed by ADD FILES. RENAME follows the FILE subcommand that specifies the file containing the variables to be renamed.

- RENAME applies only to the FILE subcommand immediately preceding it. To rename variables from more than one input file, enter a RENAME subcommand after each FILE subcommand that specifies a file with variables to be renamed.
- Specifications for RENAME consist of a left parenthesis, a list of old variable names, an equals sign, a list of new variable names, and a right parenthesis. The two variable lists must name or imply the same number of variables. If only one variable is renamed, the parentheses are optional.
- More than one such specification can be entered on a single RENAME subcommand, each enclosed in parentheses.
- The TO keyword can be used to refer to consecutive variables in the file and to generate new variable names.
- RENAME takes effect immediately. KEEP and DROP subcommands entered prior to RENAME must use the old names, while those entered after RENAME must use the new names.
- All specifications within a single set of parentheses take effect simultaneously. For example, the specification RENAME (A, B = B, A) swaps the names of the two variables.
- Variables cannot be renamed to scratch variables.
- Input data files are not changed on disk; only the copy of the file being combined is affected.

### Example

```
ADD FILES FILE="/data/clients.sav" /RENAME=(TEL_NO, ID_NO = PHONE, ID)
        /FILE="/data/master.sav" /BY ID.
```

- ADD FILES adds new client cases from the file *clients.sav* to existing client cases in the file *master.sav*.
- Two variables on *clients.sav* are renamed prior to the match. *TEL\_NO* is renamed *PHONE* to match the name used for phone numbers in the master file. *ID\_NO* is renamed *ID* so that it will have the same name as the identification variable in the master file and can be used on the BY subcommand.
- The BY subcommand orders the resulting file according to client ID number.

## BY Subcommand

BY specifies one or more key variables that determine the order of cases in the resulting file. When BY is specified, cases from the input files are interleaved according to their values for the key variables.

- BY must follow the FILE subcommands and any associated RENAME and IN subcommands.
- The key variables specified on BY must be present and have the same names in all input files.
- Key variables can be string or numeric.
- All input files must be sorted in ascending order of the key variables. If necessary, use SORT CASES before ADD FILES.
- Cases in the resulting file are ordered by the values of the key variables. All cases from the first file with the first value for the key variable are first, followed by all cases from the second file with the same value, followed by all cases from the third file with the same value, and so forth. These cases are followed by all cases from the first file with the next value for the key variable, and so on.
- Cases with system-missing values are first in the resulting file. User-missing values are interleaved with other values.

## ***DROP and KEEP Subcommands***

DROP and KEEP are used to include only a subset of variables in the resulting file. DROP specifies a set of variables to exclude and KEEP specifies a set of variables to retain.

- DROP and KEEP do not affect the input files on disk.
- DROP and KEEP must follow all FILE and RENAME subcommands.
- DROP and KEEP must specify one or more variables. If RENAME is used to rename variables, specify the new names on DROP and KEEP.
- DROP and KEEP take effect immediately. If a variable specified on DROP or KEEP does not exist in the input files, was dropped by a previous DROP subcommand, or was not retained by a previous KEEP subcommand, the program displays an error message and does not execute the ADD FILES command.
- DROP cannot be used with variables created by the IN, FIRST, or LAST subcommands.
- KEEP can be used to change the order of variables in the resulting file. With KEEP, variables are kept in the order in which they are listed on the subcommand. If a variable is named more than once on KEEP, only the first mention of the variable is in effect; all subsequent references to that variable name are ignored.
- The keyword ALL can be specified on KEEP. ALL must be the last specification on KEEP, and it refers to all variables not previously named on that subcommand. It is useful when you want to arrange the first few variables in a specific order.

### ***Example***

```
ADD FILES FILE="/data/particle.sav" /RENAME=(PARTIC=pollute1)
  /FILE="/data/gas.sav" /RENAME=(OZONE TO SULFUR=pollut2 TO pollute4)
  /KEEP=pollute1 pollute2 pollute3 pollute4.
```

- The renamed variables are retained in the resulting file. KEEP is specified after all the FILE and RENAME subcommands, and it refers to the variables by their new names.

## ***IN Subcommand***

`IN` creates a new variable in the resulting file that indicates whether a case came from the input file named on the preceding `FILE` subcommand. `IN` applies only to the file specified on the immediately preceding `FILE` subcommand.

- `IN` has only one specification, the name of the flag variable.
- The variable created by `IN` has the value 1 for every case that came from the associated input file and the value 0 for every case that came from a different input file.
- Variables created by `IN` are automatically attached to the end of the resulting file and cannot be dropped. If `FIRST` or `LAST` are used, the variable created by `IN` precedes the variables created by `FIRST` or `LAST`.

### ***Example***

```
ADD FILES FILE="/data/week10.sav" /FILE="/data/week11.sav"  
/IN=INWEEK11 /BY=EMPID.
```

- `IN` creates the variable `INWEEK11`, which has the value 1 for all cases in the resulting file that came from the input file `week11.sav` and the value 0 for those cases that were not in the file `week11.sav`.

### ***Example***

```
ADD FILES FILE="/data/week10.sav" /FILE="/data/week11.sav"  
/IN=INWEEK11 /BY=EMPID.  
IF (NOT INWEEK11) SALARY1=0.
```

- The variable created by `IN` is used to screen partially missing cases for subsequent analyses.
- Since `IN` variables have either the value 1 or 0, they can be used as logical expressions, where 1 = true and 0 = false. The `IF` command sets the variable `SALARY1` equal to 0 for all cases that came from the file `INWEEK11`.

## ***FIRST and LAST Subcommands***

`FIRST` and `LAST` create logical variables that flag the first or last case of a group of cases with the same value on the `BY` variables. `FIRST` and `LAST` must follow all `FILE` subcommands and their associated `RENAME` and `IN` subcommands.

- `FIRST` and `LAST` have only one specification, the name of the flag variable.
- `FIRST` creates a variable with the value 1 for the first case of each group and the value 0 for all other cases.
- `LAST` creates a variable with the value 1 for the last case of each group and the value 0 for all other cases.
- Variables created by `FIRST` and `LAST` are automatically attached to the end of the resulting file and cannot be dropped.

**Example**

```
ADD FILES FILE="/data/school1.sav" /FILE="/data/school2.sav"
  /BY=GRADE /FIRST=HISCORE.
```

- The variable *HISCORE* contains the value 1 for the first case in each grade in the resulting file and the value 0 for all other cases.

**MAP Subcommand**

MAP produces a list of the variables included in the new active dataset and the file or files from which they came. Variables are listed in the order in which they exist in the resulting file. MAP has no specifications and must follow after all FILE and RENAME subcommands.

- Multiple MAP subcommands can be used. Each MAP subcommand shows the current status of the active dataset and reflects only the subcommands that precede the MAP subcommand.
- To obtain a map of the active dataset in its final state, specify MAP last.
- If a variable is renamed, its original and new names are listed. Variables created by IN, FIRST, and LAST are not included in the map, since they are automatically attached to the end of the file and cannot be dropped.

**Adding Cases from Different Data Sources**

You can add cases from any data source that IBM® SPSS® Statistics can read by defining dataset names for each data source that you read (DATASET NAME command) and then using ADD FILES to add the cases from each file. The following example merges the contents of three text data files, but it could just as easily merge the contents of a text data file, and Excel spreadsheet, and a database table.

**Example**

```
DATA LIST FILE="/data/gasdata1.txt"
  /1 OZONE 10-12 CO 20-22 SULFUR 30-32.
DATASET NAME gasdata1.
DATA LIST FILE="/data/gasdata2.txt"
  /1 OZONE 10-12 CO 20-22 SULFUR 30-32.
DATASET NAME gasdata2.
DATA LIST FILE="/data/gasdata3.txt"
  /1 OZONE 10-12 CO 20-22 SULFUR 30-32.
DATASET NAME gasdata3.
ADD FILES FILE='gasdata1'
  /FILE='gasdata2'
  /FILE='gasdata3'.
SAVE OUTFILE='/data/combined_data.sav'.
```

# ADD VALUE LABELS

```
ADD VALUE LABELS varlist value 'label' value 'label'...[/varlist...]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
ADD VALUE LABELS JOBGRADE 'P' 'Parttime Employee'  
                        'C' 'Customer Support'.
```

## **Overview**

ADD VALUE LABELS adds or alters value labels without affecting other value labels already defined for that variable. In contrast, VALUE LABELS adds or alters value labels but deletes all existing value labels for that variable when it does so.

## **Basic Specification**

The basic specification is a variable name and individual values with associated labels.

## **Syntax Rules**

- Labels can be assigned to values of any previously defined variable. It is not necessary to enter value labels for all of a variable's values.
- Each value label must be enclosed in single or double quotes.
- To specify a single quote or apostrophe within a quoted string, either enclose the entire string in double quotes or double the single quote/apostrophe.
- Value labels can contain any characters, including blanks.
- The same labels can be assigned to the same values of different variables by specifying a list of variable names. For string variables, the variables on the list must have the same defined width (for example, A8).
- Multiple sets of variable names and value labels can be specified on one ADD VALUE LABELS command as long as each set is separated from the previous one by a slash.
- To continue a label from one command line to the next, specify a plus sign (+) before the continuation of the label and enclose each segment of the label, including the blank between them, in single or double quotes.

## **Operations**

- Unlike most transformations, ADD VALUE LABELS takes effect as soon as it is encountered in the command sequence. Thus, special attention should be paid to its position among commands.
- The added value labels are stored in the active dataset dictionary.

- ADD VALUE LABELS can be used for variables that have no previously assigned value labels.
- Adding labels to some values does not affect labels previously assigned to other values.

### **Limitations**

- Value labels cannot exceed 120 bytes.

## **Examples**

### **Adding Value Labels**

```
ADD VALUE LABELS V1 TO V3 1 'Officials & Managers'
                        6 'Service Workers'
/V4 'N' 'New Employee'.
```

- Labels are assigned to the values 1 and 6 of the variables between and including *V1* and *V3* in the active dataset.
- Following the required slash, a label for the value *N* for the variable *V4* is specified. *N* is a string value and must be enclosed in single or double quotes.
- If labels already exist for these values, they are changed in the dictionary. If labels do not exist for these values, new labels are added to the dictionary.
- Existing labels for other values for these variables are not affected.

### **Specifying a Label on Multiple Lines**

```
ADD VALUE LABELS OFFICE88 1 "EMPLOYEE'S OFFICE ASSIGNMENT PRIOR"
+ " TO 1988".
```

- The label for the value 1 for *OFFICE88* is specified on two command lines. The plus sign concatenates the two string segments, and a blank is included at the beginning of the second string in order to maintain correct spacing in the label.

## **Value Labels for String Variables**

- For string variables, the values and the labels must be enclosed in single or double quotes.
- If a specified value is longer than the defined width of the variable, the program displays a warning and truncates the value. The added label will be associated with the truncated value.
- If a specified value is shorter than the defined width of the variable, the program adds blanks to right-pad the value without warning. The added label will be associated with the padded value.
- If a single set of labels is to be assigned to a list of string variables, the variables must have the same defined width (for example, A8).

### **Example**

```
ADD VALUE LABELS STATE 'TEX' 'TEXAS' 'TEN' 'TENNESSEE'
                        'MIN' 'MINNESOTA'.
```

- ADD VALUE LABELS assigns labels to three values of the variable *STATE*. Each value and each label is specified in quotes.

- Assuming that the variable *STATE* is defined as three characters wide, the labels *TEXAS*, *TENNESSEE*, and *MINNESOTA* will be appropriately associated with the values *TEX*, *TEN*, and *MIN*. However, if *STATE* was defined as two characters wide, the program would truncate the specified values to two characters and would not be able to associate the labels correctly. Both *TEX* and *TEN* would be truncated to *TE* and would first be assigned the label *TEXAS*, which would then be changed to *TENNESSEE* by the second specification.

**Example**

```
ADD VALUE LABELS STATE REGION "U" "UNKNOWN".
```

- The label *UNKNOWN* is assigned to the value *U* for both *STATE* and *REGION*.
- *STATE* and *REGION* must have the same defined width. If they do not, a separate specification must be made for each, as in the following:

```
ADD VALUE LABELS STATE "U" "UNKNOWN" / REGION "U" "UNKNOWN".
```



# ADP

ADP is available in the Data Preparation option.

```
ADP
  /FIELDS [TARGET=targetField]
        INPUT=predictorFieldlist
        [ANALYSISWEIGHT=analysisweightField]
  [/PREPDATE TIME
    [DATEDURATION={YES** (REFERENCE={CURRENT**
                                }
                                {YMD(datespec)})
                    UNIT={AUTO**
                        {YEARS[(SUFFIX={'_years' })] }
                        {suffixname}
                        {MONTHS[(SUFFIX={'_months' })] }
                        {suffixname}
                        {DAYS[(SUFFIX={'_days' })] }
                        {suffixname}
                    }
                    {NO
                    }
    [TIMEDURATION={YES** (REFERENCE={CURRENT**
                                }
                                {HMS(timespec)})
                    UNIT={AUTO**
                        {HOURS[(SUFFIX={'_hours' })] }
                        {suffixname}
                        {MINUTES[(SUFFIX={'_minutes' })] }
                        {suffixname}
                        {SECONDS[(SUFFIX={'_seconds' })] }
                        {suffixname}
                    }
                    {NO
                    }
    [EXTRACTYEAR={YES[(SUFFIX={'_year' })] }
                    {suffixname}
                    {NO**
                    }
    [EXTRACTMONTH={YES[(SUFFIX={'_month' })] }
                    {suffixname}
                    {NO**
                    }
    [EXTRACTDAY={YES(SUFFIX={'_day' })]
                    {suffixname}
                    {NO**
                    }
    [EXTRACTHOUR={YES(SUFFIX={'_hour' })]
                    {suffixname}
                    {NO**
                    }
    [EXTRACTMINUTE={YES(SUFFIX={'_minute' })]
                    {suffixname}
                    {NO**
                    }
    [EXTRACTSECOND={YES(SUFFIX={'_second' })]
                    {suffixname}
                    {NO**
                    }
  [/SCREENING [PCTMISSING={YES** (MAXPCT={50**})}
                    {value}
                    {NO
                    }
    [UNIQUECAT={YES** (MAXCAT={100** })}
                    {integer}
                    {NO
                    }
    [SINGLECAT={YES (MAXPCT={95** })}
                    {value}
                    {NO**
                    }
  [/ADJUSTLEVEL [INPUT={YES**}] [TARGET={YES**}]
    {NO } {NO }
    [MAXVALORDINAL={10** }
                    {integer}
    [MINVALCONTINUOUS={5** }
                    {integer}
  [/OUTLIERHANDLING [INPUT={YES**}] [TARGET={YES**}]
    {NO } {NO }
    [CUTOFF=SD({3** })]
                    {value}
    [REPLACEWITH={CUTOFFVALUE**}]
                    {MISSING
                    }
  [/REPLACEMISSING [INPUT={YES** [(EXCLUDE (CONTINUOUS) [NOMINAL] [ORDINAL])]}]
```

```

                {NO
                [TARGET={YES [(EXCLUDE ([CONTINUOUS] [NOMINAL] [ORDINAL]))]}]}
                {NO**
                }
[/REORDERNOMINAL [INPUT={YES } ] [TARGET={YES } ]
                {NO**}                {NO**}
[/RESCALE [INPUT={ZSCORE** ([MEAN={0** } ] [SD={1** } ])}]
                {value}                {value}
                {MINMAX ([MIN={0** } ] [MAX={100** } ])}
                {value}                {value}
                {NONE
                }
                [TARGET={BOXCOX** ([MEAN={0** } ] [SD={1** } ])}]
                {value}                {value}
                {NONE
                }
[/TRANSFORM [MERGESUPERVISED={YES** (PVALUE={0.05**})}]
                {value }
                {NO
                }
                [MERGEUNSUPERVISED={YES{([ORDINAL] [NOMINAL] [MINPCT={10** } ])}]}
                {value}
                {NO**
                }
                [BINNING={SUPERVISED** (PVALUE={0.05**})}]
                {value }
                {NONE
                }
                [SELECTION={YES** (PVALUE={0.05**})}]
                {NO
                }
                [CONSTRUCTION={YES } (ROOT={feature } )]
                {rootname}
                {NO**}
[/CRITERIA [SUFFIX (TARGET={'_transformed'} INPUT={'_transformed'})]
                {suffixname }                {suffixname }
/OUTFILE PREPXML='filespec'

```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### Release History

Release 18

- Command introduced.

### Example

```

ADP
/FIELDS TARGET=targetVar
      INPUT=var1 var2 var3
/OUTFILE PREPXML='file.xml'.

```

## Overview

Automated Data Preparation helps to prepare data for analysis by automating tedious and repetitive data preparation tasks that would otherwise be done manually. The operations it performs improve analysis speed, predictive power, and robustness. A key capability of the component is feature space construction—the discovery of useful sets of predictors from the data through transformation and combination of existing fields. Feature *selection* offers the ability to narrow the attribute space by screening out irrelevant fields, but Automated Data Preparation pairs selection and construction capabilities in order to automatically remove irrelevant fields that slow down or confuse algorithms and create new fields that boost predictive power.

Note that supported operations are performed without knowing what algorithms will be run on the data in further analyses—it is not a generalized data cleaner, nor does it have an understanding of business rules. Basic cleaning and integrity checks can be done using the IBM® SPSS® Statistics Data Validation procedure.

### **Options**

**Date and Time Handling.** The year, month, and day can be extracted from fields containing dates, and new fields containing the durations since a reference date computed. Likewise, the hour, minute, and second can be extracted from fields containing times, and new fields containing the time since a reference time computed.

**Screening.** Fields with too many missing values, and categorical fields with too many unique values, or too many values concentrated in a single value, can be screened and excluded from further analysis.

**Rescaling.** Continuous inputs can optionally be rescaled using a *z* score or min-max transformation. A continuous target can optionally be rescaled using a Box-Cox transformation.

**Transformations.** The procedure can suggest transformations used to merge similar categories of categorical inputs, bin values of continuous inputs, and construct and select new input fields from continuous inputs using principal components analysis.

**Other Target and Input Handling.** The procedure can apply rules for handling outliers, replace missing values, recode the categories of nominal fields, and adjust the measurement level of continuous and ordinal fields.

**Output.** The procedure creates an XML file containing suggested operations. This can be merged with a model XML file using the Merge Model XML dialog (Utilities>Merge Model XML) or transformed into command syntax using `TMS IMPORT`.

### **Basic Specification**

The basic specification is the `ADP` command with a `FIELDS` subcommand specifying the inputs and optionally a target, and an `OUTFILE` subcommand specifying where the transformation rules should be saved.

### **Syntax Rules**

- The `VARIABLES` and `OUTFILE` subcommands are required; all other subcommands are optional.
- Subcommands may be specified in any order.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not allowed.

**Limitations**

- SPLIT FILE is ignored by this command.

**Examples****Basic Specification**

```
ADP
  /FIELDS TARGET=targetField
      INPUT=field1 field2 field3
  /OUTFILE PREPXML='file.xml'.
```

- ADP processes the target and input fields using the default settings.
- OUTFILE saves an XML file containing suggested operations. This can be merged with a model XML file using the Merge Model XML dialog (Utilities>Merge Model XML) or transformed into command syntax using TMS IMPORT.

**Using ADP with TMS IMPORT**

```
*Automatic Data Preparation.
ADP
  /FIELDS TARGET=sales INPUT=resale type price engine_s horsepower wheelbas width length
  curb_wgt fuel_cap mpg
  /PREPDATE TIME DATEDURATION=YES (REFERENCE=YMD('2009-06-04') UNIT=AUTO)
  TIMEDURATION=YES (REFERENCE=HMS('08:43:35') UNIT=AUTO) EXTRACTYEAR=YES (SUFFIX='_year')
  EXTRACTMONTH=YES (SUFFIX='_month') EXTRACTDAY=YES (SUFFIX='_day')
  EXTRACTHOUR=YES (SUFFIX='_hour') EXTRACTMINUTE=YES (SUFFIX='_minute')
  EXTRACTSECOND=YES (SUFFIX='_second')
  /SCREENING PCTMISSING=YES (MAXPCT=50) UNIQUECAT=YES (MAXCAT=100) SINGLECAT=NO
  /ADJUSTLEVEL INPUT=YES TARGET=YES MAXVALORDINAL=10 MINVALCONTINUOUS=5
  /OUTLIERHANDLING INPUT=YES TARGET=NO CUTOFF=SD(3) REPLACEWITH=CUTOFFVALUE
  /REPLACEMISSING INPUT=YES TARGET=NO
  /REORDERNOMINAL INPUT=YES TARGET=NO
  /RESCALE INPUT=ZSCORE (MEAN=0 SD=1) TARGET=BOXCOX (MEAN=0 SD=1)
  /TRANSFORM MERGESUPERVISED=NO MERGEUNSUPERVISED=NO BINNING=NONE SELECTION=NO
  CONSTRUCTION=NO
  /CRITERIA SUFFIX (TARGET='_transformed' INPUT='_transformed')
  /OUTFILE PREPXML='/workingDirectory/car_sales_transformations.xml'.
TMS IMPORT
  /INFILE TRANSFORMATIONS='/workingDirectory/car_sales_transformations.xml'
  MODE=FORWARD (ROLES=UPDATE)
  /SAVE TRANSFORMED=YES.
EXECUTE.
```

- The ADP command prepares the target field *sales* and the input fields *resale* through *mpg*.
- The PREPDATE TIME subcommand is specified, but not used because none of the fields are date or time fields.
- The ADJUSTLEVEL subcommand recasts ordinal fields with more than 10 values as continuous and continuous fields with fewer than 5 values as ordinal.
- The OUTLIERHANDLING subcommand replaces values of continuous inputs (not the target) that are more than 3 standard deviations from the mean with the value that is 3 standard deviations from the mean.
- The REPLACEMISSING subcommand replaces values of inputs (not the target) that are missing.

- The `REORDERNOMINAL` subcommand recodes the values of nominal inputs from least frequently occurring to most frequently occurring.
- The `RESCALE` subcommand standardizes continuous inputs to have mean 0 and standard deviation 1 using a z-score transformation, and standardizes the continuous target to have mean 0 and standard deviation 1 using a Box-Cox transformation.
- The `TRANSFORM` subcommand turns off all default operations specified by this subcommand.
- The `CRITERIA` subcommand specifies the default suffixes for transformations of the target and inputs.
- The `OUTFILE` subcommand specifies that the transformations should be saved to `/workingDirectory/car_sales_transformations.xml`, where `/workingDirectory` is the path to where you want to save `car_sales_transformations.xml`.
- The `TMS IMPORT` command reads the transformations in `car_sales_transformations.xml` and applies them to the active dataset, updating the roles of existing fields that are transformed.
- The `EXECUTE` command causes the transformations to be processed. When using this as part of a longer stream of syntax, you may be able to remove the `EXECUTE` command to save some processing time.

## ***FIELDS Subcommand***

The `FIELDS` subcommand is used to specify the target, inputs, and optional weights.

- The `FIELDS` subcommand and the `INPUT` keyword are required.

### ***TARGET Keyword***

Specify a single field that will be used as a target in further analyses. The target field is processed based upon its defined measurement level; nominal, ordinal, or continuous. Use the `VARIABLE LEVEL` command to change a target field's measurement level.

### ***INPUT Keyword***

Specify one or more fields that will be used as inputs in further analyses. Input fields are processed based upon their defined measurement level; nominal, ordinal, or continuous. Use the `VARIABLE LEVEL` command to change an input field's measurement level.

### ***ANALYSISWEIGHT Keyword***

Specify a variable containing analysis (regression) weights. The procedure incorporates analysis weights where appropriate in operations used to prepare the data.

The analysis weight variable must be numeric. Cases with a negative or zero analysis weight are ignored.

## ***PREPDATE TIME Subcommand***

The `PREPDATE TIME` subcommand specifies handling of date and time fields.

- If `PREPDATE` is not specified, by default the procedure computes date and time durations since the current date and time.
- The original date and time fields will not be recommended as model inputs following automated data preparation.

### **DATEDURATION Keyword**

The `DATEDURATION` keyword computes the number of years/months/days since a reference date for each variable containing dates.

#### **REFERENCE=CURRENT | YMD('datespec')**

*Reference date.* Specify `CURRENT` to use the current date as the reference date. Use `YMD` to specify a custom reference date with the year, month, and day, in that order, in parentheses using a valid date format in quotes. The default is `CURRENT`.

#### **UNIT=AUTO | YEARS | MONTHS | DAYS**

*Date units for computed durations.* Specify the units for the computed durations. `AUTO` determines the units based on the following rules. The default is `AUTO`.

- If the minimum number of elapsed days is less than 31, then the duration is returned in days.
- If the minimum number of elapsed days is less than 366 but greater than or equal to 31, then the duration is returned in months. The number of months between two dates is calculated based on average number of days in a month (30.4375):  $months = days / 30.4375$ .
- If the minimum number of elapsed days greater than or equal to 366, then the duration is returned in years. The number of years between two dates is calculated based on average number of days in a year (365.25):  $years = days / 365.25$ .

Explicitly specifying `YEARS`, `MONTHS`, or `DAYS` returns the duration in years, months, or days, respectively. Optionally, in parentheses, specify `SUFFIX=` with a suffix in quotes. The default suffix depends upon the unit; `YEARS`, `MONTHS`, and `DAYS` have defaults `_years`, `_months`, and `_days`, respectively.

### **TIMEDURATION Keyword**

The `TIMEDURATION` keyword computes the number of hours/minutes/seconds since a reference time for each variable containing times.

#### **REFERENCE=CURRENT | HMS('timespec')**

*Reference date.* Specify `CURRENT` to use the current time as the reference time or use `HMS` and the hour, minute, and second, in that order, in parentheses using a valid time format in quotes. The default is `CURRENT`.

#### **UNIT = AUTO | HOURS | MINUTES | SECONDS**

*Date units for computed durations.* Specify the units for the computed durations. `AUTO` determines the units based on the following rules. The default is `AUTO`.

- If the minimum number of elapsed seconds is less than 60, then the duration is returned in seconds.
- If the minimum number of elapsed seconds is larger than or equal to 60 but less than 3600, then the duration is returned in minutes.
- If the minimum number of elapsed seconds is larger than or equal to 3600, then the duration is returned in hours.

Explicitly specifying `HOURS`, `MINUTES`, or `SECONDS` returns the duration in hours, minutes, or seconds, respectively. Optionally, in parentheses, specify `SUFFIX=` with a suffix in quotes. The default suffix depends upon the unit; `HOURS`, `MINUTES`, and `SECONDS` have defaults `_hours`, `_minutes`, and `_seconds`, respectively.

### ***EXTRACTYEAR Keyword***

The `EXTRACTYEAR` keyword extracts the year element from a date variable.

Optionally specify the `SUFFIX` keyword in parentheses with a suffix in quotes. The default suffix is `_year`.

### ***EXTRACTMONTH Keyword***

The `EXTRACTMONTH` keyword extracts the month element from a date variable.

Optionally specify the `SUFFIX` keyword in parentheses with a suffix in quotes. The default suffix is `_month`.

### ***EXTRACTDAY Keyword***

The `EXTRACTDAY` keyword extracts the day element from a date variable.

Optionally specify the `SUFFIX` keyword in parentheses with a suffix in quotes. The default suffix is `_day`.

### ***EXTRACTHOUR Keyword***

The `EXTRACTHOUR` keyword extracts the hour element from a time variable.

Optionally specify the `SUFFIX` keyword in parentheses with a suffix in quotes. The default suffix is `_hour`.

### ***EXTRACTMINUTE Keyword***

The `EXTRACTMINUTE` keyword extracts the minute element from a time variable.

Optionally specify the `SUFFIX` keyword in parentheses with a suffix in quotes. The default suffix is `_minute`.

### ***EXTRACTSECOND Keyword***

The `EXTRACTSECOND` keyword extracts the second element from a time variable.

Optionally specify the `SUFFIX` keyword in parentheses with a suffix in quotes. The default suffix is `_second`.

## **SCREENING Subcommand**

The `SCREENING` subcommand specifies settings for excluding unsuitable fields.

**PCTMISSING=YES(MAXPCT=value) | NO**

*Screen out fields with too many missing values.* Fields with more than `MAXPCT` missing values are removed from further analysis. Specify a value greater than or equal to 0, which is equivalent to deselecting this option, and less than or equal to 100, though fields with all missing values are automatically excluded. The default is 50.

**UNIQUECAT=YES(MAXCAT=integer) | NO**

*Screen out nominal fields with too many unique categories.* Nominal fields with more than `MAXCAT` categories are removed from further analysis. Specify a positive integer. The default is 100.

**SINGLECAT=YES(MAXPCT=value) | NO**

*Screen out categorical fields that are nearly constant.* Ordinal and nominal fields with a category that contains more than `MAXPCT` of the records are removed from further analysis. Specify a value greater than or equal to 0, equivalent to deselecting this option, and less than or equal to 100, though constant fields are automatically excluded. The default is 95.

## **ADJUSTLEVEL Subcommand**

The `ADJUSTLEVEL` subcommand recasts ordinal fields with too many categories as continuous and continuous fields with too few unique values as ordinal. By default, the measurement levels of ordinal fields with more than 10 categories and continuous fields with fewer than 5 unique values are adjusted.

**INPUT=YES | NO**

*Check inputs and adjust measurement level if necessary.* By default, inputs are checked.

**TARGET=YES | NO**

*Check target and adjust measurement level if necessary.* By default, the target is checked.

**MAXVALORDINAL=integer**

*Maximum number of categories allowed for ordinal fields.* Ordinal fields with more than `MAXVALORDINAL` categories are recast as continuous fields. Specify a positive integer. The default is 10. The value of `MAXVALORDINAL` must be greater than or equal to `MINVALCONTINUOUS`.

**MINVALCONTINUOUS= integer**

*Minimum number of unique values allowed for continuous fields.* Continuous fields with less than `MINVALCONTINUOUS` unique values are recast as ordinal fields. Specify a positive integer. The default is 5. The value of `MINVALCONTINUOUS` must be less than or equal to `MAXVALORDINAL`.



## **OUTLIERHANDLING Subcommand**

The `OUTLIERHANDLING` subcommand checks fields for outliers and replaces the outlying values with less extreme values or missing values.

**INPUT=YES | NO**

*Check inputs for outliers.* By default, inputs are checked.

**TARGET=YES | NO**

*Check target for outliers.* By default, the target is checked.

**CUTOFF=SD(value)**

*Cutoff for determining outliers.* If a value is more than SD “robust” standard deviations from the mean value for a field, then it is considered an outlier. Specify a positive number. The default is 3 standard deviations.

**REPLACEWITH=CUTOFFVALUE | MISSING**

*Value to replace outliers with.* `CUTOFFVALUE` replaces outliers with the cutoff for determining outliers. `MISSING` replaces outliers with the system-missing value. These missing values can be further handled by the `REPLACEMISSING` subcommand. The default is `CUTOFFVALUE`.

## **REPLACEMISSING Subcommand**

The `REPLACEMISSING` subcommand replaces missing values in continuous, ordinal, and nominal fields with the mean, median, or mode, respectively.

**INPUT=YES | NO**

*Replace missing values in input fields.* By default, missing values are replaced in inputs. Optionally specify the keyword `EXCLUDE` and a list in parentheses of the field measurement levels to determine which input fields should be checked; for example:

`INPUT=YES`

causes the procedure to replace missing values in all input fields, while

`INPUT=YES (EXCLUDE (CONTINUOUS NOMINAL) )`

causes the procedure to replace missing values in fields with the ordinal measurement level.

**TARGET= NO | YES**

*Replace missing values in the target.* By default, missing values are not replaced in the target. When replacing missing values in the target, optionally specify the keyword `EXCLUDE` as described for the `INPUT` keyword above.

## **REORDERNOMINAL Subcommand**

The `REORDERNOMINAL` subcommand recodes the values of nominal fields from least frequently occurring to most frequently occurring. The new field values start with 0 as the least frequent category. Note that the new field will be numeric even if the original field is a string. For example,

if a nominal field's data values are "A", "A", "A", "B", "C", "C", then automated data preparation would recode "B" into 0, "C" into 1, and "A" into 2.

**INPUT=NO | YES**

*Reorder values of inputs.* By default, values of nominal inputs are not reordered. This specification is ignored if there are no nominal inputs.

**TARGET=NO | YES**

*Reorder values of the target.* By default, values of a nominal target are not reordered. This specification is ignored if the target is not nominal.

## **RESCALE Subcommand**

The RESCALE subcommand is used to rescale continuous fields. Different methods are available for inputs and the target.

### **INPUT Keyword**

The INPUT keyword specifies the method for rescaling continuous input fields.

- Z score rescaling is performed by default with a mean of 0 and standard deviation of 1.
- If there are no continuous inputs, INPUT is ignored.

**ZSCORE(MEAN=value SD=value)**

*Z score transformation.* Using the observed mean and standard deviation as population parameter estimates, the fields are standardized and then the z scores are mapped to the corresponding values of a normal distribution with the specified MEAN and SD. Specify a number for MEAN and a positive number for SD. The defaults are 0 and 1, respectively, corresponding to standardized rescaling.

**MINMAX(MIN=value MAX=value)**

*Min-Max transformation.* Using the observed minimum and maximum as population parameter estimates, the fields are mapped to the corresponding values of a uniform distribution with the specified MIN and MAX. Specify numbers with MAX greater than MIN.

**NONE**

*Do not rescale inputs.*

### **TARGET Keyword**

The TARGET keyword specifies the method for rescaling a continuous target.

- Box-Cox rescaling is performed by default with a target mean of 0 and target standard deviation of 1.
- If there is no target, or it is not continuous, TARGET is ignored.

**BOXCOX(MEAN=value SD=value)**

*Box-Cox transformation.* This transforms a continuous target using the Box-Cox transformation into a field that has an approximately normal distribution with the specified MEAN and SD. Specify a number for MEAN and a positive number for SD. The defaults are 0 and 1, respectively.

NONE

*Do not rescale target.*

## **TRANSFORM Subcommand**

The TRANSFORM subcommand is used to merge similar categories of categorical inputs, bin values of continuous inputs, and construct and select new input fields from continuous inputs using principal components analysis.

### **MERGESUPERVISED Keyword**

The MERGESUPERVISED keyword specifies how to merge similar categories of a nominal or ordinal input in the presence of a target.

- If there are no categorical inputs, MERGESUPERVISED is ignored.
- If there is no target specified on the FIELDS subcommand, MERGESUPERVISED is ignored.

**YES(PVALUE=value)**

*Supervised merge.* Similar categories are identified based upon the relationship between the input and the target. Categories that are not significantly different; that is, having a  $p$ -value greater than the value of PVALUE, are merged. Specify a value greater than 0 and less than or equal to 1. The default is 0.05. YES is the default.

**NO**

*Do not merge categories.*

### **MERGEUNSUPERVISED Keyword**

The MERGEUNSUPERVISED keyword specifies how to merge similar categories of a nominal or ordinal input when there is no target.

- If there are no categorical inputs, MERGEUNSUPERVISED is ignored.
- If there is a target specified on the FIELDS subcommand, MERGEUNSUPERVISED is ignored.

**YES(ORDINAL NOMINAL MINPCT=value)**

*Unsupervised merge.* The equal frequency method is used to merge categories with less than MINPCT of the total number of records. Specify a value greater than or equal to 0 and less than or equal to 100. The default is 10 if MINPCT is not specified. If YES is specified without ORDINAL or NOMINAL, then no merging is performed.

**NO**

*Do not merge categories. NO is the default.*

**BINNING Keyword**

The `BINNING` keyword specifies how to discretize continuous inputs in the presence of a categorical target.

**SUPERVISED(PVALUE=value)**

*Supervised binning.* Bins are created based upon the properties of “homogeneous subsets”, which are identified by the Scheffe method using `PVALUE` as the alpha for the critical value for determining homogeneous subsets. `SUPERVISED` is the default. Specify a value greater than 0 and less than or equal to 1. The default is 0.05.

If there is no target specified on the `FIELDS` subcommand, or the target is not categorical, or there are no continuous inputs, then `SUPERVISED` is ignored.

**NONE**

*Do not bin values of continuous inputs.*

**SELECTION Keyword**

The `SELECTION` keyword specifies how to perform feature selection for continuous inputs in the presence of a continuous target.

**YES(PVALUE=value)**

*Perform feature selection.* A continuous input is removed from the analysis if the  $p$ -value for its correlation with the target is greater than `PVALUE`. `YES` is the default.

If there is no target specified on the `FIELDS` subcommand, or the target is not continuous, or there are no continuous inputs, then `YES` is ignored.

**NO**

*Do not perform feature selection.*

**CONSTRUCTION Keyword**

The `CONSTRUCTION` keyword specifies how to perform feature construction for continuous inputs in the presence of a continuous target.

**YES(ROOT=rootname)**

*Perform feature construction.* New predictors are constructed from groups of “similar” predictors using principal component analysis. Optionally specify the rootname for constructed predictors using `ROOT` in parentheses. Specify a rootname (no quotes). The default is *feature*.

If there is no target specified on the `FIELDS` subcommand, or the target is not continuous, or there are no continuous inputs, then `YES` is ignored.

**NO**

*Do not perform feature construction.* `NO` is the default.

**CRITERIA Subcommand**

The `CRITERIA` subcommand is used to specify the suffixes applied to transformed target and inputs.

### **SUFFIX Keyword**

The `SUFFIX` keyword specifies the suffixes applied to transformed target and inputs.

#### **TARGET=suffixname**

*Suffix for transformed target.* Specify a suffix in quotes. The default is `_transformed`.

If there is no target specified on the `FIELDS` subcommand, `TARGET` is ignored.

#### **INPUT=suffixname**

*Suffix for transformed inputs.* Specify a suffix in quotes. The default is `_transformed`.

## **OUTFILE Subcommand**

The `OUTFILE` subcommand saves an XML-format file containing the rules for preparing the data.

- The `OUTFILE` subcommand is required.
- File names must be specified in full. `ADP` does not supply extensions.

#### **PREPXML='filespec'**

*Save rules for preparing data to an XML file.* The rules are saved in an XML format to the specified file. This file can be merged with model PMML using `TMS MERGE` or transformed into command syntax using `TMS IMPORT`

# AGGREGATE

```
AGGREGATE [OUTFILE={'savfile'|'dataset'}
          {*          } [MODE={REPLACE  } ] [OVERWRITE={NO  }]
          {ADDVARIABLES}                {YES}

[/MISSING=COLUMNWISE] [/DOCUMENT]

[/PRESORTED] [/BREAK=[varlist[({A})]][varlist...]]
              {D}

/aggvar['label'] aggvar['label']...=function(arguments)

[/aggvar ...]
```

## Available functions:

SUM	Sum	MEAN	Mean
SD	Standard deviation	MAX	Maximum
MIN	Minimum	PGT	% of cases greater than value
PLT	% of cases less than value	PIN	% of cases between values
POUT	% of cases not in range	FGT	Fraction greater than value
FLT	Fraction less than value	FIN	Fraction between values
FOUT	Fraction not in range	N	Weighted number of cases
NU	Unweighted number of cases	NMISS	Weighted number of missing cases
NUMISS	Unweighted number of missing cases	FIRST	First nonmissing value
LAST	Last nonmissing value	MEDIAN	Median

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

### Release 13.0

- MODE keyword introduced.
- OVERWRITE keyword introduced.

### Release 17.0

- AGGREGATE runs without a break variable.

## Example

```
AGGREGATE
  /OUTFILE=' /temp/temp.sav'
  /BREAK=gender
  /age_mean=MEAN(age) .
```

## Overview

AGGREGATE aggregates groups of cases in the active dataset into single cases and creates a new aggregated file or creates new variables in the active dataset that contain aggregated data. The values of one or more variables in the active dataset define the case groups. These variables are called **break variables**. A set of cases with identical values for each break variable is called a **break group**. If no break variables are specified, then the entire dataset is a single break group. Aggregate functions are applied to **source variables** in the active dataset to create new aggregated variables that have one value for each break group.

### Options

**Data.** You can create new variables in the active dataset that contain aggregated data, replace the active dataset with aggregated results, or create a new data file that contains the aggregated results.

**Documentary Text.** You can copy documentary text from the original file into the aggregated file using the `DOCUMENT` subcommand. By default, documentary text is dropped.

**Aggregated Variables.** You can create aggregated variables using any of 19 aggregate functions. The functions `SUM`, `MEAN`, and `SD` can aggregate only numeric variables. All other functions can use both numeric and string variables.

**Labels and Formats.** You can specify variable labels for the aggregated variables. Variables created with the functions `MAX`, `MIN`, `FIRST`, and `LAST` assume the formats and value labels of their source variables. All other variables assume the default formats described under Aggregate Functions on p. 166.

### Basic Specification

The basic specification is at least one aggregate function and source variable. The aggregate function creates a new aggregated variable in the active dataset.

### Subcommand Order

- If specified, `OUTFILE` must be specified first.
- If specified, `DOCUMENT` and `PRESORTED` must precede `BREAK`. No other subcommand can be specified between these two subcommands.
- `MISSING`, if specified, must immediately follow `OUTFILE`.
- The aggregate functions must be specified last.

### Operations

- When replacing the active dataset or creating a new data file, the aggregated file contains the break variables plus the variables created by the aggregate functions.
- `AGGREGATE` excludes cases with missing values from all aggregate calculations except those involving the functions `N`, `NU`, `NMISS`, and `NUMISS`.
- Unless otherwise specified, `AGGREGATE` sorts cases in the aggregated file in ascending order of the values of the grouping variables.

- `PRESORTED` uses a faster, less memory-intensive algorithm that assumes the data are already sorted into the desired groups.
- `AGGREGATE` ignores split-file processing. To achieve the same effect, name the variable or variables used to split the file as break variables before any other break variables. `AGGREGATE` produces one file, but the aggregated cases will then be in the same order as the split files.

## Example

```
AGGREGATE
  /OUTFILE=' /temp/temp.sav'
  /BREAK=gender marital
  /age_mean=MEAN(age)
  /age_median=MEDIAN(age)
  /income_median=MEDIAN(income) .
```

- `AGGREGATE` creates a new IBM® SPSS® Statistics data file, *temp.sav*, that contains two break variables (*gender* and *marital*) and all of the new aggregate variables.
- `BREAK` specifies *gender* and *marital* as the break variables. In the aggregated file, cases are sorted in ascending order of *gender* and in ascending order of *marital* within *gender*. The active dataset remains unsorted.
- Three aggregated variables are created: *age\_mean* contains the mean age for each group defined by the two break variables; *age\_median* contains the median age; and *income\_median* contains the median income.

## OUTFILE Subcommand

`OUTFILE` specifies the handling of the aggregated results. It must be the first subcommand on the `AGGREGATE` command.

- `OUTFILE='file specification'` saves the aggregated data to a new file, leaving the active dataset unaffected. The file contains the new aggregated variables and the break variables that define the aggregated cases.
- A defined dataset name can be used for the file specification, saving the aggregated data to a dataset in the current session. The dataset must be defined before being used in the `AGGREGATE` command. [For more information, see the topic DATASET DECLARE on p. 580.](#)
- `OUTFILE=*` with no additional keywords on the `OUTFILE` subcommand will replace the active dataset with the aggregated results.
- `OUTFILE=* MODE=ADDVARIABLES` appends the new variables with the aggregated data to the active dataset (instead of replacing the active dataset with the aggregated data).
- `OUTFILE=* MODE=ADDVARIABLES OVERWRITE=YES` overwrites variables in the active dataset if those variable names are the same as the aggregate variable names specified on the `AGGREGATE` command.
- `MODE` and `OVERWRITE` can be used only with `OUTFILE=*`; they are invalid with `OUTFILE='file specification'`.
- Omission of the `OUTFILE` subcommand is equivalent to `OUTFILE=* MODE=ADDVARIABLES`.



**Example**

```
AGGREGATE
  /BREAK=region
  /sales_mean = MEAN(var1)
  /sales_median = MEDIAN(var1)
  /sales_sum = SUM(var1).
```

- The aggregated variables are appended to the end of each case in the active data file. No existing cases or variables are deleted.
- For each case, the new aggregated variable values represent the mean, median, and total (sum) sales values for its region.

**Creating a New Aggregated Data File versus Appending Aggregated Variables**

When you create a new aggregated data file with `OUTFILE='file specification'` or `OUTFILE=* MODE=REPLACE`, the new file contains:

- The break variables from the original data file and the new aggregate variables defined by the aggregate functions. Original variables other than the break variables are not retained.
- One case for each group defined by the break variables. If there is one break variable with two values, the new data file will contain only two cases.

When you append aggregate variables to the active dataset with `OUTFILE=* MODE=ADDVARIABLES`, the modified data file contains:

- All of the original variables plus all of the new variables defined by the aggregate functions, with the aggregate variables appended to the end of each case.
- The same number of cases as the original data file. The data file itself is not aggregated. Each case with the same value(s) of the break variable(s) receives the same values for the new aggregate variables. For example, if *gender* is the only break variable, all males would receive the same value for a new aggregate variable that represents the average age.

**Example**

```
DATA LIST FREE /age (F2) gender (F2).
BEGIN DATA
25 1
35 1
20 2
30 2
60 2
END DATA.
*create new file with aggregated results.
AGGREGATE
  /OUTFILE=' /temp/temp.sav'
  /BREAK=gender
  /age_mean=MEAN(age)
  /groupSize=N.
*append aggregated variables to active dataset.
AGGREGATE
  /OUTFILE=* MODE=ADDVARIABLES
  /BREAK=gender
  /age_mean=MEAN(age)
  /groupSize=N.
```

Figure 9-1  
New aggregated data file

	gender	age_mean	groupSize	var	var	va
1	1	30.00	2			
2	2	36.67	3			
3						
4						
5						
6						

Figure 9-2  
Aggregate variables appended to active dataset

	age	gender	age_mean	groupSize	var	va
1	25	1	30.00	2		
2	35	1	30.00	2		
3	20	2	36.67	3		
4	30	2	36.67	3		
5	60	2	36.67	3		
6						

## ***BREAK Subcommand***

BREAK lists the optional grouping variables, also called break variables. Each unique combination of values of the break variables defines one break group.

- The variables named on BREAK can be any combination of variables in the active dataset.
- Unless PRESORTED is specified, aggregated variables are appended to the active dataset (OUTFILE=\* MODE=ADDVARIABLES), AGGREGATE sorts cases after aggregating. By default, cases are sorted in ascending order of the values of the break variables. AGGREGATE sorts first on the first break variable, then on the second break variable within the groups created by the first, and so on.
- Sort order can be controlled by specifying an A (for ascending) or D (for descending) in parentheses after any break variables.
- The designations A and D apply to all preceding undesignated variables.
- The subcommand PRESORTED overrides all sorting specifications, and no sorting is performed with OUTFILE=\* MODE=ADDVARIABLES.

```
AGGREGATE
  /BREAK=region
  /sales_mean = MEAN(var1)
  /sales_median = MEDIAN(var1)
  /sales_sum = SUM(var1).
```

For each case, the new aggregated variable values represent the mean, median, and total (sum) sales values for its region.

```
AGGREGATE
  /sales_mean = MEAN(var1)
  /sales_median = MEDIAN(var1)
  /sales_sum = SUM(var1).
```

For each case, the new aggregated variable values represent the mean, median, and total (sum) sales values for the entire dataset.

## ***DOCUMENT Subcommand***

DOCUMENT copies documentation from the original file into the aggregated file.

- DOCUMENT must appear after OUTFILE but before BREAK.
- By default, documents from the original data file are not retained with the aggregated data file when creating a new aggregated data file with either OUTFILE='file specification' or OUTFILE=\* MODE=REPLACE. The DOCUMENT subcommand retains the original data file documents.
- Appending variables with OUTFILE=\* MODE=ADDVARIABLES has no effect on data file documents, and the DOCUMENT subcommand is ignored. If the data file previously had documents, they are retained.

## ***PRESORTED Subcommand***

If the data are already sorted into the desired groups, you can reduce run time and memory requirements by using the PRESORTED subcommand.

- If specified, PRESORTED must precede BREAK. The only specification is the keyword PRESORTED. PRESORTED has no additional specifications.
- When PRESORTED is specified, the program forms an aggregate case out of each group of *adjacent* cases with the same values for the break variables. Unless the cases are sorted by the break variables, the results will be quite different from what would be produced if PRESORTED were not specified.
- When PRESORTED is specified, if AGGREGATE is appending new variables to the active dataset rather than writing a new file or replacing the active dataset, the cases must be sorted in ascending order by the BREAK variables.

### ***Example***

```
AGGREGATE OUTFILE=' /temp/temp.sav'
  /PRESORTED
  /BREAK=gender marital
  /mean_age=MEAN(age).
```

## Aggregate Functions

An aggregated variable is created by applying an aggregate function to a variable in the active dataset. The variable in the active dataset is called the **source** variable, and the new aggregated variable is the **target** variable.

- The aggregate functions must be specified last on AGGREGATE.
- The simplest specification is a target variable list, followed by an equals sign, a function name, and a list of source variables.
- The number of target variables named must match the number of source variables.
- When several aggregate variables are defined at once, the first-named target variable is based on the first-named source variable, the second-named target is based on the second-named source, and so on.
- Only the functions MAX, MIN, FIRST, and LAST copy complete dictionary information from the source variable. For all other functions, new variables do not have labels and are assigned default dictionary print and write formats. The default format for a variable depends on the function used to create it (see the list of available functions below).
- You can provide a variable label for a new variable by specifying the label in single or double quotes immediately following the new variable name. Value labels cannot be assigned in AGGREGATE.
- To change formats or add value labels to an active dataset created by AGGREGATE, use the PRINT FORMATS, WRITE FORMATS, FORMATS, or VALUE LABELS command. If the aggregate file is written to disk, first retrieve the file using GET, specify the new labels and formats, and resave the file.

The following is a list of available functions:

<b>SUM(varlist)</b>	<i>Sum across cases.</i> Default formats are F8.2.
<b>MEAN(varlist)</b>	<i>Mean across cases.</i> Default formats are F8.2.
<b>MEDIAN(varlist)</b>	<i>Median across cases.</i> Default formats are F8.2.
<b>SD(varlist)</b>	<i>Standard deviation across cases.</i> Default formats are F8.2.
<b>MAX(varlist)</b>	<i>Maximum value across cases.</i> Complete dictionary information is copied from the source variables to the target variables.
<b>MIN(varlist)</b>	<i>Minimum value across cases.</i> Complete dictionary information is copied from the source variables to the target variables.
<b>PGT(varlist,value)</b>	<i>Percentage of cases greater than the specified value.</i> Default formats are F5.1.
<b>PLT(varlist,value)</b>	<i>Percentage of cases less than the specified value.</i> Default formats are F5.1.
<b>PIN(varlist,value1,value2)</b>	<i>Percentage of cases between value1 and value2, inclusive.</i> Default formats are F5.1.
<b>POUT(varlist,value1,value2)</b>	<i>Percentage of cases not between value1 and value2.</i> Cases where the source variable equals value1 or value2 are not counted. Default formats are F5.1.
<b>FGT(varlist,value)</b>	<i>Fraction of cases greater than the specified value.</i> Default formats are F5.3.
<b>FLT(varlist,value)</b>	<i>Fraction of cases less than the specified value.</i> Default formats are F5.3.

<b>FIN(varlist,value1,value2)</b>	<i>Fraction of cases between value1 and value2, inclusive.</i> Default formats are F5.3.
<b>FOUT(varlist,value1,value2)</b>	<i>Fraction of cases not between value1 and value2.</i> Cases where the source variable equals <i>value1</i> or <i>value2</i> are not counted. Default formats are F5.3.
<b>N(varlist)</b>	<i>Weighted number of cases in break group.</i> Default formats are F7.0 for unweighted files and F8.2 for weighted files.
<b>NU(varlist)</b>	<i>Unweighted number of cases in break group.</i> Default formats are F7.0.
<b>NMISS(varlist)</b>	<i>Weighted number of missing cases.</i> Default formats are F7.0 for unweighted files and F8.2 for weighted files.
<b>NUMISS(varlist)</b>	<i>Unweighted number of missing cases.</i> Default formats are F7.0.
<b>FIRST(varlist)</b>	<i>First nonmissing observed value in break group.</i> Complete dictionary information is copied from the source variables to the target variables.
<b>LAST(varlist)</b>	<i>Last nonmissing observed value in break group.</i> Complete dictionary information is copied from the source variables to the target variables.

- The functions SUM, MEAN, and SD can be applied only to numeric source variables. All other functions can use short and long string variables as well as numeric ones.
- The N and NU functions do not require arguments. Without arguments, they return the number of weighted and unweighted valid cases in a break group. If you supply a variable list, they return the number of weighted and unweighted valid cases for the variables specified.
- For several functions, the argument includes values as well as a source variable designation. Either blanks or commas can be used to separate the components of an argument list.
- For PIN, POUT, FIN, and FOUT, the first value should be less than or equal to the second. If the first is greater, AGGREGATE automatically reverses them and prints a warning message. If the two values are equal, PIN and FIN calculate the percentages and fractions of values equal to the argument. POUT and FOUT calculate the percentages and fractions of values not equal to the argument.
- String values specified in an argument should be enclosed in quotes.

### **Using the MEAN Function**

```
AGGREGATE OUTFILE='AGGEMP.SAV' /BREAK=LOCATN
  /AVGSAL 'Average Salary' AVGRAISE = MEAN(SALARY RAISE).
```

- AGGREGATE defines two aggregate variables, *AVGSAL* and *AVGRAISE*.
- *AVGSAL* is the mean of *SALARY* for each break group, and *AVGRAISE* is the mean of *RAISE*.
- The label *Average Salary* is assigned to *AVGSAL*.

### **Using the PLT Function**

```
AGGREGATE OUTFILE=* /BREAK=DEPT
  /LOWVAC, LOWSICK = PLT (VACDAY SICKDAY, 10).
```

- AGGREGATE creates two aggregated variables: *LOWVAC* and *LOWSICK*. *LOWVAC* is the percentage of cases with values less than 10 for *VACDAY*, and *LOWSICK* is the percentage of cases with values less than 10 for *SICKDAY*.

**Using the FIN Function**

```
AGGREGATE OUTFILE='GROUPS.SAV' /BREAK=OCCGROUP
/COLLEGE = FIN(EDUC,13,16).
```

- AGGREGATE creates the variable *COLLEGE*, which is the fraction of cases with 13 to 16 years of education (variable *EDUC*).

**Using the PIN Function**

```
AGGREGATE OUTFILE=* /BREAK=CLASS
/LOCAL = PIN(STATE, 'IL', 'IO').
```

- AGGREGATE creates the variable *LOCAL*, which is the percentage of cases in each break group whose two-letter state code represents Illinois, Indiana, or Iowa. (The abbreviation for Indiana, IN, is between IL and IO in an alphabetical sort sequence.)

**MISSING Subcommand**

By default, AGGREGATE uses all nonmissing values of the source variable to calculate aggregated variables. An aggregated variable will have a missing value only if the source variable is missing for every case in the break group. You can alter the default missing-value treatment by using the MISSING subcommand. You can also specify the inclusion of user-missing values on any function.

- MISSING must immediately follow OUTFILE.
- COLUMNWISE is the only specification available for MISSING.
- If COLUMNWISE is specified, the value of an aggregated variable is missing for a break group if the source variable is missing for any case in the group.
- COLUMNWISE does not affect the calculation of the N, NU, NMISS, or NUMISS functions.
- COLUMNWISE does not apply to break variables. If a break variable has a missing value, cases in that group are processed and the break variable is saved in the file with the missing value. Use SELECT IF if you want to eliminate cases with missing values for the break variables.

**Including Missing Values**

You can force a function to include user-missing values in its calculations by specifying a period after the function name.

- AGGREGATE ignores periods used with the functions N, NU, NMISS, and NUMISS if these functions have no arguments.
- User-missing values are treated as valid when these four functions are followed by a period and have a variable as an argument. NMISS.(AGE) treats user-missing values as valid and thus gives the number of cases for which AGE has the system-missing value only.

The effect of specifying a period with N, NU, NMISS, and NUMISS is illustrated by the following:

```
N = N. = N(AGE) + NMISS(AGE) = N.(AGE) + NMISS.(AGE)
```

```
NU = NU. = NU(AGE) + NUMISS(AGE) = NU.(AGE) + NUMISS.(AGE)
```

- The function `N` (the same as `N.` with no argument) yields a value for each break group that equals the number of cases with valid values (`N(AGE)`) plus the number of cases with user- or system-missing values (`NMISS(AGE)`).
- This in turn equals the number of cases with either valid or user-missing values (`N.(AGE)`) plus the number with system-missing values (`NMISS.(AGE)`).
- The same identities hold for the `NU`, `NMISS`, and `NUMISS` functions.

### Default Treatment of Missing Values

```
AGGREGATE OUTFILE='AGGEMP.SAV' /MISSING=COLUMNWISE /BREAK=LOCATN
  /AVGSAL = MEAN(SALARY) .
```

- `AVGSAL` is missing for an aggregated case if `SALARY` is missing for any case in the break group.

### Including User-Missing Values

```
AGGREGATE OUTFILE=* /BREAK=DEPT
  /LOVAC = PLT.(VACDAY,10) .
```

- `LOVAC` is the percentage of cases within each break group with values less than 10 for `VACDAY`, even if some of those values are defined as user missing.

### Aggregated Values that Retain Missing-Value Status

```
AGGREGATE OUTFILE='CLASS.SAV' /BREAK=GRADE
  /FIRSTAGE = FIRST.(AGE) .
```

- The first value of `AGE` in each break group is assigned to the variable `FIRSTAGE`.
- If the first value of `AGE` in a break group is user missing, that value will be assigned to `FIRSTAGE`. However, the value will retain its missing-value status, since variables created with `FIRST` take dictionary information from their source variables.

## Comparing Missing-Value Treatments

The table below demonstrates the effects of specifying the `MISSING` subcommand and a period after the function name. Each entry in the table is the number of cases used to compute the specified function for the variable `EDUC`, which has 10 nonmissing cases, 5 user-missing cases, and 2 system-missing cases for the group. Note that columnwise treatment produces the same results as the default for every function except the `MEAN` function.

Table 9-1  
Default versus columnwise missing-value treatments

Function	Default	Columnwise
<code>N</code>	17	17
<code>N.</code>	17	17
<code>N(EDUC)</code>	10	10
<code>N.(EDUC)</code>	15	15
<code>MEAN(EDUC)</code>	10	0

---

*AGGREGATE*

<b>Function</b>	<b>Default</b>	<b>Columnwise</b>
MEAN.(EDUC)	15	0
NMISS(EDUC)	7	7
NMISS.(EDUC)	2	2



# AIM

AIM is available in the Statistics Base option.

AIM grouping-var

```
[/CATEGORICAL varlist]
[/CONTINUOUS varlist]
[/CRITERIA [ADJUST = {BONFERRONI**}] [CI = {95**}]
           {NONE           }           {value}
           [HIDENOTSIG = {NO**}] [SHOWREFLINE = {NO } ] ]
           {YES }                 {YES**}
[/MISSING {EXCLUDE**} ]
         {INCLUDE }
[/PLOT [CATEGORY] [CLUSTER [(TYPE = {BAR*})]] [ERRORBAR]
      {PIE }
      [IMPORTANCE [(X = {GROUP*}] [Y = {TEST*}])] ]
      {VARIABLE}           {PVALUE}
```

\* Default if the keyword is omitted.

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
AIM TSC_1
  /CATEGORICAL type
  /CONTINUOUS price engine_s horsepow wheelbas width length
              curb_wgt fuel_cap mpg
  /PLOT CLUSTER.
```

## Overview

AIM provides graphical output to show the relative importance of categorical and scale variables to the formation of clusters of cases as indicated by the grouping variable.

### Basic Specification

The basic specification is a grouping variable, a CATEGORICAL or CONTINUOUS subcommand, and a PLOT subcommand.

### Subcommand Order

- The grouping variable must be specified first.
- Subcommands can be specified in any order.

**Syntax Rules**

- All subcommands should be specified only once. If a subcommand is repeated, only the last specification will be used.

**Limitations**

The `WEIGHT` variable, if specified, is ignored by this procedure.

**Grouping Variable**

- The grouping variable must be the first specification after the procedure name.
- The grouping variable can be of any type (numeric or string).

**Example**

```
AIM clu_id  
  /CONTINUOUS age work salary.
```

- This is a typical example where `CLU_ID` is the cluster membership saved from a clustering procedure (say TwoStep Cluster) where `AGE`, `WORK`, and `SALARY` are the variables used to find the clusters.

**CATEGORICAL Subcommand**

Variables that are specified in this subcommand are treated as categorical variables, regardless of their defined measurement level.

- There is no restriction on the types of variables that can be specified on this subcommand.
- The grouping variable cannot be specified on this subcommand.

**CONTINUOUS Subcommand**

Variables that are specified in this subcommand are treated as scale variables, regardless of their defined measurement level.

- Variables specified on this subcommand must be numeric.
- The grouping variable cannot be specified on this subcommand.

**CRITERIA Subcommand**

The `CRITERIA` subcommand offers the following options in producing graphs.

**ADJUST = BONFERRONI | NONE**

*Adjust the confidence level for simultaneous confidence intervals or the tolerance level for simultaneous tests. BONFERRONI uses Bonferroni adjustments. This is the default. NONE specifies that no adjustments should be applied.*

**CI = number** *Confidence Interval.* This option controls the confidence level. Specify a value greater than 0 and less than 100. The default value is 95.

**HIDENOTSIG = NO | YES**

*Hide groups or variables that are determined to be not significant.* YES specifies that all confidence intervals and all test results should be shown. This is the default. NO specifies that only the significant confidence intervals and test results should be shown.

**SHOWREFLINE = NO | YES**

*Display reference lines that are the critical values or the tolerance levels in tests.* YES specifies that the appropriate reference lines should be shown. This is the default. NO specifies that reference lines should not be shown.

## **MISSING Subcommand**

The MISSING subcommand specifies the way to handle cases with user-missing values.

- A case is never used if it contains system-missing values in the grouping variable, categorical variable list, or the continuous variable list.
- If this subcommand is not specified, the default is EXCLUDE.

**EXCLUDE** *Exclude both user-missing and system-missing values.* This is the default.

**INCLUDE** *User-missing values are treated as valid.* Only system-missing values are not included in the analysis.

## **PLOT Subcommand**

The PLOT subcommand specifies which graphs to produce.

**CATEGORY** *Within Cluster Percentages.* This option displays a clustered bar chart for each categorical variable. The bars represent percentages of categories in each cluster. The cluster marginal count is used as the base for the percentages.

**CLUSTER (TYPE=BAR | PIE)**

*Cluster frequency charts.* Displays a bar or pie chart, depending upon the option selected, representing the frequency of each level of the grouping variable.

**ERRORBAR** *Error Bar.* This option displays an error bar by group ID for each continuous variable.

**IMPORTANCE (X=GROUP | VARIABLE Y=TEST | PVALUE)**

*Attribute Importance.* This option displays a bar chart that shows the relative importance of the attributes/variables. The specified options further control the display.

X = GROUP causes values of the grouping variable to be displayed on the  $x$  axis. A separate chart is produced for each variable. X = VARIABLE causes variable names to be displayed on the  $x$  axis. A separate chart is produced for each value of the grouping variable.

Y = TEST causes test statistics to be displayed on the  $y$  axis. Student's  $t$  statistics are displayed for scale variables, and chi-square statistics are displayed for categorical variables. Y = PVALUE causes  $p$ -value-related measures to be displayed on the  $y$  axis. Specifically,  $-\log_{10}(pvalue)$  is shown so that in both cases larger values indicate "more significant" results.

### **Example: Importance Charts by Group**

```
AIM clu_id
  /CONTINUOUS age work salary
  /CATEGORICAL minority
  /PLOT CATEGORY CLUSTER (TYPE = PIE) IMPORTANCE (X=GROUP Y=TEST).
```

- A frequency pie chart is requested.
- Student's  $t$  statistics are plotted against the group ID for each scale variable, and chi-square statistics are plotted against the group ID for each categorical variable.

### **Example: Importance Charts by Variable**

```
AIM clu_id
  /CONTINUOUS age work salary
  /CATEGORICAL minority
  /CRITERIA HIDENOTSIG=YES CI=95 ADJUST=NONE
  /PLOT CATEGORY CLUSTER (TYPE = BAR)
      IMPORTANCE (X = VARIABLE, Y = PVALUE).
```

- A frequency bar chart is requested.
- $-\log_{10}(pvalue)$  values are plotted against variables, both scale and categorical, for each level of the grouping variable.
- In addition, bars are not shown if their  $p$  values exceed 0.05.

# ALTER TYPE

```
ALTER TYPE varlist([input format = ] {output format   }) [varlist...]
                                     {AMIN [+ [n[%]]   }
                                     {AHEXMIN [+ [n[%]]}
[/PRINT {[ALTEREDTYPES*] [ALTEREDVALUES]}]
        {NONE              }
```

\* Default if subcommand omitted.

## Release History

Release 16.0

- Command introduced.

## Example

```
ALTER TYPE StringDate1 to StringDate4 (Date11).
ALTER TYPE ALL (A=AMIN).
```

## Overview

ALTER TYPE can be used to change the fundamental type (string or numeric) or format of variables, including changing the defined width of string variables.

### Options

- You can use the TO keyword to specify a list of variables or the ALL keyword to specify all variables in the active dataset.
- The optional input format specification restricts the type modification to only variables in the list that match the input format. If the input format doesn't include a width specification, all variables that match the basic format are included. An input format specification without a width specification includes all variables that match the basic format, regardless of defined width.
- AMIN or AHEXMIN can be used as the output format specification to change the defined width of a string variable to the minimum width necessary to display all observed values of that variable without truncation.
- AMIN + n or AHEXMIN + n sets the width of string variables to the minimum necessary width plus *n* bytes.
- AMIN + n% or AHEXMIN + n% sets the width of string variables to the minimum necessary width plus *n* percent of that width. The result is rounded to an integer.

**Basic Specification**

The basic specification is the name of a variable in the active dataset followed by an output format specification enclosed in parentheses, as in:

```
ALTER TYPE StringVar (A4).
```

**Syntax Rules**

- All variables specified or implied in the variable list(s) must exist in the active dataset.
- Each variable list must be followed by a format specification enclosed in parentheses.
- Format specifications must be valid IBM® SPSS® Statistics formats. For information on valid format specifications, see [Variable Types and Formats](#).
- If specified, the optional input format must be followed by an equals sign and then the output format.
- If a variable is included in more than one variable list on the same ALTER TYPE command, only the format specification associated with the last instance of the variable name will be applied. (If you want to “chain” multiple modifications for the same variable, use multiple ALTER TYPE commands.)

**Operations**

- If the command does not include any AMIN or AHEXMIN format specifications and does not include ALTEREDVALUES on the PRINT subcommand, the command takes effect immediately. It does not read the active dataset or execute pending transformations.
- If the command includes one or more AMIN or AHEXMIN format specifications or includes ALTEREDVALUES on the PRINT subcommand, the command reads the active dataset and causes execution of any pending transformations.
- Converting a numeric variable to string will result in truncated values if the numeric value cannot be represented in the specified string width.
- Converting a string variable to numeric will result in a system-missing value if the string contains characters that would be invalid for the specified numeric format.

**Examples**

```
DATA LIST FREE
  /Numvar1 (F2) Numvar2 (F1)
  StringVar1 (A20) StringVar2 (A30)
  StringDate1 (A11) StringDate2 (A10) StringDate3 (A10).
BEGIN DATA
1 23 a234 b2345 28-Oct-2007 10/28/2007 10/29/2008
END DATA.
ALTER TYPE Numvar1 (F5.2) Numvar2 (F3).
ALTER TYPE
  StringDate1 to StringDate3 (A11 = DATE11).
ALTER TYPE
  StringDate1 to StringDate3 (A10 = ADATE10).
ALTER TYPE ALL (A=AMIN).
```

- The first ALTER TYPE command changes the formats of *Numvar1* and *Numvar2* from F2 and F1 to F5.2 and F3.

- The next `ALTER TYPE` command converts all string variables between *StringDate1* and *StringDate3* (in file order) with a defined string width of 11 to the numeric date format `DATE11` (dd-mmm-yyyy). The only variable that meets these criteria is *StringDate1*; so that is the only variable converted.
- The third `ALTER TYPE` command converts all string variables between *StringDate1* and *StringDate3* with a defined string width of 10 to the numeric date format `ADATE11` (mm/dd/yyyy). In this example, this conversion is applied to *StringDate2* and *StringDate3*.
- The last `ALTER TYPE` command changes the defined width of all remaining string variables to the minimum width necessary for each variable to avoid truncation of any values. In this example, *StringVar1* changes from `A20` to `A4` and *StringVar2* changes from `A30` to `A5`. This command reads the data and executes any pending transformation commands.

## ***PRINT Subcommand***

The optional `PRINT` subcommand controls the display of information about the variables modified by the `ALTER TYPE` command. The following options are available:

**ALTEREDTYPES.** *Display a list of variables for which the formats were changed and the old and new formats.* This is the default.

**ALTEREDVALUES.** *Display a report of values that were changed if the fundamental type (string or numeric) was changed or the defined string width was changed.* This report is limited to the first 25 values that were changed for each variable.

**NONE.** *Don't display any summary information.* This is an alternative to `ALTEREDTYPES` and/or `ALTEREDVALUES` and cannot be used in combination with them.

# ALSCAL

ALSCAL is available in the Statistics Base option.

```
ALSCAL VARIABLES=varlist

[/FILE='savfile'|'dataset']
    [CONFIG [({INITIAL})] [ROWCONF [({INITIAL})]
        {FIXED } {FIXED }

        [COLCONF [({INITIAL})] [SUBJWGHT[({INITIAL})]
            {FIXED } {FIXED }

            [STIMWGHT[({INITIAL})]
                {FIXED }

[/INPUT=ROWS ({ALL**})]
    { n }

[/SHAPE={SYMMETRIC**}]
    {ASYMMETRIC }
    {RECTANGULAR}

[/LEVEL={ORDINAL** [({[UNTIE] [SIMILAR]})]
    {INTERVAL[({1})]
    { {n} }
    {RATIO[({1})]
    { {n} }
    {NOMINAL }

[/CONDITION={MATRIX** }
    {ROW }
    {UNCONDITIONAL}

[/{MODEL }={EUCLID**}]
    {METHOD} {INDSCAL }
    {ASCAL }
    {AINDS }
    {GEMSCAL }

[/CRITERIA=[NEGATIVE] [CUTOFF({0**})] [CONVERGE({.001})]
    { n } { n }

    [ITER({30})] [STRESSMIN({.005})] [NOULB]
    { n } { n }

    [DIMENS({2** })] [DIRECTIONS(n)]
    {min[,max]}

    [CONSTRAIN] [TIESTORE(n)]

[/PRINT=[DATA] [HEADER]] [/PLOT=[DEFAULT] [ALL]]

[/OUTFILE='savfile'|'dataset']

[/MATRIX=IN({'savfile'|'dataset'})]
    {* }
}
```

**\*\*Default if the subcommand or keyword is omitted.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)



**Example**

```
ALSCAL VARIABLES=ATLANTA TO TAMPA.
```

*ALSCAL was originally designed and programmed by Forrest W. Young, Yoshio Takane, and Rostyslaw J. Lewyckij of the Psychometric Laboratory, University of North Carolina.*

## Overview

ALSCAL uses an alternating least-squares algorithm to perform multidimensional scaling (MDS) and multidimensional unfolding (MDU). You can select one of the five models to obtain stimulus coordinates and/or weights in multidimensional space.

**Options**

**Data Input.** You can read inline data matrices, including all types of two- or three-way data, such as a single matrix or a matrix for each of several subjects, using the `INPUT` subcommand. You can read square (symmetrical or asymmetrical) or rectangular matrices of proximities with the `SHAPE` subcommand and proximity matrices created by `PROXIMITIES` and `CLUSTER` with the `MATRIX` subcommand. You can also read a file of coordinates and/or weights to provide initial or fixed values for the scaling process with the `FILE` subcommand.

**Methodological Assumptions.** You can specify data as matrix-conditional, row-conditional, or unconditional on the `CONDITION` subcommand. You can treat data as nonmetric (nominal or ordinal) or as metric (interval or ratio) using the `LEVEL` subcommand. You can also use `LEVEL` to identify ordinal-level proximity data as measures of similarity or dissimilarity, and you can specify tied observations as untied (continuous) or leave them tied (discrete).

**Model Selection.** You can specify the most commonly used multidimensional scaling models by selecting the correct combination of ALSCAL subcommands, keywords, and criteria. In addition to the default Euclidean distance model, the `MODEL` subcommand offers the individual differences (weighted) Euclidean distance model (`INDSCAL`), the asymmetric Euclidean distance model (`ASCAL`), the asymmetric individual differences Euclidean distance model (`AINDS`), and the generalized Euclidean metric individual differences model (`GEMSCAL`).

**Output.** You can produce output that includes raw and scaled input data, missing-value patterns, normalized data with means, squared data with additive constants, each subject's scalar product and individual weight space, plots of linear or nonlinear fit, and plots of the data transformations using the `PRINT` and `PLOT` subcommands.

**Basic Specification**

The basic specification is `VARIABLES` followed by a variable list. By default, ALSCAL produces a two-dimensional nonmetric Euclidean multidimensional scaling solution. Input is assumed to be one or more square symmetric matrices with data elements that are dissimilarities at the ordinal level of measurement. Ties are not untied, and conditionality is by subject. Values less than 0 are treated as missing. The default output includes the improvement in Young's S-stress for successive iterations, two measures of fit for each input matrix (Kruskal's stress and the squared correlation, RSQ), and the derived configurations for each of the dimensions.

**Subcommand Order**

Subcommands can be named in any order.

**Operations**

- ALSCAL calculates the number of input matrices by dividing the total number of observations in the dataset by the number of rows in each matrix. All matrices must contain the same number of rows. This number is determined by the settings on `SHAPE` and `INPUT` (if used). For square matrix data, the number of rows in the matrix equals the number of variables. For rectangular matrix data, it equals the number of rows specified or implied. For additional information, see the `INPUT` and `SHAPE` subcommands below.
- ALSCAL ignores user-missing specifications in all variables in the configuration/weights file. [For more information, see the topic FILE Subcommand on p. 183.](#) The system-missing value is converted to 0.
- With split-file data, ALSCAL reads initial or fixed configurations from the configuration/weights file for each split-file group. [For more information, see the topic FILE Subcommand on p. 183.](#) If there is only one initial configuration in the file, ALSCAL rereads these initial or fixed values for successive split-file groups.
- By default, ALSCAL estimates upper and lower bounds on missing values in the active dataset in order to compute the initial configuration. To prevent this, specify `CRITERIA=NOULB`. Missing values are always ignored during the iterative process.

**Limitations**

- A maximum of 100 variables on the `VARIABLES` subcommand.
- A maximum of six dimensions can be scaled.
- ALSCAL does not recognize data weights created by the `WEIGHT` command.
- ALSCAL analyses can include no more than 32,767 values in each of the input matrices. Large analyses may require significant computing time.

**Example**

```
* Air distances among U.S. cities.
* Data are from Johnson and Wichern (1982), page 563.
DATA LIST
  /ATLANTA BOSTON CINCNATI COLUMBUS DALLAS INDNPLIS
  LITTROCK LOSANGEL MEMPHIS STLOUIS SPOKANE TAMPA 1-60.
BEGIN DATA
  0
1068  0
  461  867  0
  549  769  107  0
  805  1819  943  1050  0
  508  941  108  172  882  0
  505  1494  618  725  325  562  0
2197  3052  2186  2245  1403  2080  1701  0
  366  1355  502  586  464  436  137  1831  0
  558  1178  338  409  645  234  353  1848  294  0
2467  2747  2067  2131  1891  1959  1988  1227  2042  1820  0
  467  1379  928  985  1077  975  912  2480  779  1016  2821  0
END DATA.
```

```
ALSCAL VARIABLES=ATLANTA TO TAMPA
/PLOT.
```

- By default, ALSCAL assumes a symmetric matrix of dissimilarities for ordinal-level variables. Only values below the diagonal are used. The upper triangle can be left blank. The 12 cities form the rows and columns of the matrix.
- The result is a classical MDS analysis that reproduces a map of the United States when the output is rotated to a north-south by east-west orientation.

## ***VARIABLES Subcommand***

VARIABLES identifies the columns in the proximity matrix or matrices that ALSCAL reads.

- VARIABLES is required and can name only numeric variables.
- Each matrix must have at least four rows and four columns.

## ***INPUT Subcommand***

ALSCAL reads data row by row, with each case in the active dataset representing a single row in the data matrix. (VARIABLES specifies the columns.) Use INPUT when reading rectangular data matrices to specify how many rows are in each matrix.

- The specification on INPUT is ROWS. If INPUT is not specified or is specified without ROWS, the default is ROWS (ALL). ALSCAL assumes that each case in the active dataset represents one row of a single input matrix and that the result is a square matrix.
- You can specify the number of rows ( $n$ ) in each matrix in parentheses after the keyword ROWS. The number of matrices equals the number of observations divided by the number specified.
- The number specified on ROWS must be at least 4 and must divide evenly into the total number of rows in the data.
- With split-file data,  $n$  refers to the number of cases in each split-file group. All split-file groups must have the same number of rows.

### ***Example***

```
ALSCAL VARIABLES=V1 to V7 /INPUT=ROWS(8).
```

- INPUT indicates that there are eight rows per matrix, with each case in the active dataset representing one row.
- The total number of cases must be divisible by 8.

## ***SHAPE Subcommand***

Use SHAPE to specify the structure of the input data matrix or matrices.

- You can specify one of the three keywords listed below.

- Both **SYMMETRIC** and **ASYMMETRIC** refer to square matrix data.

<b>SYMMETRIC</b>	<i>Symmetric data matrix or matrices.</i> For a symmetric matrix, <b>ALSCAL</b> looks only at the values below the diagonal. Values on and above the diagonal can be omitted. This is the default.
<b>ASYMMETRIC</b>	<i>Asymmetric data matrix or matrices.</i> The corresponding values in the upper and lower triangles are not all equal. The diagonal is ignored.
<b>RECTANGULAR</b>	<i>Rectangular data matrix or matrices.</i> The rows and columns represent different sets of items.

### Example

```
ALSCAL VAR=V1 TO V8 /SHAPE=RECTANGULAR.
```

- **ALSCAL** performs a classical MDU analysis, treating the rows and columns as separate sets of items.

## LEVEL Subcommand

**LEVEL** identifies the level of measurement for the values in the data matrix or matrices. You can specify one of the keywords defined below.

<b>ORDINAL</b>	<i>Ordinal-level data.</i> This specification is the default. It treats the data as ordinal, using Kruskal's least-squares monotonic transformation (Kruskal, 1964). The analysis is nonmetric. By default, the data are treated as discrete dissimilarities. Ties in the data remain tied throughout the analysis. To change the default, specify <b>UNTIE</b> and/or <b>SIMILAR</b> in parentheses. <b>UNTIE</b> treats the data as continuous and resolves ties in an optimal fashion; <b>SIMILAR</b> treats the data as similarities. <b>UNTIE</b> and <b>SIMILAR</b> cannot be used with the other levels of measurement.
<b>INTERVAL(n)</b>	<i>Interval-level data.</i> This specification produces a metric analysis of the data using classical regression techniques. You can specify any integer from 1 to 4 in parentheses for the degree of polynomial transformation to be fit to the data. The default is 1.
<b>RATIO(n)</b>	<i>Ratio-level data.</i> This specification produces a metric analysis. You can specify an integer from 1 to 4 in parentheses for the degree of polynomial transformation. The default is 1.
<b>NOMINAL</b>	<i>Nominal-level data.</i> This specification treats the data as nominal by using a least-squares categorical transformation (Takane, Young, and de Leeuw, 1977). This option produces a nonmetric analysis of nominal data. It is useful when there are few observed categories, when there are many observations in each category, and when the order of the categories is not known.

### Example

```
ALSCAL VAR=ATLANTA TO TAMPA /LEVEL=INTERVAL(2).
```

- This example identifies the distances between U.S. cities as interval-level data. The 2 in parentheses indicates a polynomial transformation with linear and quadratic terms.

## CONDITION Subcommand

CONDITION specifies which numbers in a dataset are comparable.

<b>MATRIX</b>	<i>Only numbers within each matrix are comparable.</i> If each matrix represents a different subject, this specification makes comparisons conditional by subject. This is the default.
<b>ROW</b>	<i>Only numbers within the same row are comparable.</i> This specification is appropriate only for asymmetric or rectangular data. They cannot be used when ASCAL or AINDS is specified on MODEL.
<b>UNCONDITIONAL</b>	<i>All numbers are comparable.</i> Comparisons can be made among any values in the input matrix or matrices.

### Example

```
ALSCAL VAR=V1 TO V8 /SHAPE=RECTANGULAR /CONDITION=ROW.
```

- ALSCAL performs a Euclidean MDU analysis conditional on comparisons within rows.

## FILE Subcommand

ALSCAL can read proximity data from the active dataset or, with the MATRIX subcommand, from a matrix data file created by PROXIMITIES or CLUSTER. The FILE subcommand reads a file containing additional data—an initial or fixed configuration for the coordinates of the stimuli and/or weights for the matrices being scaled. This file can be created with the OUTFILE subcommand on ALSCAL or with an input program (created with the INPUT PROGRAM command).

- The minimum specification is the file that contains the configurations and/or weights.
- FILE can include additional specifications that define the structure of the configuration/weights file.
- The variables in the configuration/weights file that correspond to successive ALSCAL dimensions must have the names *DIM1*, *DIM2*, ..., *DIMr*, where *r* is the maximum number of ALSCAL dimensions. The file must also contain the short string variable *TYPE\_* to identify the types of values in all rows.
- Values for the variable *TYPE\_* can be CONFIG, ROWCONF, COLCONF, SUBJWGHT, and STIMWGHT, in that order. Each value can be truncated to the first three letters. Stimulus coordinate values are specified as CONFIG; row stimulus coordinates, as ROWCONF; column stimulus coordinates, as COLCONF; and subject and stimulus weights, as SUBJWGHT and STIMWGHT, respectively. ALSCAL accepts CONFIG and ROWCONF interchangeably.
- ALSCAL skips unneeded types as long as they appear in the file in their proper order. Generalized weights (GEM) and flattened subject weights (FLA) cannot be initialized or fixed and will always be skipped. (These weights can be generated by ALSCAL but cannot be used as input.)

The following list summarizes the optional specifications that can be used on FILE to define the structure of the configuration/weights file:

- Each specification can be further identified with the option INITIAL or FIXED in parentheses.

- `INITIAL` is the default. `INITIAL` indicates that the external configuration or weights are to be used as initial coordinates and are to be modified during each iteration.
- `FIXED` forces `ALSCAL` to use the externally defined structure without modification to calculate the best values for all unfixed portions of the structure.

<b>CONFIG</b>	<i>Read stimulus configuration.</i> The configuration/weights file contains initial stimulus coordinates. Input of this type is appropriate when <code>SHAPE=SYMMETRIC</code> or <code>SHAPE=ASYMMETRIC</code> , or when the number of variables in a matrix equals the number of variables on the <code>ALSCAL</code> command. The value of the <code>TYPE_</code> variable must be either <code>CON</code> or <code>ROW</code> for all stimulus coordinates for the configuration.
<b>ROWCONF</b>	<i>Read row stimulus configuration.</i> The configuration/weights file contains initial row stimulus coordinates. This specification is appropriate if <code>SHAPE=RECTANGULAR</code> and if the number of <code>ROWCONF</code> rows in the matrix equals the number of rows specified on the <code>INPUT</code> subcommand (or, if <code>INPUT</code> is omitted, the number of cases in the active dataset). The value of <code>TYPE_</code> must be either <code>ROW</code> or <code>CON</code> for the set of coordinates for each row.
<b>COLCONF</b>	<i>Read column stimulus configuration.</i> The configuration/weights file contains initial column stimulus coordinates. This kind of file can be used only if <code>SHAPE=RECTANGULAR</code> and if the number of <code>COLCONF</code> rows in the matrix equals the number of variables on the <code>ALSCAL</code> command. The value of <code>TYPE_</code> must be <code>COL</code> for the set of coordinates for each column.
<b>SUBJWGHT</b>	<i>Read subject (matrix) weights.</i> The configuration/weights file contains subject weights. The number of observations in a subject-weights matrix must equal the number of matrices in the proximity file. Subject weights can be used only if the model is <code>INDSCAL</code> , <code>AINDS</code> , or <code>GEMSCAL</code> . The value of <code>TYPE_</code> for each set of weights must be <code>SUB</code> .
<b>STIMWGHT</b>	<i>Read stimulus weights.</i> The configuration/weights file contains stimulus weights. The number of observations in the configuration/weights file must equal the number of matrices in the proximity file. Stimulus weights can be used only if the model is <code>AINDS</code> or <code>ASCAL</code> . The value of <code>TYPE_</code> for each set of weights must be <code>STI</code> .

If the optional specifications for the configuration/weights file are not specified on `FILE`, `ALSCAL` sequentially reads the `TYPE_` values appropriate to the model and shape according to the defaults in the table below.

### Example

```
ALSCAL VAR=V1 TO V8 /FILE=ONE CON(FIXED) STI(INITIAL) .
```

- `ALSCAL` reads the configuration/weights file `ONE`.
- The stimulus coordinates are read as fixed values, and the stimulus weights are read as initial values.

Table 12-1  
Default specifications for the `FILE` subcommand

Shape	Model	Default specifications
SYMMETRIC	EUCLID	CONFIG (or ROWCONF)
	INDSCAL	CONFIG (or ROWCONF) SUBJWGHT
	GEMSCAL	CONFIG (or ROWCONF) SUBJWGHT
ASYMMETRIC	EUCLID	CONFIG (or ROWCONF)

Shape	Model	Default specifications
	INDSCAL	CONFIG (or ROWCONF) SUBJWGHT
	GEMSCAL	CONFIG (or ROWCONF) SUBJWGHT
	ASCAL	CONFIG (or ROWCONF) STIMWGHT
	AINDS	CONFIG (or ROWCONF) SUBJWGHT STIMWGHT
RECTANGULAR	EUCLID	ROWCONF (or CONFIG) COLCONF
	INDSCAL	ROWCONF (or CONFIG) COLCONF SUBJWGHT
	GEMSCAL	ROWCONF (or CONFIG) COLCONF SUBJWGHT

## MODEL Subcommand

MODEL (alias METHOD) defines the scaling model for the analysis. The only specification is MODEL (or METHOD) and any one of the five scaling and unfolding model types. EUCLID is the default.

- EUCLID**      *Euclidean distance model.* This model can be used with any type of proximity matrix and is the default.
- INDSCAL**      *Individual differences (weighted) Euclidean distance model.* ALSCAL scales the data using the weighted individual differences Euclidean distance model (Carroll and Chang, 1970). This type of analysis can be specified only if the analysis involves more than one data matrix and more than one dimension is specified on CRITERIA.
- ASCAL**      *Asymmetric Euclidean distance model.* This model (Young, 1975) can be used only if SHAPE=ASYMMETRIC and more than one dimension is requested on CRITERIA.
- AINDS**      *Asymmetric individual differences Euclidean distance model.* This option combines Young's asymmetric Euclidean model (Young et al., 1975) with the individual differences model (Carroll et al., 1970). This model can be used only when SHAPE=ASYMMETRIC, the analysis involves more than one data matrix, and more than one dimension is specified on CRITERIA.
- GEMSCAL**      *Generalized Euclidean metric individual differences model.* The number of directions for this model is set with the DIRECTIONS option on CRITERIA. The number of directions specified can be equal to but cannot exceed the group space dimensionality. By default, the number of directions equals the number of dimensions in the solution.

### Example

```
ALSCAL VARIABLES = V1 TO V6
/SHAPE = ASYMMETRIC
/CONDITION = ROW
/MODEL = GEMSCAL
/CRITERIA = DIM(4) DIRECTIONS(4) .
```

- In this example, the number of directions in the GEMSCAL model is set to 4.

## **CRITERIA Subcommand**

Use **CRITERIA** to control features of the scaling model and to set convergence criteria for the solution. You can specify one or more of the following:

<b>CONVERGE(n)</b>	<i>Stop iterations if the change in S-stress is less than n.</i> S-stress is a goodness-of-fit index. By default, $n=0.001$ . To increase the precision of a solution, specify a smaller value, for example, 0.0001. To obtain a less precise solution (perhaps to reduce computing time), specify a larger value, for example, 0.05. Negative values are not allowed. If $n=0$ , the algorithm will iterate 30 times unless a value is specified with the <b>ITER</b> option.
<b>ITER(n)</b>	<i>Set the maximum number of iterations to n.</i> The default value is 30. A higher value will give a more precise solution but will take longer to compute.
<b>STRESSMIN(n)</b>	<i>Set the minimum stress value to n.</i> By default, <b>ALSCAL</b> stops iterating when the value of S-stress is 0.005 or less. <b>STRESSMIN</b> can be assigned any value from 0 to 1.
<b>NEGATIVE</b>	<i>Allow negative weights in individual differences models.</i> By default, <b>ALSCAL</b> does not permit the weights to be negative. Weighted models include <b>INDSCAL</b> , <b>ASCAL</b> , <b>AINDS</b> , and <b>GEMSCAL</b> . The <b>NEGATIVE</b> option is ignored if the model is <b>EUCLID</b> .
<b>CUTOFF(n)</b>	<i>Set the cutoff value for treating distances as missing to n.</i> By default, <b>ALSCAL</b> treats all negative similarities (or dissimilarities) as missing and 0 and positive similarities as nonmissing ( $n=0$ ). Changing the <b>CUTOFF</b> value causes <b>ALSCAL</b> to treat similarities greater than or equal to that value as nonmissing. User- and system-missing values are considered missing regardless of the <b>CUTOFF</b> specification.
<b>NOULB</b>	<i>Do not estimate upper and lower bounds on missing values.</i> By default, <b>ALSCAL</b> estimates the upper and lower bounds on missing values in order to compute the initial configuration. This specification has no effect during the iterative process, when missing values are ignored.
<b>DIMENS(min[,max])</b>	<i>Set the minimum and maximum number of dimensions in the scaling solution.</i> By default, <b>ALSCAL</b> calculates a solution with two dimensions. To obtain solutions for more than two dimensions, specify the minimum and the maximum number of dimensions in parentheses after <b>DIMENS</b> . The minimum and maximum can be integers between 2 and 6. A single value represents both the minimum and the maximum. For example, <b>DIMENS(3)</b> is equivalent to <b>DIMENS(3,3)</b> . The minimum number of dimensions can be set to 1 only if <b>MODEL=EUCLID</b> .
<b>DIRECTIONS(n)</b>	<i>Set the number of principal directions in the generalized Euclidean model to n.</i> This option has no effect for models other than <b>GEMSCAL</b> . The number of principal directions can be any positive integer between 1 and the number of dimensions specified on the <b>DIMENS</b> option. By default, the number of directions equals the number of dimensions.
<b>TIESTORE(n)</b>	<i>Set the amount of storage needed for ties to n.</i> This option estimates the amount of storage needed to deal with ties in ordinal data. By default, the amount of storage is set to 1000 or the number of cells in a matrix, whichever is smaller. Should this be insufficient, <b>ALSCAL</b> terminates and displays a message that more space is needed.
<b>CONSTRAIN</b>	<i>Constrain multidimensional unfolding solution.</i> This option can be used to keep the initial constraints throughout the analysis.



## ***PRINT Subcommand***

PRINT requests output not available by default. You can specify the following:

<b>DATA</b>	<i>Display input data.</i> The display includes both the initial data and the scaled data for each subject according to the structure specified on SHAPE.
<b>HEADER</b>	<i>Display a header page.</i> The header includes the model, output, algorithmic, and data options in effect for the analysis.

- Data options listed by PRINT=HEADER include the number of rows and columns, number of matrices, measurement level, shape of the data matrix, type of data (similarity or dissimilarity), whether ties are tied or untied, conditionality, and data cutoff value.
- Model options listed by PRINT=HEADER are the type of model specified (EUCLID, INDSCAL, ASCAL, AINDS, or GEMSCAL), minimum and maximum dimensionality, and whether or not negative weights are permitted.
- Output options listed by PRINT=HEADER indicate whether the output includes the header page and input data, whether ALSCAL plotted configurations and transformations, whether an output dataset was created, and whether initial stimulus coordinates, initial column stimulus coordinates, initial subject weights, and initial stimulus weights were computed.
- Algorithmic options listed by PRINT=HEADER include the maximum number of iterations permitted, the convergence criterion, the maximum S-stress value, whether or not missing data are estimated by upper and lower bounds, and the amount of storage allotted for ties in ordinal data.

### ***Example***

```
ALSCAL VAR=ATLANTA TO TAMPA /PRINT=DATA.
```

- In addition to scaled data, ALSCAL will display initial data.

## ***PLOT Subcommand***

PLOT controls the display of plots. The minimum specification is simply PLOT to produce the defaults.

<b>DEFAULT</b>	<i>Default plots.</i> Default plots include plots of stimulus coordinates, matrix weights (if the model is INDSCAL, AINDS, or GEMSCAL), and stimulus weights (if the model is AINDS or ASCAL). The default also includes a scatterplot of the linear fit between the data and the model and, for certain types of data, scatterplots of the nonlinear fit and the data transformation.
<b>ALL</b>	<i>Transformation plots in addition to the default plots.</i> A separate plot is produced for each subject if CONDITION=MATRIX and a separate plot for each row if CONDITION=ROW. For interval and ratio data, PLOT=ALL has the same effect as PLOT=DEFAULT. This option can generate voluminous output, particularly when CONDITION=ROW.

**Example**

```
ALSCAL VAR=V1 TO V8 /INPUT=ROWS(8) /PLOT=ALL.
```

- This command produces all of the default plots. It also produces a separate plot for each subject's data transformation and a plot of *V1* through *V8* in a two-dimensional space for each subject.

**OUTFILE Subcommand**

OUTFILE saves coordinate and weight matrices to a data file in IBM® SPSS® Statistics format. The only specification is a name for the output file.

- The output data file has an alphanumeric (short string) variable named *TYPE\_* that identifies the kind of values in each row, a numeric variable named *DIMENS* that specifies the number of dimensions, a numeric variable named *MATNUM* that indicates the subject (matrix) to which each set of coordinates corresponds, and variables named *DIM1*, *DIM2*, ..., *DIMn* that correspond to the *n* dimensions in the model.
- The values of any split-file variables are also included in the output file.
- The file created by OUTFILE can be used by subsequent ALSICAL commands as initial data.

The following are the types of configurations and weights that can be included in the output file:

<b>CONFIG</b>	<i>Stimulus configuration coordinates.</i>
<b>ROWCONF</b>	<i>Row stimulus configuration coordinates.</i>
<b>COLCONF</b>	<i>Column stimulus configuration coordinates.</i>
<b>SUBJWGHT</b>	<i>Subject (matrix) weights.</i>
<b>FLATWGHT</b>	<i>Flattened subject (matrix) weights.</i>
<b>GEMWGHT</b>	<i>Generalized weights.</i>
<b>STIMWGHT</b>	<i>Stimulus weights.</i>

Only the first three characters of each identifier are written to the variable *TYPE\_* in the file. For example, CONFIG becomes CON. The structure of the file is determined by the SHAPE and MODEL subcommands, as shown in the following table.

Table 12-2  
Types of configurations and/or weights in output files

Shape	Model	TYPE_
SYMMETRIC	EUCLID	CON
	INDSCAL	CON SUB FLA
	GEMSCAL	CON SUB FLA GEM
ASYMMETRIC	EUCLID	CON
	INDSCAL	CON SUB FLA

Shape	Model	TYPE_
	GEMSCAL	CON SUB FLA GEM
	ASCAL	CON STI
	AINDS	CON SUB FLA STI
RECTANGULAR	EUCLID	ROW COL
	INDSCAL	ROW COL SUB FLA
	GEMSCAL	ROW COL SUB FLA GEM

### Example

ALSCAL VAR=ATLANTA TO TAMPA /OUTFILE=ONE.

- `OUTFILE` creates the configuration/weights file *ONE* from the example of air distances between cities.

## MATRIX Subcommand

`MATRIX` reads matrix data files. It can read a matrix written by either `PROXIMITIES` or `CLUSTER`.

- Generally, data read by `ALSCAL` are already in matrix form. If the matrix materials are in the active dataset, you do not need to use `MATRIX` to read them. Simply use the `VARIABLES` subcommand to indicate the variables (or columns) to be used. However, if the matrix materials are not in the active dataset, `MATRIX` must be used to specify the matrix data file that contains the matrix.
- The proximity matrices that `ALSCAL` reads have `ROWTYPE_` values of `PROX`. No additional statistics should be included with these matrix materials.
- `ALSCAL` ignores unrecognized `ROWTYPE_` values in the matrix file. In addition, it ignores variables present in the matrix file that are not specified on the `VARIABLES` subcommand in `ALSCAL`. The order of rows and columns in the matrix is unimportant.
- Since `ALSCAL` does not support case labeling, it ignores values for the `ID` variable (if present) in a `CLUSTER` or `PROXIMITIES` matrix.
- If split-file processing was in effect when the matrix was written, the same split file must be in effect when `ALSCAL` reads that matrix.
- The specification on `MATRIX` is the keyword `IN` and the matrix file in parentheses.

- `MATRIX=IN` cannot be used unless a active dataset has already been defined. To read an existing matrix data file at the beginning of a session, first use `GET` to retrieve the matrix file and then specify `IN(*)` on `MATRIX`.

**IN (filename)**      *Read a matrix data file.* If the matrix data file is the active dataset, specify an asterisk in parentheses (\*). If the matrix data file is another file, specify the filename in parentheses. A matrix file read from an external file does not replace the active dataset.

### Example

```
PROXIMITIES V1 TO V8 /ID=NAMEVAR /MATRIX=OUT(*) .
ALSCAL VAR=CASE1 TO CASE10 /MATRIX=IN(*) .
```

- `PROXIMITIES` uses `V1` through `V8` in the active dataset to generate a matrix file of Euclidean distances between each pair of cases based on the eight variables. The number of rows and columns in the resulting matrix equals the number of cases. `MATRIX=OUT` then replaces the active dataset with this new matrix data file.
- `MATRIX=IN` on `ALSCAL` reads the matrix data file, which is the new active dataset. In this instance, `MATRIX` is optional because the matrix materials are in the active dataset.
- If there were 10 cases in the original active dataset, `ALSCAL` performs a multidimensional scaling analysis in two dimensions on `CASE1` through `CASE10`.

### Example

```
GET FILE PROXMTX .
ALSCAL VAR=CASE1 TO CASE10 /MATRIX=IN(*) .
```

- `GET` retrieves the matrix data file `PROXMTX`.
- `MATRIX=IN` specifies an asterisk because the active dataset is the matrix. `MATRIX` is optional, however, since the matrix materials are in the active dataset.

### Example

```
GET FILE PRSNNL .
FREQUENCIES VARIABLE=AGE .
ALSCAL VAR=CASE1 TO CASE10 /MATRIX=IN(PROXMTX) .
```

- This example performs a frequencies analysis on the file `PRSNNL` and then uses a different file containing matrix data for `ALSCAL`. The file is an existing matrix data file.
- `MATRIX=IN` is required because the matrix data file, `PROXMTX`, is not the active dataset. `PROXMTX` does not replace `PRSNNL` as the active dataset.

## Specification of Analyses

The following tables summarize the analyses that can be performed for the major types of proximity matrices that you can use with `ALSCAL`, list the specifications needed to produce these analyses for nonmetric models, and list the specifications for metric models. You can include additional specifications to control the precision of your analysis with `CRITERIA`.

Table 12-3  
Models for types of matrix input

Matrix mode	Matrix form	Model class	Single matrix	Replications of single matrix	Two or more individual matrices
Object by object	Symmetric	Multidimensional scaling	CMDS Classical multidimensional scaling	RMDS Replicated multidimensional scaling	WMDS(INDSCAL) Weighted multidimensional scaling
	Asymmetric single process	Multidimensional scaling	CMDS(row conditional) Classical row conditional multidimensional scaling	RMDS(row conditional) Replicated row conditional multi dimensional scaling	WMDS(row conditional) Weighted row conditional multidimensional scaling
	Asymmetric multiple process	Internal asymmetric multidimensional scaling	CAMDS Classical asymmetric multidimensional scaling	RAMDS Replicated asymmetric multidimensional scaling	WAMDS Weighted asymmetric multidimensional scaling
		External asymmetric multidimensional scaling	CAMDS(external) Classical external asymmetric multidimensional scaling	RAMDS(external) Replicated external asymmetric multidimensional scaling	WAMDS(external) Weighted external asymmetric multidimensional scaling
Object by attribute	Rectangular	Internal unfolding	CMDU Classical internal multidimensional unfolding	RMDU Replicated internal multidimensional unfolding	WMDU Weighted internal multidimensional unfolding
		External unfolding	CMDU(external) Classical external multidimensional unfolding	RMDU(external) Replicated external multidimensional unfolding	WMDU(external) Weighted external multidimensional unfolding

Table 12-4  
ALSCAL specifications for nonmetric models

Matrix mode	Matrix form	Model class	Single matrix	Replications of single matrix	Two or more individual matrices	
Object by object	Symmetric	Multidimensional scaling	ALSCAL VAR= varlist.	ALSCAL VAR= varlist.	ALSCAL VAR= varlist /MODEL=INDSCAL.	
	Asymmetric single process	Multidimensional scaling	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /CONDITION=ROW.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /CONDITION=ROW.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /CONDITION=ROW /MODEL=INDSCAL.	
	Asymmetric multiple process	Internal asymmetric multidimensional scaling	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /MODEL=ASCAL.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /MODEL=ASCAL.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /MODEL=ASCAL.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /MODEL=AINDS.
		External asymmetric multidimensional scaling	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /MODEL=ASCAL /FILE=file COLCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /MODEL=ASCAL /FILE=file COLCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /MODEL=ASCAL /FILE=file COLCONF (FIX) .	

Matrix mode	Matrix form	Model class	Single matrix	Replications of single matrix	Two or more individual matrices
Object by attribute	Rectangular	Internal unfolding	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW.	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION(ROW) .	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /MODEL=INDSCAL.
		External unfolding	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /FILE=file ROWCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /FILE=file ROWCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /FILE=file ROWCONF (FIX) /MODEL=INDSCAL.

Table 12-5  
ALSCAL specifications for metric models

Matrix mode	Matrix form	Model class	Single matrix	Replications of single matrix	Two or more individual matrices
Object by object	Symmetric	Multidimensional scaling	ALSCAL VAR= varlist /LEVEL=INT.	ALSCAL VAR= varlist /LEVEL=INT.	ALSCAL VAR= varlist /LEVEL=INT /MODEL=INDSCAL.
	Asymmetric single process	Multidimensional scaling	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /CONDITION=ROW /LEVEL=INT.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /CONDITION=ROW /LEVEL=INT.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /CONDITION=ROW /LEVEL=INT /MODEL=INDSCAL.
	Asymmetric multiple process	Internal asymmetric multidimensional scaling	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /LEVEL=INT /MODEL=ASCAL.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /LEVEL=INT /MODEL=ASCAL.	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /LEVEL=INT /MODEL=AINDS.
		External asymmetric multidimensional scaling	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /LEVEL=INT /MODEL=ASCAL /FILE=file COLCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /LEVEL=INT /MODEL=ASCAL /FILE=file COLCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=ASYMMETRIC /LEVEL=INT /MODEL=AINDS /FILE=file COLCONF (FIX) .
Object by attribute	Rectangular	Internal unfolding	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /LEVEL=INT.	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /LEVEL=INT.	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /LEVEL=INT /MODEL=INDSCAL.
		External unfolding	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /LEVEL=INT /FILE=file ROWCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /LEVEL=INT /FILE=file ROWCONF (FIX) .	ALSCAL VAR= varlist /SHAPE=REC /INP=ROWS /CONDITION=ROW /LEVEL=INT /FILE=file ROWCONF (FIX) /MODEL=INDSCAL.

## References

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- Johnson, R., and D. W. Wichern. 1982. *Applied multivariate statistical analysis*. Englewood Cliffs, N.J.: Prentice-Hall.
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# ANACOR

ANACOR is available in the Categories option.

```
ANACOR TABLE={row var (min, max) BY column var (min, max)}
           {ALL (# of rows, # of columns)           }

[/DIMENSION={2**  }]
           {value}

[/NORMALIZATION={CANONICAL**}]
           {PRINCIPAL  }
           {RPRINCIPAL }
           {CPRINCIPAL }
           {value      }

[/VARIANCES=[SINGULAR] [ROWS] [COLUMNS]]

[/PRINT={TABLE**} [PROFILES] [SCORES**] [CONTRIBUTIONS**]
        [DEFAULT] [PERMUTATION] [NONE]]

[/PLOT=[NDIM=({1, 2**  })]
        {value, value}
        {ALL, MAX    }
        [ROWS**[(n)]] [COLUMNS**[(n)]] [DEFAULT[(n)]]
        [TRROWS] [TRCOLUMNS] [JOINT[(n)]] [NONE]]

[/MATRIX OUT=[SCORE({*          })] [VARIANCE({*          })]]
               {'savfile'|'dataset'}          {'savfile'|'dataset'}
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
ANACOR TABLE=MENTAL(1,4) BY SES(1,6).
```

## Overview

ANACOR performs correspondence analysis, which is an isotropic graphical representation of the relationships between the rows and columns of a two-way table.

### Options

**Number of Dimensions.** You can specify how many dimensions ANACOR should compute.

**Method of Normalization.** You can specify one of five different methods for normalizing the row and column scores.

**Computation of Variances and Correlations.** You can request computation of variances and correlations for singular values, row scores, or column scores.

**Data Input.** You can analyze the usual individual casewise data or aggregated data from table cells.



**Display Output.** You can control which statistics are displayed and plotted. You can also control how many value-label characters are used on the plots.

**Writing Matrices.** You can write matrix data files containing row and column scores and variances for use in further analyses.

### **Basic Specification**

- The basic specification is ANACOR and the TABLE subcommand. By default, ANACOR computes a two-dimensional solution, displays the TABLE, SCORES, and CONTRIBUTIONS statistics, and plots the row scores and column scores of the first two dimensions.

### **Subcommand Order**

- Subcommands can appear in any order.

### **Operations**

- If a subcommand is specified more than once, only the last occurrence is executed.

### **Limitations**

- If the data within table cells contains negative values. ANACOR treats those values as 0.

## **Example**

```
ANACOR TABLE=MENTAL(1,4) BY SES(1,6)
/PRINT=SCORES CONTRIBUTIONS
/PLOT=ROWS COLUMNS.
```

- Two variables, *MENTAL* and *SES*, are specified on the TABLE subcommand. *MENTAL* has values ranging from 1 to 4, and *SES* has values ranging from 1 to 6.
- The row and column scores and the contribution of each row and column to the inertia of each dimension are displayed.
- Two plots are produced. The first one plots the first two dimensions of row scores, and the second one plots the first two dimensions of column scores.

## **TABLE Subcommand**

TABLE specifies the row and column variables, along with their value ranges for individual casewise data. For table data, TABLE specifies the keyword ALL and the number of rows and columns.

- The TABLE subcommand is required.

### **Casewise Data**

- Each variable is followed by a value range in parentheses. The value range consists of the variable's minimum value, a comma, and the variable's maximum value.
- Values outside of the specified range are not included in the analysis.

- Values do not have to be sequential. Empty categories receive scores of 0 and do not affect the rest of the computations.

### **Example**

```
DATA LIST FREE/VAR1 VAR2.
BEGIN DATA
3 1
6 1
3 1
4 2
4 2
6 3
6 3
6 3
3 2
4 2
6 3
END DATA.
ANACOR TABLE=VAR1(3,6) BY VAR2(1,3).
```

- `DATA LIST` defines two variables, *VAR1* and *VAR2*.
- *VAR1* has three levels, coded 3, 4, and 6, while *VAR2* also has three levels, coded 1, 2, and 3.
- Because a range of (3,6) is specified for *VAR1*, ANACOR defines four categories, coded 3, 4, 5, and 6. The empty category, 5, for which there is no data, receives zeros for all statistics but does not affect the analysis.

### **Table Data**

- The cells of a table can be read and analyzed directly by using the keyword `ALL` after `TABLE`.
- The columns of the input table must be specified as variables on the `DATA LIST` command. Only columns are defined, not rows.
- `ALL` is followed by the number of rows in the table, a comma, and the number of columns in the table, all enclosed in parentheses.
- If you want to analyze only a subset of the table, the specified number of rows and columns can be smaller than the actual number of rows and columns.
- The variables (columns of the table) are treated as the column categories, and the cases (rows of the table) are treated as the row categories.
- Rows cannot be labeled when you specify `TABLE=ALL`. If labels in your output are important, use the `WEIGHT` command method to enter your data (see [Analyzing Aggregated Data on p. 200](#)).

### **Example**

```
DATA LIST /COL01 TO COL07 1-21.
BEGIN DATA
50 19 26 8 18 6 2
16 40 34 18 31 8 3
12 35 65 66123 23 21
11 20 58110223 64 32
14 36114185714258189
0 6 19 40179143 71
END DATA.
```

ANACOR TABLE=ALL(6,7).

- DATA LIST defines the seven columns of the table as the variables.
- The TABLE=ALL specification indicates that the data are the cells of a table. The (6,7) specification indicates that there are six rows and seven columns.

## ***DIMENSION Subcommand***

DIMENSION specifies the number of dimensions you want ANACOR to compute.

- If you do not specify the DIMENSION subcommand, ANACOR computes two dimensions.
- DIMENSION is followed by an integer indicating the number of dimensions.
- In general, you should choose as few dimensions as needed to explain most of the variation. The minimum number of dimensions that can be specified is 1. The maximum number of dimensions that can be specified is equal to the number of levels of the variable with the least number of levels, minus 1. For example, in a table where one variable has five levels and the other has four levels, the maximum number of dimensions that can be specified is  $(4 - 1)$ , or 3. Empty categories (categories with no data, all zeros, or all missing data) are not counted toward the number of levels of a variable.
- If more than the maximum allowed number of dimensions is specified, ANACOR reduces the number of dimensions to the maximum.

## ***NORMALIZATION Subcommand***

The NORMALIZATION subcommand specifies one of five methods for normalizing the row and column scores. Only the scores and variances are affected; contributions and profiles are not changed.

The following keywords are available:

<b>CANONICAL</b>	<i>For each dimension, rows are the weighted average of columns divided by the matching singular value, and columns are the weighted average of rows divided by the matching singular value. This is the default if the NORMALIZATION subcommand is not specified. DEFAULT is an alias for CANONICAL. Use this normalization method if you are primarily interested in differences or similarities between variables.</i>
<b>PRINCIPAL</b>	<i>Distances between row points and column points are approximations of chi-square distances. The distances represent the distance between the row or column and its corresponding average row or column profile. Use this normalization method if you want to examine both differences between categories of the row variable and differences between categories of the column variable (but not differences between variables).</i>
<b>RPRINCIPAL</b>	<i>Distances between row points are approximations of chi-square distances. This method maximizes distances between row points. This is useful when you are primarily interested in differences or similarities between categories of the row variable.</i>
<b>CPRINCIPAL</b>	<i>Distances between column points are approximations of chi-square distances. This method maximizes distances between column points. This is useful when you are primarily interested in differences or similarities between categories of the column variable.</i>

The fifth method has no keyword. Instead, any value in the range  $-2$  to  $+2$  is specified after `NORMALIZATION`. A value of  $1$  is equal to the `RPRINCIPAL` method, a value of  $0$  is equal to `CANONICAL`, and a value of  $-1$  is equal to the `CPRINCIPAL` method. The inertia is spread over both row and column scores. This method is useful for interpreting joint plots.

## **VARIANCES Subcommand**

Use `VARIANCES` to display variances and correlations for the singular values, the row scores, and/or the column scores. If `VARIANCES` is not specified, variances and correlations are not included in the output.

The following keywords are available:

<b>SINGULAR</b>	<i>Variances and correlations of the singular values.</i>
<b>ROWS</b>	<i>Variances and correlations of the row scores.</i>
<b>COLUMNS</b>	<i>Variances and correlations of the column scores.</i>

## **PRINT Subcommand**

Use `PRINT` to control which correspondence statistics are displayed. If `PRINT` is not specified, displayed statistics include the numbers of rows and columns, all nontrivial singular values, proportions of inertia, and the cumulative proportion of inertia that is accounted for.

The following keywords are available:

<b>TABLE</b>	<i>A crosstabulation of the input variables showing row and column marginals.</i>
<b>PROFILES</b>	<i>The row and column profiles. <code>PRINT=PROFILES</code> is analogous to the <code>CELLS=ROW COLUMN</code> subcommand in <code>CROSSTABS</code>.</i>
<b>SCORES</b>	<i>The marginal proportions and scores of each row and column.</i>
<b>CONTRIBUTIONS</b>	<i>The contribution of each row and column to the inertia of each dimension, and the proportion of distance to the origin that is accounted for in each dimension.</i>
<b>PERMUTATION</b>	<i>The original table permuted according to the scores of the rows and columns for each dimension.</i>
<b>NONE</b>	<i>No output other than the singular values.</i>
<b>DEFAULT</b>	<i><code>TABLE</code>, <code>SCORES</code>, and <code>CONTRIBUTIONS</code>. These statistics are displayed if you omit the <code>PRINT</code> subcommand.</i>

## **PLOT Subcommand**

Use `PLOT` to produce plots of the row scores, column scores, and row and column scores, as well as to produce plots of transformations of the row scores and transformations of the column scores. If `PLOT` is not specified, plots are produced for the row scores in the first two dimensions and the column scores in the first two dimensions.

The following keywords are available:

<b>TRROWS</b>	<i>Plot of transformations of the row category values into row scores.</i>
<b>TRCOLUMNS</b>	<i>Plot of transformations of the column category values into column scores.</i>
<b>ROWS</b>	<i>Plot of row scores.</i>
<b>COLUMNS</b>	<i>Plot of column scores.</i>
<b>JOINT</b>	<i>A combined plot of the row and column scores. This plot is not available when NORMALIZATION=PRINCIPAL.</i>
<b>NONE</b>	<i>No plots.</i>
<b>DEFAULT</b>	<i>ROWS and COLUMNS.</i>

- The keywords ROWS, COLUMNS, JOINT, and DEFAULT can be followed by an integer value in parentheses to indicate how many characters of the value label are to be used on the plot. The value can range from 1 to 20; the default is 3. Spaces between words count as characters.
- TRROWS and TRCOLUMNS plots use the full value labels up to 20 characters.
- If a label is missing for any value, the actual values are used for all values of that variable.
- Value labels should be unique.
- The first letter of a label on a plot marks the place of the actual coordinate. Be careful that multiple-word labels are not interpreted as multiple points on a plot.

In addition to the plot keywords, the following keyword can be specified:

**NDIM**      *Dimension pairs to be plotted.* NDIM is followed by a pair of values in parentheses. If NDIM is not specified, plots are produced for dimension 1 by dimension 2.

- The first value indicates the dimension that is plotted against all higher dimensions. This value can be any integer from 1 to the number of dimensions minus 1.
- The second value indicates the highest dimension to be used in plotting the dimension pairs. This value can be any integer from 2 to the number of dimensions.
- Keyword ALL can be used instead of the first value to indicate that all dimensions are paired with higher dimensions.
- Keyword MAX can be used instead of the second value to indicate that plots should be produced up to, and including, the highest dimension fit by the procedure.

### **Example**

```
ANACOR TABLE=MENTAL(1,4) BY SES(1,6)
/PLOT NDIM(1,3) JOINT(5).
```

- The NDIM(1,3) specification indicates that plots should be produced for two dimension pairs—dimension 1 versus dimension 2 and dimension 1 versus dimension 3.
- JOINT requests combined plots of row and column scores. The (5) specification indicates that the first five characters of the value labels are to be used on the plots.

### **Example**

```
ANACOR TABLE=MENTAL(1,4) BY SES(1,6)
```

```
/PLOT NDIM(ALL,3) JOINT(5).
```

- This plot is the same as above except for the ALL specification following NDIM, which indicates that all possible pairs up to the second value should be plotted. Therefore, JOINT plots will be produced for dimension 1 versus dimension 2, dimension 2 versus dimension 3, and dimension 1 versus dimension 3.

## MATRIX Subcommand

Use MATRIX to write row and column scores and variances to matrix data files.

MATRIX is followed by keyword OUT, an equals sign, and one or both of the following keywords:

**SCORE** ('file'|'dataset') *Write row and column scores to a matrix data file.*

**VARIANCE** ('file'|'dataset') *Write variances to a matrix data file.*

- You can specify the file with either an asterisk (\*), to replace the active dataset, a quoted file specification or a previously declared dataset name (DATASET DECLARE command), enclosed in parentheses.
- If you specify both SCORE and VARIANCE on the same MATRIX subcommand, you must specify two different files.

The variables in the SCORE matrix data file and their values are:

**ROWTYPE\_** *String variable containing the value ROW for all rows and COLUMN for all columns.*

**LEVEL** *String variable containing the values (or value labels, if present) of each original variable.*

**VARNAME\_** *String variable containing the original variable names.*

**DIM1...DIMn** *Numeric variables containing the row and column scores for each dimension. Each variable is labeled DIMn, where n represents the dimension number.*

The variables in the VARIANCE matrix data file and their values are:

**ROWTYPE\_** *String variable containing the value COV for all cases in the file.*

**SCORE** *String variable containing the values SINGULAR, ROW, and COLUMN.*

**LEVEL** *String variable containing the system-missing value for SINGULAR and the sequential row or column number for ROW and COLUMN.*

**VARNAME\_** *String variable containing the dimension number.*

**DIM1...DIMn** *Numeric variables containing the covariances for each dimension. Each variable is labeled DIMn, where n represents the dimension number.*

## Analyzing Aggregated Data

To analyze aggregated data, such as data from a crosstabulation where cell counts are available but the original raw data are not, you can use the TABLE=ALL option or the WEIGHT command before ANACOR.

**Example**

To analyze a  $3 \times 3$  table, such as the table that is shown below, you could use these commands:

```
DATA LIST FREE/ BIRTHORD ANXIETY COUNT.
BEGIN DATA
1 1 48
1 2 27
1 3 22
2 1 33
2 2 20
2 3 39
3 1 29
3 2 42
3 3 47
END DATA.
WEIGHT BY COUNT.
ANACOR TABLE=BIRTHORD (1,3) BY ANXIETY (1,3).
```

- The `WEIGHT` command weights each case by the value of `COUNT`, as if there are 48 subjects with `BIRTHORD=1` and `ANXIETY=1`, 27 subjects with `BIRTHORD=1` and `ANXIETY=2`, and so on.
- `ANACOR` can then be used to analyze the data.
- If any table cell value equals 0, the `WEIGHT` command issues a warning, but the `ANACOR` analysis is done correctly.
- The table cell values (the `WEIGHT` values) cannot be negative. `WEIGHT` changes system-missing values and negative values to 0.
- For large aggregated tables, you can use the `TABLE=ALL` option or the transformation language to enter the table “as is.”

Table 13-1  
 $3 \times 3$  table

		Anxiety		
		High	Med	Low
Birth order	First	48	27	22
	Second	33	20	39
	Other	29	42	47

# ANOVA

ANOVA is available in the Statistics Base option.

```
ANOVA VARIABLES= varlist BY varlist(min,max) . . . varlist(min,max)
  [WITH varlist] [/VARIABLES=...]

  [/COVARIATES={FIRST**}]
    {WITH   }
    {AFTER  }

  [/MAXORDERS={ALL**}]
    {n      }
    {NONE   }

  [/METHOD={UNIQUE**}]
    {EXPERIMENTAL}
    {HIERARCHICAL}

  [/STATISTICS=[MCA] [REG†] [MEAN] [ALL] [NONE]]

  [/MISSING={EXCLUDE**}]
    {INCLUDE  }
```

**\*\***Default if the subcommand is omitted.

†REG (table of regression coefficients) is displayed only if the design is relevant.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
ANOVA VARIABLES=PRESTIGE BY REGION(1,9) SEX,RACE(1,2)
  /MAXORDERS=2
  /STATISTICS=MEAN.
```

## Overview

ANOVA performs analysis of variance for factorial designs. The default is the full factorial model if there are five or fewer factors. Analysis of variance tests the hypothesis that the group means of the dependent variable are equal. The dependent variable is interval-level, and one or more categorical variables define the groups. These categorical variables are termed **factors**. ANOVA also allows you to include continuous explanatory variables, termed **covariates**. Other procedures that perform analysis of variance are ONEWAY, SUMMARIZE, and GLM. To perform a comparison of two means, use TTEST.

### Options

**Specifying Covariates.** You can introduce covariates into the model using the WITH keyword on the VARIABLES subcommand.



**Order of Entry of Covariates.** By default, covariates are processed before main effects for factors. You can process covariates with or after main effects for factors using the `COVARIATES` subcommand.

**Suppressing Interaction Effects.** You can suppress the effects of various orders of interaction using the `MAXORDERS` subcommand.

**Methods for Decomposing Sums of Squares.** By default, the regression approach (keyword `UNIQUE`) is used. You can request the classic experimental or hierarchical approach using the `METHOD` subcommand.

**Statistical Display.** Using the `STATISTICS` subcommand, you can request means and counts for each dependent variable for groups defined by each factor and each combination of factors up to the fifth level. You also can request unstandardized regression coefficients for covariates and multiple classification analysis (MCA) results, which include the MCA table, the Factor Summary table, and the Model Goodness of Fit table. The MCA table shows **treatment effects** as deviations from the grand mean and includes a listing of unadjusted category effects for each factor, category effects adjusted for other factors, and category effects adjusted for all factors and covariates. The Factor Summary table displays eta and beta values. The Goodness of Fit table shows  $R$  and  $R^2$  for each model.

### ***Basic Specification***

- The basic specification is a single `VARIABLES` subcommand with an analysis list. The minimum analysis list specifies a list of dependent variables, the keyword `BY`, a list of factor variables, and the minimum and maximum integer values of the factors in parentheses.
- By default, the model includes all interaction terms up to five-way interactions. The sums of squares are decomposed using the regression approach, in which all effects are assessed simultaneously, with each effect adjusted for all other effects in the model. A case that has a missing value for any variable in an analysis list is omitted from the analysis.

### ***Subcommand Order***

- The subcommands can be named in any order.

### ***Operations***

A separate analysis of variance is performed for each dependent variable in an analysis list, using the same factors and covariates.

### ***Limitations***

- A maximum of 5 analysis lists.
- A maximum of 5 dependent variables per analysis list.
- A maximum of 10 factor variables per analysis list.
- A maximum of 10 covariates per analysis list.
- A maximum of 5 interaction levels.
- A maximum of 25 value labels per variable displayed in the MCA table.

- The combined number of categories for all factors in an analysis list plus the number of covariates must be less than the sample size.

## Examples

```
ANOVA VARIABLES=PRESTIGE BY REGION(1,9) SEX, RACE(1,2)
/MAXORDERS=2
/STATISTICS=MEAN.
```

- `VARIABLES` specifies a three-way analysis of variance—*PRESTIGE* by *REGION*, *SEX*, and *RACE*.
- The variables *SEX* and *RACE* each have two categories, with values 1 and 2 included in the analysis. *REGION* has nine categories, valued 1 through 9.
- `MAXORDERS` examines interaction effects up to and including the second order. All three-way interaction terms are pooled into the error sum of squares.
- `STATISTICS` requests a table of means of *PRESTIGE* within the combined categories of *REGION*, *SEX*, and *RACE*.

### Example: Specifying Multiple Analyses

```
ANOVA VARIABLES=PRESTIGE BY REGION(1,9) SEX, RACE(1,2)
/RINCOME BY SEX, RACE(1,2).
```

- `ANOVA` specifies a three-way analysis of variance of *PRESTIGE* by *REGION*, *SEX*, and *RACE*, and a two-way analysis of variance of *RINCOME* by *SEX* and *RACE*.

## VARIABLES Subcommand

`VARIABLES` specifies the analysis list.

- More than one design can be specified on the same `ANOVA` command by separating the analysis lists with a slash.
- Variables named before the keyword `BY` are dependent variables. Value ranges are not specified for dependent variables.
- Variables named after `BY` are factor (independent) variables.
- Every factor variable must have a value range indicating its minimum and maximum values. The values must be separated by a space or a comma and enclosed in parentheses.
- Factor variables must have integer values. Non-integer values for factors are truncated.
- Cases with values outside the range specified for a factor are excluded from the analysis.
- If two or more factors have the same value range, you can specify the value range once following the last factor to which it applies. You can specify a single range that encompasses the ranges of all factors on the list. For example, if you have two factors, one with values 1 and 2 and the other with values 1 through 4, you can specify the range for both as 1,4. However, this may reduce performance and cause memory problems if the specified range is larger than some of the actual ranges.

- Variables named after the keyword `WITH` are covariates.
- Each analysis list can include only one `BY` and one `WITH` keyword.

## ***COVARIATES Subcommand***

`COVARIATES` specifies the order for assessing blocks of covariates and factor main effects.

- The order of entry is irrelevant when `METHOD=UNIQUE`.

<b>FIRST</b>	<i>Process covariates before factor main effects.</i> This is the default.
<b>WITH</b>	<i>Process covariates concurrently with factor main effects.</i>
<b>AFTER</b>	<i>Process covariates after factor main effects.</i>

## ***MAXORDERS Subcommand***

`MAXORDERS` suppresses the effects of various orders of interaction.

<b>ALL</b>	<i>Examine all interaction effects up to and including the fifth order.</i> This is the default.
<b>n</b>	<i>Examine all interaction effects up to and including the nth order.</i> For example, <code>MAXORDERS=3</code> examines all interaction effects up to and including the third order. All higher-order interaction sums of squares are pooled into the error term.
<b>NONE</b>	<i>Delete all interaction terms from the model.</i> All interaction sums of squares are pooled into the error sum of squares. Only main and covariate effects appear in the ANOVA table.

## ***METHOD Subcommand***

`METHOD` controls the method for decomposing sums of squares.

<b>UNIQUE</b>	<i>Regression approach.</i> <code>UNIQUE</code> overrides any keywords on the <code>COVARIATES</code> subcommand. All effects are assessed simultaneously for their partial contribution. The <code>MCA</code> and <code>MEAN</code> specifications on the <code>STATISTICS</code> subcommand are not available with the regression approach. This is the default if <code>METHOD</code> is omitted.
<b>EXPERIMENTAL</b>	<i>Classic experimental approach.</i> Covariates, main effects, and ascending orders of interaction are assessed separately in that order.
<b>HIERARCHICAL</b>	<i>Hierarchical approach.</i>

### ***Regression Approach***

All effects are assessed simultaneously, with each effect adjusted for all other effects in the model. This is the default when the `METHOD` subcommand is omitted. Since `MCA` tables cannot be produced when the regression approach is used, specifying `MCA` or `ALL` on `STATISTICS` with the default method triggers a warning.

Some restrictions apply to the use of the regression approach:

- The lowest specified categories of all the independent variables must have a marginal frequency of at least 1, since the lowest specified category is used as the reference category. If this rule is violated, no ANOVA table is produced and a message identifying the first offending variable is displayed.
- Given an  $n$ -way crosstabulation of the independent variables, there must be no empty cells defined by the lowest specified category of any of the independent variables. If this restriction is violated, one or more levels of interaction effects are suppressed and a warning message is issued. However, this constraint does not apply to categories defined for an independent variable but not occurring in the data. For example, given two independent variables, each with categories of 1, 2, and 4, the (1,1), (1,2), (1,4), (2,1), and (4,1) cells must not be empty. The (1,3) and (3,1) cells will be empty but the restriction on empty cells will not be violated. The (2,2), (2,4), (4,2), and (4,4) cells may be empty, although the degrees of freedom will be reduced accordingly.

To comply with these restrictions, specify precisely the lowest non-empty category of each independent variable. Specifying a value range of (0,9) for a variable that actually has values of 1 through 9 results in an error, and no ANOVA table is produced.

### ***Classic Experimental Approach***

Each type of effect is assessed separately in the following order (unless WITH or AFTER is specified on the COVARIATES subcommand):

- Effects of covariates
- Main effects of factors
- Two-way interaction effects
- Three-way interaction effects
- Four-way interaction effects
- Five-way interaction effects

The effects within each type are adjusted for all other effects of that type and also for the effects of all prior types. (See [Table 14-1](#) on p. 207.)

### ***Hierarchical Approach***

The hierarchical approach differs from the classic experimental approach only in the way it handles covariate and factor main effects. In the hierarchical approach, factor main effects and covariate effects are assessed hierarchically—factor main effects are adjusted only for the factor main effects already assessed, and covariate effects are adjusted only for the covariates already assessed. (See [Table 14-1](#) on p. 207.) The order in which factors are listed on the ANOVA command determines the order in which they are assessed.

## Example

The following analysis list specifies three factor variables named *A*, *B*, and *C*:

```
ANOVA VARIABLES=Y BY A,B,C(0,3).
```

The following table summarizes the three methods for decomposing sums of squares for this example.

- With the default *regression* approach, each factor or interaction is assessed with all other factors and interactions held constant.
- With the *classic experimental* approach, each main effect is assessed with the two other main effects held constant, and two-way interactions are assessed with all main effects and other two-way interactions held constant. The three-way interaction is assessed with all main effects and two-way interactions held constant.
- With the *hierarchical* approach, the factor main effects *A*, *B*, and *C* are assessed with all prior main effects held constant. The order in which the factors and covariates are listed on the ANOVA command determines the order in which they are assessed in the hierarchical analysis. The interaction effects are assessed the same way as in the experimental approach.

Table 14-1  
Terms adjusted for under each option

Effect	Regression (UNIQUE)	Experimental	Hierarchical
A	All others	B,C	None
B	All others	A,C	A
C	All others	A,B	A,B
AB	All others	A,B,C,AC,BC	A,B,C,AC,BC
AC	All others	A,B,C,AB,BC	A,B,C,AB,BC
BC	All others	A,B,C,AB,AC	A,B,C,AB,AC
ABC	All others	A,B,C,AB,AC,BC	A,B,C,AB,AC,BC

## Summary of Analysis Methods

The following table describes the results obtained with various combinations of methods for controlling the entry of covariates and decomposing the sums of squares.

Table 14-2  
Combinations of COVARIATES and METHOD subcommands

Method	Assessments between types of effects	Assessments within the same type of effect
METHOD=UNIQUE	<b>Covariates, Factors, and Interactions</b> simultaneously	<b>Covariates:</b> adjust for factors, interactions, and all other covariates <b>Factors:</b> adjust for covariates, interactions, and all other factors <b>Interactions:</b> adjust for covariates, factors, and all other interactions

Method	Assessments between types of effects	Assessments within the same type of effect
METHOD=EXPERIMENTAL	<b>Covariates</b> then <b>Factors</b> then <b>Interactions</b>	<b>Covariates:</b> adjust for all other covariates <b>Factors:</b> adjust for covariates and all other factors <b>Interactions:</b> adjust for covariates, factors, and all other interactions of the same and lower orders
METHOD=HIERARCHICAL	<b>Covariates</b> then <b>Factors</b> then <b>Interactions</b>	<b>Covariates:</b> adjust for covariates that are preceding in the list <b>Factors:</b> adjust for covariates and factors preceding in the list <b>Interactions:</b> adjust for covariates, factors, and all other interactions of the same and lower orders
COVARIATES=WITH and METHOD=EXPERIMENTAL	<b>Factors and Covariates</b> concurrently then <b>Interactions</b>	<b>Covariates:</b> adjust for factors and all other covariates <b>Factors:</b> adjust for covariates and all other factors <b>Interactions:</b> adjust for covariates, factors, and all other interactions of the same and lower orders
COVARIATES=WITH and METHOD=HIERARCHICAL	<b>Factors and Covariates</b> concurrently then <b>Interactions</b>	<b>Factors:</b> adjust only for preceding factors <b>Covariates:</b> adjust for factors and preceding covariates <b>Interactions:</b> adjust for covariates, factors, and all other interactions of the same and lower orders
COVARIATES=AFTER and METHOD=EXPERIMENTAL	<b>Factors</b> then <b>Covariates</b> then <b>Interactions</b>	<b>Factors:</b> adjust for all other factors <b>Covariates:</b> adjust for factors and all other covariates <b>Interactions:</b> adjust for covariates, factors, and all other interactions of the same and lower orders
COVARIATES=AFTER and METHOD=HIERARCHICAL	<b>Factors</b> then <b>Covariates</b> then <b>Interactions</b>	<b>Factors:</b> adjust only for preceding factors <b>Covariates:</b> adjust factors and preceding covariates <b>Interactions:</b> adjust for covariates, factors, and all other interactions of the same and lower orders

## ***STATISTICS Subcommand***

STATISTICS requests additional statistics. STATISTICS can be specified by itself or with one or more keywords.

- If you specify `STATISTICS` without keywords, ANOVA calculates `MEAN` and `REG` (each defined below).
- If you specify a keyword or keywords on the `STATISTICS` subcommand, ANOVA calculates only the additional statistics you request.

<b>MEAN</b>	<i>Means and counts table.</i> This statistic is not available when <code>METHOD</code> is omitted or when <code>METHOD=UNIQUE</code> . See “Cell Means” below.
<b>REG</b>	<i>Unstandardized regression coefficients.</i> Displays unstandardized regression coefficients for the covariates. For more information, see the topic <a href="#">Regression Coefficients for the Covariates</a> on p. 209.
<b>MCA</b>	<i>Multiple classification analysis.</i> The MCA, the Factor Summary, and the Goodness of Fit tables are not produced when <code>METHOD</code> is omitted or when <code>METHOD=UNIQUE</code> . For more information, see the topic <a href="#">Multiple Classification Analysis</a> on p. 209.
<b>ALL</b>	<i>Means and counts table, unstandardized regression coefficients, and multiple classification analysis.</i>
<b>NONE</b>	<i>No additional statistics.</i> ANOVA calculates only the statistics needed for analysis of variance. This is the default if the <code>STATISTICS</code> subcommand is omitted.

## Cell Means

`STATISTICS=MEAN` displays the Cell Means table.

- This statistic is not available with `METHOD=UNIQUE`.
- The Cell Means table shows the means and counts of each dependent variable for each cell defined by the factors and combinations of factors. Dependent variables and factors appear in their order on the `VARIABLES` subcommand.
- If `MAXORDERS` is used to suppress higher-order interactions, cell means corresponding to suppressed interaction terms are not displayed.
- The means displayed are the observed means in each cell, and they are produced only for dependent variables, not for covariates.

## Regression Coefficients for the Covariates

`STATISTICS=REG` requests the unstandardized regression coefficients for the covariates.

- The regression coefficients are computed at the point where the covariates are entered into the equation. Thus, their values depend on the type of design specified by the `COVARIATES` or `METHOD` subcommand.
- The coefficients are displayed in the ANOVA table.

## Multiple Classification Analysis

`STATISTICS=MCA` displays the MCA, the Factor Summary, and the Model Goodness of Fit tables.

- The MCA table presents counts, predicted means, and deviations of predicted means from the grand mean for each level of each factor. The predicted and deviation means each appear in up to three forms: unadjusted, adjusted for other factors, and adjusted for other factors and covariates.

- The Factor Summary table displays the correlation ratio ( $\eta$ ) with the unadjusted deviations (the square of  $\eta$  indicates the proportion of variance explained by all categories of the factor), a partial beta equivalent to the standardized partial regression coefficient that would be obtained by assigning the unadjusted deviations to each factor category and regressing the dependent variable on the resulting variables, and the parallel partial betas from a regression that includes covariates in addition to the factors.
- The Model Goodness of Fit table shows  $R$  and  $R^2$  for each model.
- The tables cannot be produced if `METHOD` is omitted or if `METHOD=UNIQUE`. When produced, the MCA table does not display the values adjusted for factors if `COVARIATES` is omitted, if `COVARIATES=FIRST`, or if `COVARIATES=WITH` and `METHOD=EXPERIMENTAL`. A full MCA table is produced only if `METHOD=HIERARCHICAL` or if `METHOD=EXPERIMENTAL` and `COVARIATES=AFTER`.

## ***MISSING Subcommand***

By default, a case that has a missing value for any variable named in the analysis list is deleted for all analyses specified by that list. Use `MISSING` to include cases with user-missing data.

- EXCLUDE**            *Exclude cases with missing data. This is the default.*  
**INCLUDE**            *Include cases with user-defined missing data.*

## ***References***

Andrews, F., J. Morgan, J. Sonquist, and L. Klein. 1973. *Multiple classification analysis*, 2nd ed. Ann Arbor: University of Michigan.



# APPLY DICTIONARY

```
APPLY DICTIONARY FROM [{'savfile'|'dataset'}]
                        {*          }

[/SOURCE VARIABLES = varlist]
[/TARGET VARIABLES = varlist]
[/NEWVARS]
[/FILEINFO [ATTRIBUTES = [{REPLACE}]]
           {MERGE  }

           [DOCUMENTS = [{REPLACE}]]  ]
           {MERGE  }

           [FILELABEL]
           [MRSETS = [{REPLACE}]]
           {MERGE  }

           [VARSETS = [{REPLACE}]]
           {MERGE  }

           [WEIGHT**]

           [ALL]

[/VARINFO [ALIGNMENT**]  ]

           [ATTRIBUTES = [{REPLACE}]]
           {MERGE  }
           [FORMATS**]
           [LEVEL**]
           [MISSING**]
           [ROLE**]
           [VALLABELS = [{REPLACE**}]]
           {MERGE  }
           [VARLABEL**]
           [WIDTH**]
           [ALL]
```

\*\*Default if the subcommand is not specified.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- ATTRIBUTES keyword introduced on FILEINFO and VARINFO subcommands.

Release 18

- ROLE keyword introduced on VARINFO subcommands.

**Example**

```
APPLY DICTIONARY FROM = 'lastmonth.sav'.
```

**Overview**

APPLY DICTIONARY can apply variable and file-based dictionary information from an external IBM® SPSS® Statistics data file or open dataset to the current active dataset. Variable-based dictionary information in the current active dataset can be applied to other variables in the current active dataset.

- The applied variable information includes variable and value labels, missing-value flags, alignments, variable print and write formats, measurement levels, and widths.
- The applied file information includes variable and multiple response sets, documents, file label, and weight.
- APPLY DICTIONARY can apply information selectively to variables and can apply selective file-based dictionary information.
- Individual variable attributes can be applied to individual and multiple variables of the same type (strings of the same character length or numeric).
- APPLY DICTIONARY can add new variables but cannot remove variables, change data, or change a variable's name or type.
- Undefined (empty) attributes in the source dataset do not overwrite defined attributes in the active dataset.

**Basic Specification**

The basic specification is the FROM subcommand and the name of an external SPSS Statistics data file or open dataset. The file specification should be enclosed in quotation marks.

**Subcommand Order**

The subcommands can be specified in any order.

**Syntax Rules**

- The file containing the dictionary information to be applied (the **source file**) must be an external SPSS Statistics data file or a currently open dataset.
- The file to which the dictionary information is applied (the **target file**) must be the active dataset. You cannot specify another file.
- If a subcommand is issued more than once, APPLY DICTIONARY will ignore all but the last instance of the subcommand.
- Equals signs displayed in the syntax chart and in the examples presented here are required elements; they are not optional.

### **Matching Variable Type**

APPLY DICTIONARY considers two variables to have a matching variable type if:

- Both variables are numeric. This includes all numeric, currency, and date formats.
- Both variables are string (alphanumeric).

## **FROM Subcommand**

FROM specifies an external IBM® SPSS® Statistics data file or an open dataset as the source file whose dictionary information is to be applied to the active dataset.

- FROM is required.
- Only one SPSS Statistics data file or open dataset (including the active dataset) can be specified on FROM.
- The file specification should be enclosed in quotation marks.
- The active dataset can be specified in the FROM subcommand by using an asterisk (\*) as the value. File-based dictionary information (FILEINFO subcommand) is ignored when the active dataset is used as the source file.

### **Example**

```
APPLY DICTIONARY FROM "lastmonth.sav".
```

- This will apply variable information from *lastmonth.sav* to matching variables in the active dataset.
- The default variable information applied from the source file includes variable labels, value labels, missing values, level of measurement, alignment, column width (for Data Editor display), and print and write formats.
- If weighting is on in the source dataset and a matching weight variable exists in the active (target) dataset, weighting by that variable is turned on in the active dataset. No other file information (documents, file label, multiple response sets) from the source file is applied to the active dataset.

## **NEWVARS Subcommand**

NEWVARS is required to create new variables in the active (target) dataset.

### **Example**

```
APPLY DICTIONARY FROM "lastmonth.sav"  
/NEWVARS.
```

- For a new, blank active dataset, all variables with all of their variable definition attributes are copied from the source dataset, creating a new dataset with an identical set of variables (but no data values).
- For an active dataset that contains any variables, variable definition attributes from the source dataset are applied to the matching variables in the active (target) dataset. If the source dataset contains any variables that are not present in the active dataset (determined by variable name), these variables are created in the active dataset.

## ***SOURCE and TARGET Subcommands***

The **SOURCE** subcommand is used to specify variables in the source file from which to apply variable definition attributes. The **TARGET** subcommand is used to specify variables in the active dataset to which to apply variable definition attributes.

- All variables specified in the **SOURCE** subcommand must exist in the source file.
- If the **TARGET** subcommand is specified without the **SOURCE** subcommand, all variables specified must exist in the source file.
- If the **NEWVARS** subcommand is specified, variables that are specified in the **SOURCE** subcommand that exist in the source file but not in the target file will be created in the target file as new variables using the variable definition attributes (variable and value labels, missing values, etc.) from the source variable.
- For variables with matching name and type, variable definition attributes from the source variable are applied to the matching target variable.
- If both **SOURCE** and **TARGET** are specified, the **SOURCE** subcommand can specify only one variable. Variable definition attributes from that single variable in the **SOURCE** subcommand are applied to all variables of the matching type. When applying the attributes of one variable to many variables, all variables specified in the **SOURCE** and **TARGET** subcommands must be of the same type.
- For variables with matching names but different types, only variable labels are applied to the target variables.

Table 15-1  
Variable mapping for **SOURCE** and **TARGET** subcommands

<b>SOURCE subcommand</b>	<b>TARGET subcommand</b>	<b>Variable mapping</b>
none	none	Variable definition attributes from the source dataset are applied to matching variables in the active (target) dataset. New variables may be created if the <b>NEWVARS</b> subcommand is specified.
many	none	Variable definition attributes for the specified variables are copied from the source dataset to the matching variables in the active (target) dataset. All specified variables must exist in the source dataset. New variables may be created if the <b>NEWVARS</b> subcommand is specified.
none	many	Variable definition attributes for the specified variables are copied from the source dataset to the matching variables in the active (target) dataset. All specified variables must exist in the source dataset. New variables may be created if the <b>NEWVARS</b> subcommand is specified.

SOURCE subcommand	TARGET subcommand	Variable mapping
one	many	Variable definition attributes for the specified variable in the source dataset are applied to all specified variables in the active (target) dataset that have a matching type. New variables may be created if the NEWVARS subcommand is specified.
many	many	Invalid. Command not executed.

**Example**

```
APPLY DICTIONARY from *
  /SOURCE VARIABLES = var1
  /TARGET VARIABLES = var2 var3 var4
  /NEWVARS.
```

- Variable definition attributes for *var1* in the active dataset are copied to *var2*, *var3*, and *var4* in the same dataset if they have a matching type.
- Any variables specified in the TARGET subcommand that do not already exist are created, using the variable definition attributes of the variable specified in the SOURCE subcommand.

**Example**

```
APPLY DICTIONARY from "lastmonth.sav"
  /SOURCE VARIABLES = var1, var2, var3.
```

- Variable definition attributes from the specified variables in the source dataset are applied to the matching variables in the active dataset.
- For variables with matching names but different types, only variable labels from the source variable are copied to the target variable.
- In the absence of a NEWVARS subcommand, no new variables will be created.

**FILEINFO Subcommand**

FILEINFO applies global file definition attributes from the source dataset to the active (target) dataset.

- File definition attributes in the active dataset that are undefined in the source dataset are not affected.
- This subcommand is ignored if the source dataset is the active dataset.
- This subcommand is ignored if no keywords are specified.
- For keywords that contain an associated value, the equals sign between the keyword and the value is required—for example, DOCUMENTS = MERGE.

**ATTRIBUTES**

*Applies file attributes defined by the DATAFILE ATTRIBUTE command. You can REPLACE or MERGE file attributes.*

**DOCUMENTS**

*Applies documents (defined with the DOCUMENTS command) from the source dataset to the active (target) dataset. You can REPLACE or MERGE documents. DOCUMENTS = REPLACE replaces any documents in the active dataset, deleting preexisting documents in the file. This is the default if DOCUMENTS is specified without a value.*

DOCUMENTS = MERGE merges documents from the source and active datasets. Unique documents in the source file that don't exist in the active dataset are added to the active dataset. All documents are then sorted by date.

**FILELABEL**

*Replaces the file label (defined with the FILE LABEL command).*

**MRSETS**

*Applies multiple response set definitions from the source dataset to the active dataset. Multiple response sets that contain no variables in the active dataset (including variables added by the same APPLY DICTIONARY command) are ignored. You can REPLACE or MERGE multiple response sets.*

MRSETS = REPLACE deletes any existing multiple response sets in the active dataset, replacing them with multiple response sets from the source dataset.

MRSETS = MERGE adds multiple response sets from the source dataset to the collection of multiple response sets in the active dataset. If a set with the same name exists in both files, the existing set in the active dataset is unchanged.

**VARSETS**

*Applies variable set definitions from the source dataset to the active dataset. Variable sets are used to control the list of variables that are displayed in dialog boxes. Variable sets are defined by selecting Define Variable Sets from the Utilities menu. Sets in the source data file that don't contain any variables in the active dataset are ignored unless those variables are created by the same APPLY DICTIONARY command. You can REPLACE or MERGE variable sets.*

VARSETS = REPLACE deletes any existing variable sets in the active dataset, replacing them with variable sets from the source dataset.

VARSETS = MERGE adds variable sets from the source dataset to the collection of variable sets in the active dataset. If a set with the same name exists in both files, the existing set in the active dataset is unchanged.

**WEIGHT**

*Weights cases by the variable specified in the source file if there's a matching variable in the target file. This is the default if the subcommand is omitted.*

**ALL**

*Applies all file information from the source dataset to the active dataset. Documents, multiple response sets, and variable sets are merged, not replaced. File definition attributes in the active dataset that are undefined in the source data file are not affected.*

**Example**

```
APPLY DICTIONARY FROM "lastmonth.sav"
  /FILEINFO DOCUMENTS = REPLACE MRSETS = MERGE.
```

- Documents in the source dataset replace documents in the active dataset unless there are no defined documents in the source dataset.
- Multiple response sets from the source dataset are added to the collection of defined multiple response sets in the active dataset. Sets in the source dataset that contain variables that don't exist in the active dataset are ignored. If the same set name exists in both datasets, the set in the active dataset remains unchanged.

## **VARINFO Subcommand**

VARINFO applies variable definition attributes from the source dataset to the matching variables in the active dataset. With the exception of VALLABELS, all keywords replace the variable definition attributes in the active dataset with the attributes from the matching variables in the source dataset.

<b>ALIGNMENT</b>	<i>Applies variable alignment for Data Editor display.</i> This setting affects alignment (left, right, center) only in the Data View display of the Data Editor.
<b>ATTRIBUTES</b>	<i>Applies variable attributes defined by the VARIABLE ATTRIBUTE command.</i> You can REPLACE or MERGE variable attributes.
<b>FORMATS</b>	<i>Applies variable print and write formats.</i> This is the same variable definition attribute that can be defined with the FORMATS command. This setting is primarily applicable only to numeric variables. For string variables, this affects only the formats if the source or target variable is AHEX format and the other is A format.
<b>LEVEL</b>	<i>Applies variable measurement level (nominal, ordinal, scale).</i> This is the same variable definition attribute that can be defined with the VARIABLE LEVEL command.
<b>MISSING</b>	<i>Applies variable missing value definitions.</i> Any existing defined missing values in the matching variables in the active dataset are deleted. This is the same variable definition attribute that can be defined with the MISSING VALUES command. Missing values definitions are not applied to string variables if the source variable contains missing values of a longer width than the defined width of the target variable.
<b>MISSING</b>	<i>Applies variable missing value definitions.</i> Any existing defined missing values in the matching variables in the active dataset are deleted. This is the same variable definition attribute that can be defined with the MISSING VALUES command. Missing values definitions are not applied to string variables if the source variable contains missing values of a longer width than the defined width of the target variable.
<b>MISSING</b>	<i>Applies variable missing value definitions.</i> Any existing defined missing values in the matching variables in the active dataset are deleted. This is the same variable definition attribute that can be defined with the MISSING VALUES command. Missing values definitions are not applied to string variables if the source variable contains missing values of a longer width than the defined width of the target variable.
<b>MISSING</b>	<i>Applies variable missing value definitions.</i> Any existing defined missing values in the matching variables in the active dataset are deleted. This is the same variable definition attribute that can be defined with the MISSING VALUES command. Missing values definitions are not applied to string variables if the source variable contains missing values of a longer width than the defined width of the target variable.
<b>ROLE</b>	<i>Applies role assignments.</i> <a href="#">For more information, see the topic Overview on p. 2133.</a>
<b>VALLABELS</b>	<i>Applies value label definitions.</i> Value labels are not applied to string variables if the source variable contains defined value labels for values longer than the defined width of the target variable. You can REPLACE or MERGE value labels.  VALLABELS = REPLACE replaces any defined value labels from variable in the active dataset with the value labels from the matching variable in the source dataset.  VALLABELS = MERGE merges defined value labels for matching variables. If the same value has a defined value label in both the source and active datasets, the value label in the active dataset is unchanged.
<b>WIDTH</b>	<i>Display column width in the Data Editor.</i> This affects only column width in Data View in the Data Editor. It has no effect on the defined width of the variable.

**Example**

```
APPLY DICTIONARY from "lastmonth.sav"  
  /VARINFO LEVEL MISSING VALLABELS = MERGE.
```

- The level of measurement and defined missing values from the source dataset are applied to the matching variables in the active (target) dataset. Any existing missing values definitions for those variables in the active dataset are deleted.
- Value labels for matching variables in the two datasets are merged. If the same value has a defined value label in both the source and active datasets, the value label in the active dataset is unchanged.



# AUTORECODE

```
AUTORECODE VARIABLES=varlist  
  
  /INTO new varlist  
  
  [/BLANK={VALID**}  
    {MISSING}]  
  
  [/GROUP]  
  
  [/APPLY TEMPLATE='filespec']  
  
  [/SAVE TEMPLATE='filespec']  
  
  [/DESCENDING]  
  
  [/PRINT]
```

**\*\***Default if the subcommand omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 13.0

- BLANK subcommand introduced.
- GROUP subcommand introduced.
- APPLY TEMPLATE and SAVE TEMPLATE subcommands introduced.

## **Example**

```
AUTORECODE VARIABLES=Company /INTO Rcompany.
```

## **Overview**

AUTORECODE recodes the values of string and numeric variables to consecutive integers and puts the recoded values into a new variable called a **target variable**. The value labels or values of the original variable are used as value labels for the target variable. AUTORECODE is useful for creating numeric independent (grouping) variables from string variables for procedures such as ONEWAY and DISCRIMINANT. AUTORECODE can also recode the values of factor variables to consecutive integers, which may be required by some procedures and which reduces the amount of workspace needed by some statistical procedures.

## **Basic Specification**

The basic specification is VARIABLES and INTO. VARIABLES specifies the variables to be recoded. INTO provides names for the target variables that store the new values. VARIABLES and INTO must name or imply the same number of variables.

**Subcommand Order**

- VARIABLES must be specified first.
- INTO must immediately follow VARIABLES.
- All other subcommands can be specified in any order.

**Syntax Rules**

- A variable cannot be recoded into itself. More generally, target variable names cannot duplicate any variable names already in the working file.
- If the GROUP or APPLY TEMPLATE subcommand is specified, all variables on the VARIABLES subcommand must be the same type (numeric or string).
- If APPLY TEMPLATE is specified, all variables on the VARIABLES subcommand must be the same type (numeric or string) as the type defined in the template.
- File specifications on the APPLY TEMPLATE and SAVE TEMPLATE subcommands follow the normal conventions for file specifications. Enclosing file specifications in quotation marks is recommended.

**Operations**

- The values of each variable to be recoded are sorted and then assigned numeric values. By default, the values are assigned in ascending order: 1 is assigned to the lowest nonmissing value of the original variable; 2, to the second-lowest nonmissing value; and so on, for each value of the original variable.
- Values of the original variables are unchanged.
- Missing values are recoded into values higher than any nonmissing values, with their order preserved. For example, if the original variable has 10 nonmissing values, the first missing value is recoded as 11 and retains its user-missing status. System-missing values remain system-missing. (See the GROUP, APPLY TEMPLATE, and SAVE TEMPLATE subcommands for additional rules for user-missing values.)
- AUTORECODE does not sort the cases in the working file. As a result, the consecutive numbers assigned to the target variables may not be in order in the file.
- Target variables are assigned the same variable labels as the original source variables. To change the variable labels, use the VARIABLE LABELS command after AUTORECODE.
- Value labels are automatically generated for each value of the target variables. If the original value had a label, that label is used for the corresponding new value. If the original value did not have a label, the old value itself is used as the value label for the new value. The defined print format of the old value is used to create the new value label.
- AUTORECODE ignores SPLIT FILE specifications. However, any SELECT IF specifications are in effect for AUTORECODE.

**Example**

```
DATA LIST / COMPANY 1-21 (A) SALES 24-28.  
BEGIN DATA  
CATFOOD JOY           10000  
OLD FASHIONED CATFOOD 11200
```

```

. . .
PRIME CATFOOD          10900
CHOICE CATFOOD        14600
END DATA.

AUTORECODE VARIABLES=COMPANY /INTO=RCOMPANY /PRINT.

TABLES TABLE = SALES BY RCOMPANY
  /TTITLE='CATFOOD SALES BY COMPANY'.

```

- AUTORECODE recodes *COMPANY* into a numeric variable *RCOMPANY*. Values of *RCOMPANY* are consecutive integers beginning with 1 and ending with the number of different values entered for *COMPANY*. The values of *COMPANY* are used as value labels for *RCOMPANY*'s numeric values. The *PRINT* subcommand displays a table of the original and recoded values.

## VARIABLES Subcommand

VARIABLES specifies the variables to be recoded. VARIABLES is required and must be specified first. The actual keyword VARIABLES is optional.

- Values from the specified variables are recoded and stored in the target variables listed on INTO. Values of the original variables are unchanged.

## INTO Subcommand

INTO provides names for the target variables that store the new values. INTO is required and must immediately follow VARIABLES.

- The number of target variables named or implied on INTO must equal the number of source variables listed on VARIABLES.

### Example

```
AUTORECODE VARIABLES=V1 V2 V3 /INTO=NEWV1 TO NEWV3 /PRINT.
```

- AUTORECODE stores the recoded values of *V1*, *V2*, and *V3* into target variables named *NEWV1*, *NEWV2*, and *NEWV3*.

## BLANK Subcommand

The BLANK subcommand specifies how to autorecode blank string values.

- BLANK is followed by an equals sign (=) and the keyword VALID or MISSING.
- The BLANK subcommand applies only to string variables (both short and long strings). System-missing numeric values remain system-missing in the new, autorecoded variable(s).

- The `BLANK` subcommand has no effect if there are no string variables specified on the `VARIABLES` subcommand.

**VALID**            *Blank string values are treated as valid, nonmissing values and are autorecoded into nonmissing values. This is the default.*

**MISSING**        *Blank string values are autorecoded into a user-missing value higher than the highest nonmissing value.*

### **Example**

```
DATA LIST /stringVar (A1).
BEGIN DATA
a
b

c
d
END DATA.
AUTORECODE
  VARIABLES=stringVar /INTO NumericVar
  /BLANK=MISSING.
```

- The values *a*, *b*, *c*, and *d* are autorecoded into the numeric values 1 through 4.
- The blank value is autorecoded to 5, and 5 is defined as user-missing.

## **GROUP Subcommand**

The subcommand `GROUP` allows you to specify that a single autorecoding scheme should be generated for all the specified variables, yielding consistent coding for all of the variables.

- The `GROUP` subcommand has no additional keywords or specifications. By default, variables are not grouped for autorecoding.
- All variables must be the same type (numeric or string).
- All observed values for all specified variables are used to create a sorted order of values to recode into sequential integers.
- String variables can be of any length and can be of unequal length.
- User-missing values for the target variables are based on the first variable in the original variable list with defined user-missing values. All other values from other original variables, except for system-missing, are treated as valid.
- If only one variable is specified on the `VARIABLES` subcommand, the `GROUP` subcommand is ignored.
- If `GROUP` and `APPLY TEMPLATE` are used on the same `AUTORECODE` command, value mappings from the template are applied first. All remaining values are recoded into values higher than the last value in the template, with user-missing values (based on the first variable in the list with defined user-missing values) recoded into values higher than the last valid value. See the `APPLY TEMPLATE` subcommand for more information.

**Example**

```
DATA LIST FREE /var1 (a1) var2 (a1).
BEGIN DATA
a d
b e
c f
END DATA.
MISSING VALUES var1 ("c") var2 ("f").
AUTORECODE VARIABLES=var1 var2
/INTO newvar1 newvar2
/GROUP.
```

- A single autorecoding scheme is created and applied to both new variables.
- The user-missing value "c" from *var1* is autorecoded into a user-missing value.
- The user-missing value "f" from *var2* is autorecoded into a valid value.

Table 16-1

Original and recoded values

Original value	Autorecoded value
a	1
b	2
c	6 (user-missing)
d	3
e	4
f	5

**SAVE TEMPLATE Subcommand**

The `SAVE TEMPLATE` subcommand allows you to save the autorecode scheme used by the current `AUTORECODE` command to an external template file, which you can then use when autorecoding other variables using the `APPLY TEMPLATE` subcommand.

- `SAVE TEMPLATE` is followed by an equals sign (=) and a quoted file specification. The default file extension for autorecode templates is *.sat*.
- The template contains information that maps the original nonmissing values to the recoded values.
- Only information for nonmissing values is saved in the template. User-missing value information is not retained.
- If more than one variable is specified on the `VARIABLES` subcommand, the first variable specified is used for the template, unless `GROUP` or `APPLY TEMPLATE` is also specified, in which case a common autorecoding scheme for all variables is saved in the template.

**Example**

```
DATA LIST FREE /var1 (a1) var2 (a1).
BEGIN DATA
a d
b e
c f
END DATA.
```

```
MISSING VALUES var1 ("c") var2 ("f").
AUTORECODE VARIABLES=var1 var2
  /INTO newvar1 newvar2
  /SAVE TEMPLATE='/temp/var1_template.sat'.
```

- The saved template contains an autorecode scheme that maps the string values of "a" and "b" from *var1* to the numeric values 1 and 2, respectively.
- The template contains no information for the value of "c" for *var1* because it is defined as user-missing.
- The template contains no information for values associated with *var2* because the `GROUP` subcommand was not specified.

### **Template File Format**

An autorecode template file is actually a data file in IBM® SPSS® Statistics format that contains two variables: *Source\_* contains the original, unrecoded valid values, and *Target\_* contains the corresponding recoded values. Together these two variables provide a mapping of original and recoded values.

You can therefore, theoretically, build your own custom template files, or simply include the two mapping variables in an existing data file—but this type of use has not been tested.

## **APPLY TEMPLATE Subcommand**

The `APPLY TEMPLATE` subcommand allows you to apply a previously saved autorecode template to the variables in the current `AUTORECODE` command, appending any additional values found in the variables to the end of the scheme, preserving the relationship between the original and autorecode values stored in the saved scheme.

- `APPLY TEMPLATE` is followed by an equals sign (=) and a quoted file specification.
- All variables on the `VARIABLES` subcommand must be the same type (numeric or string), and that type must match the type defined in the template.
- Templates do not contain any information on user-missing values. User-missing values for the target variables are based on the first variable in the original variable list with defined user-missing values. All other values from other original variables, except for system-missing, are treated as valid.
- Value mappings from the template are applied first. All remaining values are recoded into values higher than the last value in the template, with user-missing values (based on the first variable in the list with defined user-missing values) recoded into values higher than the last valid value.
- If multiple variables are specified on the `VARIABLES` subcommand, `APPLY TEMPLATE` generates a grouped recoding scheme, with or without an explicit `GROUP` subcommand.

### **Example**

```
DATA LIST FREE /var1 (a1).
BEGIN DATA
a b d
```

```

END DATA.
AUTORECODE VARIABLES=var1
  /INTO newvar1
  /SAVE TEMPLATE='/temp/var1_template.sat'.
DATA LIST FREE /var2 (a1).
BEGIN DATA
a b c
END DATA.
AUTORECODE VARIABLES=var2
  /INTO newvar2
  /APPLY TEMPLATE='/temp/var1_template.sat'.

```

- The template file *var1\_template.sat* maps the string values a, b, and d to the numeric values 1, 2, and 3, respectively.
- When the template is applied to the variable *var2* with the string values a, b, and c, the autorecoded values for *newvar2* are 1, 2, and 4, respectively. The string value “c” is autorecoded to 4 because the template maps 3 to the string value “d”.
- The data dictionary contains defined value labels for all four values—the three from the template and the one new value read from the file.

Table 16-2  
Defined value labels for *newvar2*

Value	Label
1	a
2	b
3	d
4	c

## ***Interaction between APPLY TEMPLATE and SAVE TEMPLATE***

- If `APPLY TEMPLATE` and `SAVE TEMPLATE` are both used in the same `AUTORECODE` command, `APPLY TEMPLATE` is always processed first, regardless of subcommand order, and the autorecode scheme saved by `SAVE TEMPLATE` is the union of the original template plus any appended value definitions.
- `APPLY TEMPLATE` and `SAVE TEMPLATE` can specify the same file, resulting in the template being updated to include any newly appended value definitions.

### ***Example***

```

AUTORECODE VARIABLES=products
  /INTO productCodes
  /APPLY TEMPLATE='/mydir/product_codes.sat'
  /SAVE TEMPLATE='/mydir/product_codes.sat'.

```

- The autorecode scheme in the template file is applied for autorecoding *products* into *productCodes*.
- Any data values for *products* not defined in the template are autorecoded into values higher than the highest value in the original template.

- Any user-missing values for *products* are autorecoded into values higher than the highest nonmissing autorecoded value.
- The template saved is the autorecode scheme used to autorecode *product*—the original autorecode scheme plus any additional values in *product* that were appended to the scheme.

### ***PRINT Subcommand***

PRINT displays a correspondence table of the original values of the source variables and the new values of the target variables. The new value labels are also displayed.

- The only specification is the keyword PRINT. There are no additional specifications.

### ***DESCENDING Subcommand***

By default, values for the source variable are recoded in ascending order (from lowest to highest). DESCENDING assigns the values to new variables in descending order (from highest to lowest). The largest value is assigned 1, the second-largest, 2, and so on.

- The only specification is the keyword DESCENDING. There are no additional specifications.



# BEGIN DATA-END DATA

```
BEGIN DATA
data records
END DATA
```

## **Example**

```
BEGIN DATA
1 3424 274 ABU DHABI 2
2 39932 86 AMSTERDAM 4
3 8889 232 ATHENS
4 3424 294 BOGOTA 3
END DATA.
```

## **Overview**

BEGIN DATA and END DATA are used when data are entered within the command sequence (inline data). BEGIN DATA and END DATA are also used for inline matrix data. BEGIN DATA signals the beginning of data lines and END DATA signals the end of data lines.

### **Basic Specification**

The basic specification is BEGIN DATA, the data lines, and END DATA. BEGIN DATA must be specified by itself on the line that immediately precedes the first data line. END DATA is specified by itself on the line that immediately follows the last data line.

### **Syntax Rules**

- BEGIN DATA, the data, and END DATA must precede the first procedure.
- The command terminator after BEGIN DATA is optional. It is best to leave it out so that the program will treat inline data as one continuous specification.
- END DATA must always begin in column 1. It must be spelled out in full and can have only one space between the words END and DATA. Procedures and additional transformations can follow the END DATA command.
- Data lines must *not* have a command terminator. For inline data formats, see DATA LIST.
- Inline data records are limited to a maximum of 80 columns. (On some systems, the maximum may be fewer than 80 columns.) If data records exceed 80 columns, they must be stored in an external file that is specified on the FILE subcommand of the DATA LIST (or similar) command.

### **Operations**

- When the program encounters BEGIN DATA, it begins to read and process data on the next input line. All preceding transformation commands are processed as the working file is built.
- The program continues to evaluate input lines as data until it encounters END DATA, at which point it begins evaluating input lines as commands.

---

*BEGIN DATA-END DATA*

- No other commands are recognized between `BEGIN DATA` and `END DATA`.
- The `INCLUDE` command can specify a file that contains `BEGIN DATA`, data lines, and `END DATA`. The data in such a file are treated as inline data. Thus, the `FILE` subcommand should be omitted from the `DATA LIST` (or similar) command.
- When running the program from prompts, the prompt `DATA>` appears immediately after `BEGIN DATA` is specified. After `END DATA` is specified, the command line prompt returns.

## **Examples**

```
DATA LIST /XVAR 1 YVAR ZVAR 3-12 CVAR 14-22 (A) JVAR 24 .
BEGIN DATA
1 3424 274 ABU DHABI 2
2 39932 86 AMSTERDAM 4
3 8889 232 ATHENS
4 3424 294 BOGOTA 3
5 11323 332 HONG KONG 3
6 323 232 MANILA 1
7 3234 899 CHICAGO 4
8 78998 2344 VIENNA 3
9 8870 983 ZURICH 5
END DATA .
MEANS XVAR BY JVAR .
```

- `DATA LIST` defines the names and column locations of the variables. The `FILE` subcommand is omitted because the data are inline.
- There are nine cases in the inline data. Each line of data completes a case.
- `END DATA` signals the end of data lines. It begins in column 1 and has only a single space between `END` and `DATA`.

# BEGIN GPL-END GPL

```
BEGIN GPL
  gpl specification
END GPL
```

## Release History

Release 14.0

- Command introduced.

## Example

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: count=col(source(s), name("COUNT"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Count"))
  ELEMENT: interval(position(jobcat*count))
END GPL.
```

If you are looking for more details about GPL, see the *GPL Reference Guide* on the manuals CD.

## Overview

BEGIN GPL and END GPL are used when Graphics Production Language (GPL) code is entered within the command sequence (inline graph specification). BEGIN GPL and END GPL must follow a GGRAPH command, without any blank lines between BEGIN GPL and the command terminator line for GGRAPH. Only comments are allowed between BEGIN GPL and the command terminator line for GGRAPH. BEGIN GPL must be at the start of the line on which it appears, with no preceding spaces. BEGIN GPL signals the beginning of GPL code, and END GPL signals the end of GPL code.

For more information about GGRAPH, see [GGRAPH on p. 864](#). See the *GPL Reference Guide* on the manuals CD for more details about GPL. The examples in the GPL documentation may look different compared to the syntax pasted from the Chart Builder. The main difference is when aggregation occurs. See [Working with the GPL on p. 875](#) for information about the differences. See [GPL Examples on p. 879](#) for examples with GPL that is similar to the pasted syntax.

## Syntax Rules

- Within a GPL block, only GPL statements are allowed.
- Strings in GPL are enclosed in quotation marks. You cannot use single quotes (apostrophes).

- With the IBM® SPSS® Statistics Batch Facility (available only with SPSS Statistics Server), use the `-i` switch when submitting command files that contain GPL blocks.

***Scope and Limitations***

- GPL blocks cannot be nested within GPL blocks.
- GPL blocks cannot be contained within `DEFINE- !ENDDDEFINE` macro definitions.
- GPL blocks can be contained in command syntax files run via the `INSERT` command, with the default `SYNTAX=INTERACTIVE` setting.
- GPL blocks cannot be contained within command syntax files run via the `INCLUDE` command.

# ***BEGIN PROGRAM-END PROGRAM***

BEGIN PROGRAM-END PROGRAM is available in the IBM® SPSS® Statistics Programmability Extension. It is not available in Statistical Services for SQL Server 2005.

```
BEGIN PROGRAM [programming language name].  
programming language-specific statements  
END PROGRAM.
```

## ***Release History***

Release 14.0

- Command introduced.

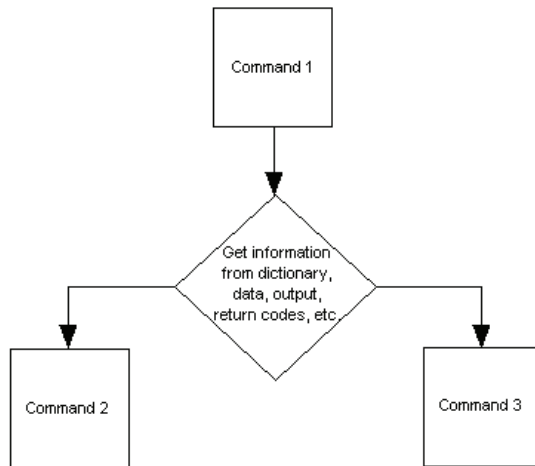
## ***Overview***

BEGIN PROGRAM-END PROGRAM provides the ability to integrate the capabilities of external programming languages with IBM® SPSS® Statistics. One of the major benefits of these program blocks is the ability to add **jobwise** flow control to the command stream. Outside of program blocks, SPSS Statistics can execute **casewise** conditional actions, based on criteria that evaluate each case, but jobwise flow control, such as running different procedures for different variables based on data type or level of measurement or determining which procedure to run next based on the results of the last procedure is much more difficult. Program blocks make jobwise flow control much easier to accomplish. With program blocks, you can control the commands that are run based on many criteria, including:

- Dictionary information (e.g., data type, measurement level, variable names)
- Data conditions
- Output values
- Error codes (that indicate if a command ran successfully or not)

You can also read data from the active dataset to perform additional computations, update the active dataset with results, create new datasets, and create custom pivot table output.

Figure 19-1  
Jobwise Flow Control



### Operations

- `BEGIN PROGRAM` signals the beginning of a set of code instructions controlled by an external programming language.
- After `BEGIN PROGRAM` is executed, other commands do not execute until `END PROGRAM` is encountered.

### Syntax Rules

- Within a program block, only statements recognized by the specified programming language are allowed.
- Command syntax generated within a program block must follow **interactive** syntax rules. [For more information, see the topic Running Commands on p. 38.](#)
- Within a program block, each line should not exceed 251 bytes (although syntax generated by those lines can be longer).
- With the SPSS Statistics Batch Facility (available only with SPSS Statistics Server), use the `-i` switch when submitting command files that contain program blocks. All command syntax (not just the program blocks) in the file must adhere to interactive syntax rules.

Within a program block, the programming language is in control, and the syntax rules for that programming language apply. Command syntax generated from within program blocks must always follow interactive syntax rules. For most practical purposes this means command strings you build in a programming block must contain a period (.) at the end of each command.

### Scope and Limitations

- Programmatic variables created in a program block cannot be used outside of program blocks.
- Program blocks cannot be contained within `DEFINE-!ENDDEFINE` macro definitions.

- Program blocks can be contained in command syntax files run via the `INSERT` command, with the default `SYNTAX=INTERACTIVE` setting.
- Program blocks cannot be contained within command syntax files run via the `INCLUDE` command.

### ***Using External Programming Languages***

Use of the IBM® SPSS® Statistics Programmability Extension requires an Integration Plug-In for an external language. Integration Plug-Ins are available for the Python and R programming languages. For information, see *How to Get Integration Plug-Ins*, available from *Core System>Frequently Asked Questions* in the Help system. Documentation for the plug-ins is available from the topics *Integration Plug-In for Python* and *Integration Plug-In for R* in the Help system once the plug-ins are installed. It is also available in PDF from *Help>Programmability* once the plug-ins are installed.

# BOOTSTRAP

BOOTSTRAP is available in the Bootstrapping option.

```
BOOTSTRAP
  [/SAMPLING METHOD={SIMPLE**          }]
    {STRATIFIED(STRATA=varlist)  }
    {RESIDUAL({RESIDUALS=varlist})}
      {PREDICTED=varlist}
    {WILD({RESIDUALS=varlist})  }
      {PREDICTED=varlist}
  [/VARIABLES [TARGET=varlist] [INPUT=varlist]]
  [/CRITERIA [CILEVEL={95** } [CITYPE={PERCENTILE**}]]]
    {value}          {BCA      }
    [NSAMPLES={1000**}]
      {int      }
  [/MISSING [USERMISSING={EXCLUDE**}]]
    {INCLUDE  }.
```

\*\* Default if the subcommand or keyword is omitted.

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 18

- Command introduced.

## **Example**

```
BOOTSTRAP.
```

## **Overview**

Bootstrapping is a method for deriving robust estimates of standard errors and confidence intervals for estimates such as the mean, median, proportion, odds ratio, correlation coefficient or regression coefficient. It may also be used for constructing hypothesis tests. Bootstrapping is most useful as an alternative to parametric estimates when the assumptions of those methods are in doubt (as in the case of regression models with heteroscedastic residuals fit to small samples), or where parametric inference is impossible or requires very complicated formulas for the calculation of standard errors (as in the case of computing confidence intervals for the median, quartiles, and other percentiles).

The `BOOTSTRAP` command signals the beginning of temporary bootstrap samples that are in effect only for the next procedure.



### **Options**

**Resampling method.** Simple, stratified, and residuals bootstrap resampling are supported. You can also specify the number of bootstrap samples to take.

**Pooling method.** Choose between percentile and BCa methods for computing confidence intervals. You can also specify the confidence level.

### **Basic Specification**

The basic specification is the `BOOTSTRAP` command.

By default, `BOOTSTRAP` draws 1000 samples using simple bootstrap resampling. When the procedure following `BOOTSTRAP` is run, the pooling algorithm produces 95% confidence intervals using the percentile method. Since no variables have been specified, no records are excluded from resampling.

### **Syntax Rules**

- All subcommands are optional.
- Subcommands may be specified in any order.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not allowed.
- Any split variable defined on the `SPLIT FILE` command may not be used on the `BOOTSTRAP` command.

### **Limitations**

- `BOOTSTRAP` does not work with multiply imputed datasets. If there is an *Imputation\_* variable in the dataset, running `BOOTSTRAP` will cause an error.
- `BOOTSTRAP` should not be used in conjunction with the `N OF CASES` command.

## **Examples**

### **Simple Resampling; Maintaining a Consistent Case Basis**

```
BOOTSTRAP.  
DESCRIPTIVES VARIABLES=var1 var2 var3  
  /MISSING=VARIABLE.
```

- The `BOOTSTRAP` command requests 1000 bootstrap samples.

- No variables are specified on the `BOOTSTRAP` command, so no records are deleted from the resampling. This allows the `DESCRIPTIVES` procedure to use variablewise deletion of missing values on the full set of records; however, the case basis will be inconsistent across bootstrap resamples, and inferences made from the results would be questionable.

```
BOOTSTRAP
/VARIABLES ANALYSIS(INPUT=var1 var2 var3).
DESCRIPTIVES VARIABLES=var1 var2 var3
/STATISTICS MEAN STDDEV MIN MAX
/MISSING=VARIABLE.
```

- This is the same as the previous analysis, but variables *var1*, *var2*, and *var3* are used to determine the case basis for resampling. Records with missing values on any of these variables are deleted from the analysis.
- The `DESCRIPTIVES` procedure following `BOOTSTRAP` is run on the bootstrap samples.
- The `STATISTICS` subcommand produces the mean, standard deviation, minimum, and maximum for variables *var1*, *var2*, and *var3* on the original data. Additionally, pooled statistics are produced for the mean and standard deviation.
- Even though the `MISSING` subcommand specifies variablewise deletion of missing values, the listwise deletion performed by `BOOTSTRAP` is what determines the case basis. In effect, the `MISSING` specification on `DESCRIPTIVES` is irrelevant here.

### **Simple Resampling; Replicating Results**

```
PRESERVE.
SET RNG=MT MTINDEX=592004.
SHOW RNG.
BOOTSTRAP
/SAMPLING METHOD=SIMPLE
/VARIABLES TARGET=prevexp
/CRITERIA CILEVEL=95 CITYPE=BCA NSAMPLES=1000
/MISSING USERMISSING=EXCLUDE.
EXAMINE VARIABLES=prevexp
/PLOT NONE
/STATISTICS DESCRIPTIVES
/CINTERVAL 95
/MISSING LISTWISE
/NOTOTAL.
RESTORE.
```

- The `PRESERVE` and `RESTORE` commands “remember” the current state of the random number generator and restore the system to that state after bootstrapping is over.
- The `SET` command sets the random number generator to the Mersenne Twister and the index to 592004, so that the bootstrapping results can be replicated exactly. The `SHOW` command displays the index in the output for reference.
- The `BOOTSTRAP` command requests 1000 bootstrap samples using simple resampling.
- The `VARIABLES` subcommand specifies that the variable *prevexp* is used to determine the case basis for resampling. Records with missing values on this variable are deleted from the analysis.
- The `CRITERIA` subcommand, in addition to requesting the number of bootstrap samples, requests bias-corrected and accelerated bootstrap confidence intervals instead of the default percentile intervals.

- The EXAMINE procedure following BOOTSTRAP is run on each of the bootstrap samples.
- The PLOT subcommand turns off plot output.
- All other options are set to their default values.

### **Stratified Resampling**

```
BOOTSTRAP
  /VARIABLES SAMPLING (STRATA=strataVar)
              ANALYSIS (INPUTS=var1) .
DESCRIPTIVES var1.
```

- The BOOTSTRAP command requests 1000 bootstrap samples stratified by *strataVar*.
- Variables *var1* and *strataVar* are used to determine the case basis for resampling. Records with missing values on these variables are deleted from the analysis.
- The DESCRIPTIVES procedure following BOOTSTRAP is run on the bootstrap samples, and produces the mean, standard deviation, minimum, and maximum for the variable *var1* on the original data. Additionally, pooled statistics are produced for the mean and standard deviation.

### **Residual Resampling**

```
COMPUTE diff=salary-salbegin.
EXECUTE.

UNIANOVA diff BY gender jobcat minority WITH jobtime prevexp
  /METHOD=SSTYPE(3)
  /INTERCEPT=INCLUDE
  /SAVE=RESID
  /CRITERIA=ALPHA(0.05)
  /DESIGN=gender jobcat minority jobtime prevexp.
```

- The COMPUTE command creates the variable containing the change in salary.
- The UNIANOVA command fits a model for *diff* using *gender*, *jobcat*, and *minority* as factors and *jobtime* and *prevexp* as covariates.
- Unstandardized residuals are saved to a new variable in the dataset named *RES\_1* by default.
- All other options are set to their default values.

```
PRESERVE.
SET RNG=MT MTINDEX=9191972.
SHOW RNG.
BOOTSTRAP
  /SAMPLING METHOD=WILD(RESIDUALS=RES_1)
  /VARIABLES TARGET=diff INPUT=gender jobcat minority jobtime prevexp
  /CRITERIA CILEVEL=95 CITYPE=PERCENTILE NSAMPLES=1000
  /MISSING USERMISSING=EXCLUDE.
UNIANOVA diff BY gender jobcat minority WITH jobtime prevexp
  /METHOD=SSTYPE(3)
  /INTERCEPT=INCLUDE
  /PRINT=PARAMETER
  /CRITERIA=ALPHA(.05)
  /DESIGN=gender jobcat minority jobtime prevexp.
RESTORE.
```

- The `PRESERVE` and `RESTORE` commands “remember” the current state of the random number generator and restore the system to that state after bootstrapping is over.
- The `SET` command sets the random number generator to the Mersenne Twister and the index to 9191972, so that the bootstrapping results can be replicated exactly. The `SHOW` command displays the index in the output for reference.
- The `BOOTSTRAP` command requests 1000 bootstrap samples using wild sampling and `RES_1` as the variable containing the residuals.
- The `VARIABLES` subcommand specifies that the *diff* is the target variable in the linear model; it and the variables *gender*, *jobcat*, *minority*, *jobtime*, and *prevexp* are used to determine the case basis for resampling. Records with missing values on these variables are deleted from the analysis.
- The `CRITERIA` subcommand, in addition to requesting the number of bootstrap samples, requests bias-corrected and accelerated bootstrap confidence intervals instead of the default percentile intervals.
- The `UNIANOVA` procedure following `BOOTSTRAP` is run on each of the bootstrap samples and produces parameter estimates for the original data. Additionally, pooled statistics are produced for the model coefficients.

## ***SAMPLING Subcommand***

The `SAMPLING` subcommand is used to specify the sampling method and any associated variables.

- If `SAMPLING` is not specified, the procedure performs simple bootstrap resampling..

### **SIMPLE**

*Simple resampling.* This performs case resampling with replacement from the original dataset. This is the default.

### **STRATIFIED(STRATA = varlist)**

*Stratified resampling.* Specify one or more variables that define strata within the dataset. This performs case resampling with replacement from the original dataset, *within* the strata defined by the cross-classification of strata variables, preserving the size of each stratum. Stratified bootstrap sampling can be useful when units within strata are relatively homogeneous while units across strata are very different.

### **RESIDUAL(RESIDUALS=varlist | PREDICTED=varlist)**

*Residual resampling.* Specify one or more variables containing residuals from fitting a model to the data. The model that produced the residuals should ideally be the same model that follows `BOOTSTRAP`. A residual sample is drawn by replacing each target variable value with that case’s predicted value plus a residual sampled from the entire original set of residuals.

Specify `PREDICTED` as an alternative to `RESIDUALS` when the model residuals are not immediately available but the predicted values are. Specify one or more variables containing predicted values from fitting a model to the data.

If `RESIDUAL` is specified, the `TARGET` keyword is required and the variables specified on `RESIDUAL` should be the residuals (or predicted values) for, and match the order of, the variables specified on `TARGET`.

### **WILD(RESIDUALS=varlist | PREDICTED=varlist)**

*Wild bootstrap resampling.* Specify one or more variables containing residuals from fitting a model to the data. The model that produced the residuals should ideally be the same model that follows `BOOTSTRAP`. A wild sample is drawn by replacing each target variable value with that case's predicted value plus either the case's residual or the negative of the case's residual.

Specify `PREDICTED` as an alternative to `RESIDUALS` when the model residuals are not immediately available but the predicted values are. Specify one or more variables containing predicted values from fitting a model to the data.

If `WILD` is specified, the `TARGET` keyword is required and the variables specified on `WILD` should be the residuals (or predicted values) for, and match the order of, the variables specified on `TARGET`.

## **VARIABLES Subcommand**

The `VARIABLES` subcommand is used to specify the target and inputs.

- If `VARIABLES` is not specified, the procedure performs bootstrap resampling on all the records in the dataset. `TARGET` is required when performing residual resampling, but these specifications are otherwise *technically* optional. However, these variables are used to determine the case basis for bootstrap resampling, so it is important to specify these variables when there are missing values in the data.

### **TARGET=varlist**

*Target variables.* Specify one or more variables that will be used as targets (responses, dependent variables) in the procedure following `BOOTSTRAP`.

### **INPUT=varlist**

*Input variables.* Specify one or more variables that will be used as inputs (factors, covariates) in the procedure following `BOOTSTRAP`.

## **CRITERIA Subcommand**

The `CRITERIA` subcommand controls pooling options and the number of bootstrap samples to take.

### **CILEVEL = number**

*Confidence interval level.* Specify a number greater than or equal to 0, and less than 100. The default value is 95. Note that bootstrapping can only support intervals up to confidence level  $100 * (1 - 2 / (NSAMPLES + 1))$ .

### **CITYPE = PERCENTILE | BCA**

*Confidence interval type.* Specify `PERCENTILE` for percentile intervals or `BCA` for BCa intervals. The default value is `PERCENTILE`.

### **NSAMPLES = integer**

*Number of bootstrap samples.* Specify a positive integer. The default value is 1000.

## ***MISSING Subcommand***

The `MISSING` subcommand is used to control whether user-missing values for categorical variables are treated as valid values. By default, user-missing values for categorical variables are treated as invalid. The setting used here should be the same as that used on the procedure following the `BOOTSTRAP` command.

- Cases with invalid values are deleted listwise.
- The `MISSING` subcommand defines categorical variables as variables with measurement level set at Ordinal or Nominal in the data dictionary. Use the `VARIABLE LEVEL` command to change a variable's measurement level.
- User-missing values for continuous variables are always treated as invalid.
- System-missing values for any variables are always treated as invalid.

### **USERMISSING=EXCLUDE**

*User-missing values for categorical variables are treated as invalid.  
This is the default.*

### **USERMISSING=INCLUDE**

*User-missing values for categorical variables are treated as valid values.*

# BREAK

BREAK

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## Overview

BREAK controls looping that cannot be fully controlled with IF clauses. Generally, BREAK is used within a DO IF-END IF structure. The expression on the DO IF command specifies the condition in which BREAK is executed.

### Basic Specification

- The only specification is the keyword BREAK. There are no additional specifications.
- BREAK must be specified within a loop structure. Otherwise, an error results.

### Operations

- A BREAK command inside a loop structure but not inside a DO IF-END IF structure terminates the first iteration of the loop for all cases, since no conditions for BREAK are specified.
- A BREAK command within an inner loop terminates only iterations in that structure, not in any outer loop structures.

## Examples

```
VECTOR          #X(10).  
LOOP            #I = 1 TO #NREC.  
+ DATA LIST   NOTABLE/ #X1 TO #X10 1-20.  
+ LOOP         #J = 1 TO 10.  
+   DO IF      SYSMIS(#X(#J)).  
+     BREAK.  
+   END IF.  
+   COMPUTE    X = #X(#J).  
+   END CASE.  
+ END LOOP.  
END LOOP.
```

- The inner loop terminates when there is a system-missing value for any of the variables #X1 to #X10.
- The outer loop continues until all records are read.

# CACHE

CACHE.

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

Although the virtual active file can vastly reduce the amount of temporary disk space required, the absence of a temporary copy of the “active” file means that the original data source has to be reread for each procedure. For data tables read from a database source, this means that the SQL query that reads the information from the database must be reexecuted for any command or procedure that needs to read the data. Since virtually all statistical analysis procedures and charting procedures need to read the data, the SQL query is reexecuted for each procedure that you run, which can result in a significant increase in processing time if you run a large number of procedures.

If you have sufficient disk space on the computer performing the analysis (either your local computer or a remote server), you can eliminate multiple SQL queries and improve processing time by creating a data cache of the active file with the `CACHE` command. The `CACHE` command copies all of the data to a temporary disk file the next time the data are passed to run a procedure. If you want the cache written immediately, use the `EXECUTE` command after the `CACHE` command.

- The only specification is the command name `CACHE`.
- A cache file will not be written during a procedure that uses temporary variables.
- A cache file will not be written if the data are already in a temporary disk file and that file has not been modified since it was written.

## **Example**

```
CACHE.  
TEMPORARY.  
RECODE alcohol(0 thru .04 = 'sober') (.04 thru .08 = 'tipsy')  
      (else = 'drunk') into state.  
FREQUENCIES var=state.  
GRAPH...
```

No cache file will be written during the `FREQUENCIES` procedure. It will be written during the `GRAPH` procedure.



# CASEPLOT

```
CASEPLOT VARIABLES=varlist

[/DIFF={1}]
      {n}

[/SDIFF={1}]
      {n}

[/PERIOD=n]

[/ {NOLOG**}]
      {LN      }

[/ID=varname]

[/MARK={varname      }]
      {date specification}

[/SPLIT {UNIFORM**}]
      {SCALE      }

[/APPLY [= 'model name']]
```

*For plots with one variable:*

```
[/FORMAT={ {NOFILL**} } [ {NOREFERENCE** } ] ]
      {LEFT      }      {REFERENCE[ (value) ] }
```

*For plots with multiple variables:*

```
[/FORMAT={NOJOIN**}]
      {JOIN      }
      {HILO      }
```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- For plots with one variable, new option to specify a value with the REFERENCE keyword on the FORMAT subcommand.

## **Example**

```
CASEPLOT VARIABLES = TICKETS
  /LN
  /DIFF
  /SDIFF
  /PERIOD=12
  /FORMAT=REFERENCE
  /MARK=Y 55 M 6.
```

## Overview

CASEPLOT produces a plot of one or more time series or sequence variables. You can request natural log and differencing transformations to produce plots of transformed variables. Several plot formats are available.

### Options

**Modifying the Variables.** You can request a natural log transformation of the variable using the `LN` subcommand and seasonal and nonseasonal differencing to any degree using the `SDIFF` and `DIFF` subcommands. With seasonal differencing, you can also specify the periodicity on the `PERIOD` subcommand.

**Plot Format.** With the `FORMAT` subcommand, you can fill in the area on one side of the plotted values on plots with one variable. You can also plot a reference line indicating the variable mean. For plots with two or more variables, you can specify whether you want to join the values for each case with a horizontal line. With the `ID` subcommand, you can label the vertical axis with the values of a specified variable. You can mark the onset of an intervention variable on the plot with the `MARK` subcommand.

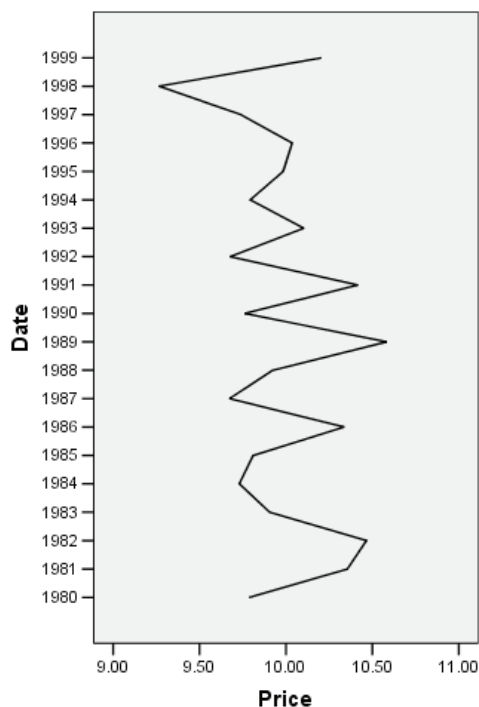
**Split-File Processing.** You can control how to plot data that have been divided into subgroups by a `SPLIT FILE` command using the `SPLIT` subcommand.

### Basic Specification

The basic specification is one or more variable names.

- If the `DATE` command has been specified, the vertical axis is labeled with the `DATE_` variable at periodic intervals. Otherwise, sequence numbers are used. The horizontal axis is labeled with the value scale determined by the plotted variables.

Figure 23-1  
CASEPLOT with DATE variable



### **Subcommand Order**

- Subcommands can be specified in any order.

### **Syntax Rules**

- VARIABLES can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each one is executed.

### **Operations**

- Subcommand specifications apply to all variables named on the CASEPLOT command.
- If the LN subcommand is specified, any differencing requested on that CASEPLOT command is done on the log-transformed variables.
- Split-file information is displayed as part of the subtitle, and transformation information is displayed as part of the footnote.

### **Limitations**

- A maximum of one VARIABLES subcommand. There is no limit on the number of variables named on the list.

## Examples

```
CASEPLOT VARIABLES = TICKETS
  /LN
  /DIFF
  /SDIFF
  /PERIOD=12
  /FORMAT=REFERENCE
  /MARK=Y 55 M 6.
```

- This example produces a plot of *TICKETS* after a natural log transformation, differencing, and seasonal differencing have been applied.
- LN transforms the data using the natural logarithm (base  $e$ ) of the variable.
- DIFF differences the variable once.
- SDIFF and PERIOD apply one degree of seasonal differencing with a periodicity of 12.
- FORMAT=REFERENCE adds a reference line at the variable mean.
- MARK provides a marker on the plot at June, 1955. The marker is displayed as a horizontal reference line.

## VARIABLES Subcommand

VARIABLES specifies the names of the variables to be plotted and is the only required subcommand.

## DIFF Subcommand

DIFF specifies the degree of differencing used to convert a nonstationary variable to a stationary one with a constant mean and variance before plotting.

- You can specify any positive integer on DIFF.
- If DIFF is specified without a value, the default is 1.
- The number of values displayed decreases by 1 for each degree of differencing.

### Example

```
CASEPLOT VARIABLES = TICKETS
  /DIFF=2.
```

- In this example, *TICKETS* is differenced twice before plotting.

## SDIFF Subcommand

If the variable exhibits a seasonal or periodic pattern, you can use the SDIFF subcommand to seasonally difference a variable before plotting.

- The specification on SDIFF indicates the degree of seasonal differencing and can be any positive integer.
- If SDIFF is specified without a value, the degree of seasonal differencing defaults to 1.

- The number of seasons displayed decreases by 1 for each degree of seasonal differencing.
- The length of the period used by `SDIFF` is specified on the `PERIOD` subcommand. If the `PERIOD` subcommand is not specified, the periodicity established on the `TSET` or `DATE` command is used (see the `PERIOD` subcommand below).

## ***PERIOD Subcommand***

`PERIOD` indicates the length of the period to be used by the `SDIFF` subcommand.

- The specification on `PERIOD` indicates how many observations are in one period or season and can be any positive integer.
- `PERIOD` is ignored if it is used without the `SDIFF` subcommand.
- If `PERIOD` is not specified, the periodicity established on `TSET` `PERIOD` is in effect. If `TSET` `PERIOD` is not specified either, the periodicity established on the `DATE` command is used. If periodicity is not established anywhere, the `SDIFF` subcommand will not be executed.

### ***Example***

```
CASEPLOT VARIABLES = TICKETS
  /SDIFF=1
  /PERIOD=12 .
```

- This command applies one degree of seasonal differencing with 12 observations per season to *TICKETS* before plotting.

## ***LN and NOLOG Subcommands***

`LN` transforms the data using the natural logarithm (base  $e$ ) of the variable and is used to remove varying amplitude over time. `NOLOG` indicates that the data should not be log transformed. `NOLOG` is the default.

- If you specify `LN` on `CASEPLOT`, any differencing requested on that command will be done on the log-transformed variable.
- There are no additional specifications on `LN` or `NOLOG`.
- Only the last `LN` or `NOLOG` subcommand on a `CASEPLOT` command is executed.
- If a natural log transformation is requested, any value less than or equal to zero is set to system-missing.
- `NOLOG` is generally used with an `APPLY` subcommand to turn off a previous `LN` specification.

### ***Example***

```
CASEPLOT VARIABLES = TICKETS
  /LN .
```

- In this example, *TICKETS* is transformed using the natural logarithm before plotting.

## ***ID Subcommand***

ID names a variable whose values will be used as the left-axis labels.

- The only specification on ID is a variable name. If you have a variable named *ID* in your active dataset, the equals sign after the subcommand is required.
- ID overrides the specification on TSET ID.
- If ID or TSET ID is not specified, the left vertical axis is labeled with the *DATE\_* variable created by the DATE command. If the *DATE\_* variable has not been created, the observation or sequence number is used as the label.

### ***Example***

```
CASEPLOT VARIABLES = VARA  
/ID=VARB.
```

- In this example, the values of the variable *VARB* will be used to label the left axis of the plot of *VARA*.

## ***FORMAT Subcommand***

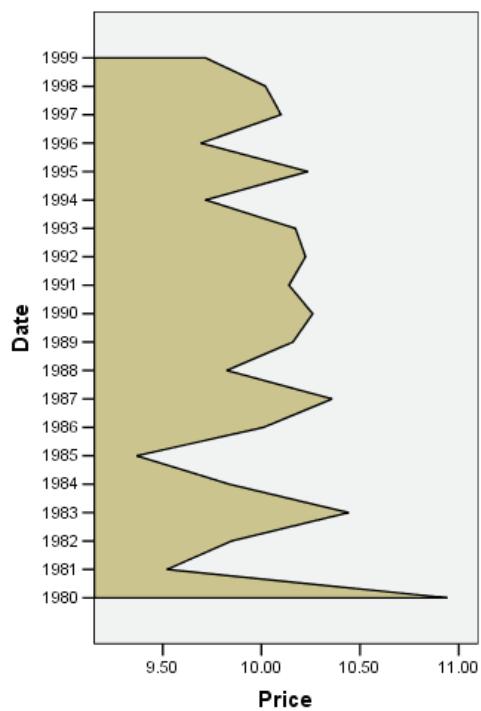
FORMAT controls the plot format.

- The specification on FORMAT is one of the keywords listed below.
- The keywords NOFILL, LEFT, NOREFERENCE, and REFERENCE apply to plots with one variable. NOFILL and LEFT are alternatives and indicate how the plot is filled. NOREFERENCE and REFERENCE are alternatives and specify whether a reference line is displayed. One keyword *from each set* can be specified. NOFILL and NOREFERENCE are the defaults.
- The keywords JOIN, NOJOIN, and HILO apply to plots with multiple variables and are alternatives. NOJOIN is the default. Only one keyword can be specified on a FORMAT subcommand for plots with two variables.

The following formats are available for plots of one variable:

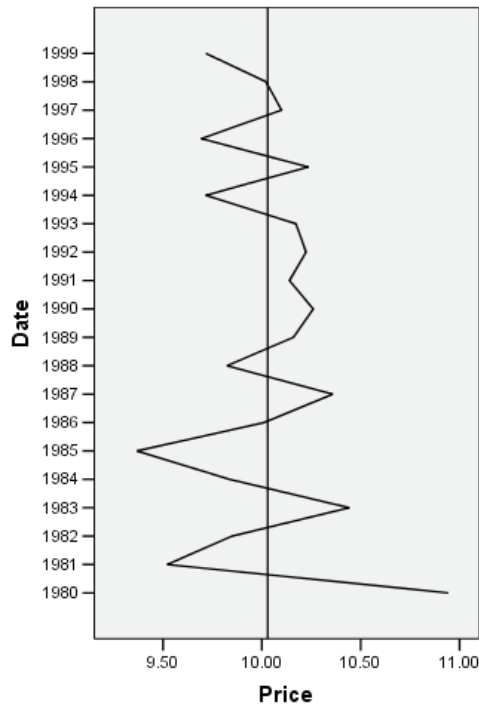
<b>NOFILL</b>	<i>Plot only the values for the variable with no fill.</i> NOFILL produces a plot with no fill to the left or right of the plotted values. This is the default format when one variable is specified.
<b>LEFT</b>	<i>Plot the values for the variable and fill in the area to the left.</i> If the plotted variable has missing or negative values, the keyword LEFT is ignored and the default NOFILL is used instead.

Figure 23-2  
FORMAT=LEFT



- NOREFERENCE** *Do not plot a reference line. This is the default when one variable is specified.*
- REFERENCE(value)** *Plot a reference line at the specified value or at the variable mean if no value is specified. A fill chart is displayed as an area chart with a reference line and a non-fill chart is displayed as a line chart with a reference line.*

Figure 23-3  
 FORMAT=REFERENCE



The following formats are available for plots of multiple variables:

- |               |  |
|---------------|--|
| <b>NOJOIN</b> | <i>Plot the values of each variable named.</i> Different colors or line patterns are used for multiple variables. Multiple occurrences of the same value for a single observation are plotted using a dollar sign (\$). This is the default format for plots of multiple variables.  |
| <b>JOIN</b>   | <i>Plot the values of each variable and join the values for each case.</i> Values are plotted as described for <b>NOJOIN</b> , and the values for each case are joined together by a line.   |
| <b>HILO</b>   | <i>Plot the highest and lowest values across variables for each case and join the two values together.</i> The high and low values are plotted as a pair of vertical bars and are joined with a dashed line. <b>HILO</b> is ignored if more than three variables are specified, and the default <b>NOJOIN</b> is used instead. |

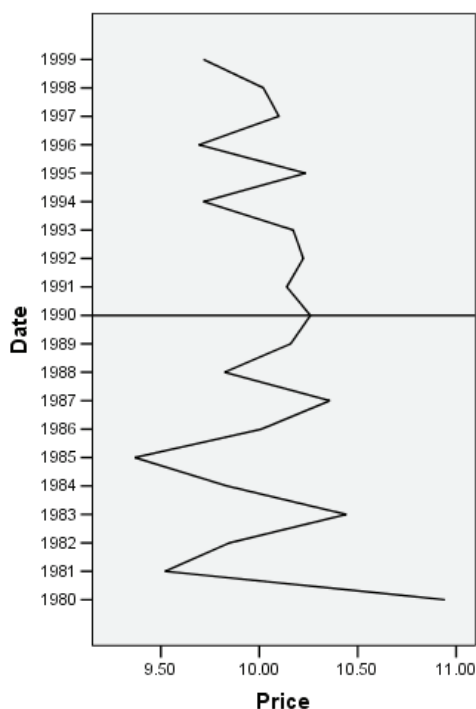
## **MARK Subcommand**

Use **MARK** to indicate the onset of an intervention variable.

- The onset date is indicated by a horizontal reference line.
- The specification on **MARK** can be either a variable name or an onset date if the *DATE\_* variable exists.
- If a variable is named, the reference line indicates where the values of that variable change.
- A date specification follows the same format as the **DATE** command—that is, a keyword followed by a value. For example, the specification for June, 1955, is `Y 1955 M 6` (or `Y 55 M 6` if only the last two digits of the year are used on **DATE**).



Figure 23-4  
MARK Y=1990



## ***SPLIT Subcommand***

**SPLIT** specifies how to plot data that have been divided into subgroups by a **SPLIT FILE** command. The specification on **SPLIT** is either **SCALE** or **UNIFORM**.

- If **FORMAT=REFERENCE** is specified when **SPLIT=SCALE**, the reference line is placed at the mean of the subgroup. If **FORMAT=REFERENCE** is specified when **SPLIT=UNIFORM**, the reference line is placed at the overall mean.

**UNIFORM**                      *Uniform scale.* The horizontal axis is scaled according to the values of the entire dataset. This is the default if **SPLIT** is not specified.

**SCALE**                              *Individual scale.* The horizontal axis is scaled according to the values of each individual subgroup.

### ***Example***

```
SPLIT FILE BY REGION.
CASEPLOT VARIABLES = TICKETS / SPLIT=SCALE.
```

- This example produces one plot for each *REGION* subgroup.
- The horizontal axis for each plot is scaled according to the values of *TICKETS* for each particular region.

## ***APPLY Subcommand***

APPLY allows you to produce a caseplot using previously defined specifications without having to repeat the CASEPLOT subcommands.

- The only specification on APPLY is the name of a previous model in quotes. If a model name is not specified, the specifications from the previous CASEPLOT command are used.
- If no variables are specified, the variables that were specified for the original plot are used.
- To change one or more plot specifications, specify the subcommands of only those portions you want to change after the APPLY subcommand.
- To plot different variables, enter new variable names before or after the APPLY subcommand.

### ***Example***

```
CASEPLOT VARIABLES = TICKETS
  /LN
  /DIFF=1
  /SDIFF=1
  /PER=12.
CASEPLOT VARIABLES = ROUNDTRP
  /APPLY.
CASEPLOT APPLY
  /NOLOG.
```

- The first command produces a plot of *TICKETS* after a natural log transformation, differencing, and seasonal differencing.
- The second command plots *ROUNDTRP* using the same transformations specified for *TICKETS*.
- The third command produces a plot of *ROUNDTRP* but this time without any natural log transformation. The variable is still differenced once and seasonally differenced with a periodicity of 12.

# CASESTOVARS

```
CASESTOVARS
[/ID = varlist]
[/FIXED = varlist]
[/AUTOFIX = {YES**}
             {NO   }]
[/VIND [ROOT = rootname]]
[/COUNT = new variable ["label"]]
[/RENAME varname=rootname varname=rootname ...]
[/SEPARATOR = {"."   }
              {"string"}]
[/INDEX = varlist]
[/GROUPBY = {VARIABLE**}
             {INDEX   }]}
[/DROP = varlist]
```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
CASESTOVARS /ID idvar /INDEX var1.
```

## **Overview**

A **variable** contains information that you want to analyze, such as a measurement or a test score. A **case** is an observation, such as an individual or an institution.

In a **simple data file**, each variable is a single **column** in your data, and each case is a single **row** in your data. So, if you were recording the score on a test for all students in a class, the scores would appear in only one column and there would be only one row for each student.

**Complex data files** store data in more than one column or row. For example, in a complex data file, information about a case could be stored in more than one row. So, if you were recording monthly test scores for all students in a class, there would be multiple rows for each student—one for each month.

CASESTOVARS restructures complex data that has multiple rows for a case. You can use it to restructure data in which repeated measurements of a single case were recorded in multiple rows (row groups) into a new data file in which each case appears as separate variables (variable groups) in a single row. It replaces the active dataset.

### ***Options***

**Automatic classification of fixed variables.** The values of **fixed variables** do not vary within a row group. You can use the `AUTOFIX` subcommand to let the procedure determine which variables are fixed and which variables are to become variable groups in the new data file.

**Naming new variables.** You can use the `RENAME`, `SEPARATOR`, and `INDEX` subcommands to control the names for the new variables.

**Ordering new variables.** You can use the `GROUPBY` subcommand to specify how to order the new variables in the new data file.

**Creating indicator variables.** You can use the `VIND` subcommand to create indicator variables. An **indicator variable** indicates the presence or absence of a value for a case. An indicator variable has the value of 1 if the case has a value; otherwise, it is 0.

**Creating a count variable.** You can use the `COUNT` subcommand to create a count variable that contains the number of rows in the original data that were used to create a row in the new data file.

**Variable selection.** You can use the `DROP` subcommand to specify which variables from the original data file are dropped from the new data file.

### ***Basic specification***

The basic specification is simply the command keyword.

- If split-file processing is in effect, the basic specification creates a row in the new data file for each combination of values of the `SPLIT FILE` variables. If split-file processing is not in effect, the basic specification results in a new data file with one row.
- Because the basic specification can create quite a few new columns in the new data file, the use of an `ID` subcommand to identify groups of cases is recommended.

### ***Subcommand order***

Subcommands can be specified in any order.

### ***Syntax rules***

Each subcommand can be specified only once.

### ***Operations***

- **Original row order.** `CASESTOVARS` assumes that the original data are sorted by `SPLIT` and `ID` variables.
- **Identifying row groups in the original file.** A **row group** consists of rows in the original data that share the same values of variables listed on the `ID` subcommand. Row groups are consolidated into a single row in the new data file. Each time a new combination of `ID` values is encountered, a new row is created.

- **Split-file processing and row groups.** If split-file processing is in effect, the split variables are automatically used to identify row groups (they are treated as though they appeared first on the `ID` subcommand). Split-file processing remains in effect in the new data file unless a variable that is used to split the file is named on the `DROP` subcommand.
- **New variable groups.** A **variable group** is a group of related columns in the new data file that is created from a variable in the original data. Each variable group contains a variable for each index value or combination of index values encountered.
- **Candidate variables.** A variable in the original data is a candidate to become a variable group in the new data file if it is not used on the `SPLIT` command or the `ID`, `FIXED`, or `DROP` subcommands and its values vary within the row group. Variables named on the `SPLIT`, `ID`, and `FIXED` subcommands are assumed to not vary within the row group and are simply copied into the new data file.
- **New variable names.** The names of the variables in a new group are constructed by the procedure. It uses the rootname specified on the `RENAME` subcommand and the string named on the `SEPARATOR` subcommand.
- **New variable formats.** With the exception of names and labels, the dictionary information for all of the new variables in a group (for example, value labels and format) is taken from the variable in the original data.
- **New variable order.** New variables are created in the order specified by the `GROUPBY` subcommand.
- **Weighted files.** The `WEIGHT` command does not affect the results of `CASESTOVARS`. If the original data are weighted, the new data file will be weighted unless the variable that is used as the weight is dropped from the new data file.
- **Selected cases.** The `FILTER` and `USE` commands do not affect the results of `CASESTOVARS`. It processes all cases.

### Limitations

The `TEMPORARY` command cannot be in effect when `CASESTOVARS` is executed.

## Examples

The following is the `LIST` output for a data file in which repeated measurements for the same case are stored on separate rows in a single variable.

insure	caseid	month	bps	bpd
BCBS	1	1	160	100
BCBS	2	1	120	70
BCBS	2	2	130	86
Prucare	1	1	160	94
Prucare	1	2	200	105
Prucare	1	3	180	105
Pruecare	2	1	135	90

The commands:

```
SPLIT FILE BY insure.
CASESTOVARS
```

```

/ID=caseid
/INDEX=month.

```

create a new variable group for *bps* and a new group for *bpd*. The LIST output for the new active dataset is as follows:

- The row groups in the original data are identified by *insure* and *caseid*.
- There are four row groups—one for each combination of the values in *insure* and *caseid*.
- The command creates four rows in the new data file, one for each row group.
- The candidate variables from the original file are *bps* and *bpd*. They vary within the row group, so they will become variable groups in the new data file.
- The command creates two new variable groups—one for *bps* and one for *bpd*.
- Each variable group contains three new variables—one for each unique value of the index variable *month*.

## ***ID subcommand***

The ID subcommand specifies variables that identify the rows from the original data that should be grouped together in the new data file.

- If the ID subcommand is omitted, only SPLIT FILE variables (if any) will be used to group rows in the original data and to identify rows in the new data file.
- CASESTOVARS expects the data to be sorted by SPLIT FILE variables and then by ID variables. If split-file processing is in effect, the original data should be sorted on the split variables in the order given on the SPLIT FILE command and then on the ID variables in the order in which they appear in the ID subcommand.
- A variable may appear on both the SPLIT FILE command and the ID subcommand.
- Variables listed on the SPLIT FILE command and on the ID subcommand are copied into the new data file with their original values and dictionary information unless they are dropped with the DROP subcommand.
- Variables listed on the ID subcommand may not appear on the FIXED or INDEX subcommands.
- Rows in the original data for which any ID variable has the system-missing value or is blank are not included in the new data file, and a warning message is displayed.
- ID variables are not candidates to become a variable group in the new data file.

## ***INDEX subcommand***

In the original data, a variable appears in a single column. In the new data file, that variable will appear in multiple new columns. The INDEX subcommand names the variables in the original data that should be used to create the new columns. INDEX variables are also used to name the new columns.

Optionally, with the GROUPBY subcommand, INDEX variables can be used to determine the order of the new columns, and, with the VIND subcommand, INDEX variables can be used to create indicator variables.

- String variables can be used as index variables. They cannot contain blank values for rows in the original data that qualify for inclusion in the new data file.
- Numeric variables can be used as index variables. They must contain only non-negative integer values and cannot have system-missing or blank values.
- Within each row group in the original file, each row must have a different combination of values of the index variables.
- If the `INDEX` subcommand is not used, the index starts with 1 within each row group and increments each time a new value is encountered in the original variable.
- Variables listed on the `INDEX` subcommand may not appear on the `ID`, `FIXED`, or `DROP` subcommands.
- Index variables are not candidates to become a variable group in the new data file.

## ***VIND subcommand***

The `VIND` subcommand creates indicator variables in the new data file. An indicator variable indicates the presence or absence of a value for a case. An indicator variable has the value of 1 if the case has a value; otherwise, it is 0.

- One new indicator variable is created for each unique value of the variables specified on the `INDEX` subcommand.
- If the `INDEX` subcommand is not used, an indicator variable is created each time a new value is encountered within a row group.
- An optional rootname can be specified after the `ROOT` keyword on the subcommand. The default rootname is *ind*.
- The format for the new indicator variables is F1.0.

### ***Example***

If the original variables are:

```
insure      caseid      month      bps      bpd
```

and the data are as shown in the first example, the commands:

```
SPLIT FILE BY insure.
CASESTOVARs
  /ID=caseid
  /INDEX=month
  /VIND
  /DROP=caseid bpd.
```

create a new file with the following data:

- The command created three new indicator variables—one for each unique value of the index variable *month*.

## ***COUNT subcommand***

CASESTOVARS consolidates row groups in the original data into a single row in the new data file. The COUNT subcommand creates a new variable that contains the number of rows in the original data that were used to generate the row in the new data file.

- One new variable is named on the COUNT subcommand. It must have a unique name.
- The label for the new variable is optional and, if specified, must be delimited by single or double quotes.
- The format of the new count variable is F4.0.

### ***Example***

If the original data are as shown in the first example, the commands:

```
SPLIT FILE BY insure.  
CASESTOVARS  
  /ID=caseid  
  /COUNT=countvar  
  /DROP=insure month bpd.
```

create a new file with the following data:

- The command created a count variable, *countvar*, which contains the number of rows in the original data that were used to generate the current row.

## ***FIXED subcommand***

The FIXED subcommand names the variables that should be copied from the original data to the new data file.

- CASESTOVARS assumes that variables named on the FIXED subcommand do not vary within row groups in the original data. If they vary, a warning message is generated and the command is executed.
- Fixed variables appear as a single column in the new data file. Their values are simply copied to the new file.
- The AUTOFIX subcommand can automatically determine which variables in the original data are fixed. By default, the AUTOFIX subcommand overrides the FIXED subcommand.

## ***AUTOFIX subcommand***

The AUTOFIX subcommand evaluates candidate variables and classifies them as either fixed or as the source of a variable group.

- A **candidate variable** is a variable in the original data that does not appear on the SPLIT command or on the ID, INDEX, and DROP subcommands.
- An original variable that does not vary within any row group is classified as a **fixed variable** and is copied into a single variable in the new data file.



- An original variable that has only a single valid value plus the system-missing value within a row group is classified as a fixed variable and is copied into a single variable in the new data file.
- An original variable that does vary within the row group is classified as the **source of a variable group**. It becomes a variable group in the new data file.
- Use `AUTOFIX=NO` to overrule the default behavior and expand all variables not marked as ID or fixed or record into a variable group.

**YES**            *Evaluate and automatically classify all candidate variables.* The procedure automatically evaluates and classifies all candidate variables. This is the default. If there is a `FIXED` subcommand, the procedure displays a warning message for each misclassified variable and automatically corrects the error. Otherwise, no warning messages are displayed. This option overrides the `FIXED` subcommand.

**NO**            *Evaluate all candidate variables and issue warnings.* The procedure evaluates all candidate variables and determines if they are fixed. If a variable is listed on the `FIXED` subcommand but it is not actually fixed (that is, it varies within the row group), a warning message is displayed and the command is not executed. If a variable is not listed on the `FIXED` subcommand but it is actually fixed (that is, it does not vary within the row group), a warning message is displayed and the command is executed. The variable is classified as the source of a variable group and becomes a variable group in the new data file.

## ***RENAME subcommand***

`CASESTOVARS` creates variable groups with new variables. The first part of the new variable name is either derived from the name of the original variable or is the rootname specified on the `RENAME` subcommand.

- The specification is the original variable name followed by a rootname.
- The named variable cannot be a `SPLIT FILE` variable and cannot appear on the `ID`, `FIXED`, `INDEX`, or `DROP` subcommands.
- A variable can be renamed only once.
- Only one `RENAME` subcommand can be used, but it can contain multiple specifications.

## ***SEPARATOR subcommand***

`CASESTOVARS` creates variable groups that contain new variables. There are two parts to the name of a new variable—a rootname and an index. The parts are separated by a string. The separator string is specified on the `SEPARATOR` subcommand.

- If a separator is not specified, the default is a period.
- A separator can contain multiple characters.
- The separator must be delimited by single or double quotes.
- You can suppress the separator by specifying `/SEPARATOR=" "`.

## **GROUPBY subcommand**

The GROUPBY subcommand controls the order of the new variables in the new data file.

<b>VARIABLE</b>	<i>Group new variables by original variable.</i> The procedure groups all variables created from an original variable together. This is the default.
<b>INDEX</b>	<i>Group new variables by index variable.</i> The procedure groups variables according to the index variables.

### **Example**

If the original variables are:

```
insure    caseid    month    bps    bpd
```

and the data are as shown in the first example, the commands:

```
SPLIT FILE BY insure.
CASESTOVARS
  /ID=caseid
  /INDEX=month
  /GROUPBY=VARIABLE.
```

create a new data file with the following variable order:

- Variables are grouped by variable group—*bps* and *bpd*.

### **Example**

Using the same original data, the commands:

```
SPLIT FILE BY insure.
CASESTOVARS
  /ID=insure caseid
  /INDEX=month
  /GROUPBY=INDEX.
```

create a new data file with the following variable order:

- Variables are grouped by values of the index variable *month*—1, 2, and 3.

## **DROP subcommand**

The DROP subcommand specifies the subset of variables to exclude from the new data file.

- You can drop variables that appear on the ID list.
- Variables listed on the DROP subcommand may not appear on the FIXED or INDEX subcommand.
- Dropped variables are not candidates to become a variable group in the new data file.
- You cannot drop all variables. The new data file is required to have at least one variable.

# CATPCA

CATPCA is available in the Categories option.

```
CATPCA VARIABLES = varlist

/ANALYSIS = varlist
  [[ (WEIGHT={1**}) [LEVEL={SPORD**}] [DEGREE={2}] [INKNOT={2}]]
    {n }
    {SPNOM } [DEGREE={2}] [INKNOT={2}]
    {ORDI }
    {NOMI }
    {MNOM }
    {NUME }

[/DISCRETIZATION = [varlist[[({GROUPING })] [({NCAT={7} )] [DISTR={NORMAL }]]]]]
  {n }
  {EQINTV={n} }
  {RANKING }
  {MULTIPLYING}

[/MISSING = [varlist [[({PASSIVE**}) [({MODEIMPU})]]]]]
  {EXTRACAT}
  {ACTIVE } {MODEIMPU}
  {EXTRACAT}
  {LISTWISE}

[/SUPPLEMENTARY = [OBJECT(varlist)] [VARIABLE(varlist)]]

[/CONFIGURATION = [({INITIAL}) (file)]
  {FIXED }

[/DIMENSION = {2**}]
  {n }

[/NORMALIZATION = {VPRINCIPAL**}]
  {OPRINCIPAL }
  {SYMMETRICAL }
  {INDEPENDENT }
  {n }

[/MAXITER = {100**}]
  {n }

[/CRITITER = {.00001**}]
  {value }

[/PRINT = [DESCRIP**[(varlist)] [VAF] [LOADING**][QUANT[(varlist)]] [HISTORY]
  [CORR**] [OCORR] [OBJECT[({(varname)}varlist)] [NONE]]

[/PLOT = [OBJECT**[(varlist)][(n)]]
  [LOADING**[(varlist [(CENTR[(varlist)])]])] [(n)]
  [CATEGORY (varlist)[(n)]]
  [JOINTCAT[({varlist})] [(n)]]
  [TRANS[(varlist[({1})])] [(n)]]
  {n}
  [BIPLOT[({LOADING}[(varlist)])[(varlist)]] [(n)]]
  {CENTR }
  [TRIPLLOT[(varlist[(varlist)])] [(n)]]
  [RESID(varlist[({1})])] [(n)]]
  {n}
  [PROJCENTR(varname, varlist)[(n)]] [NONE]]
  [NDIM(value,value)]

[/SAVE = [TRDATA[({TRA }[(n)])] [OBJECT[({OBSCO }[(n)])]] ]
  {rootname} {rootname}
  [APPROX[({APP })] ]
  {rootname}
```

```
[/OUTFILE = [TRDATA[('savfile' | 'dataset')]] [DISCRDATA[('savfile' | 'dataset')]]
[OBJECT[('savfile' | 'dataset')]] [APPROX[('savfile' | 'dataset')]]].
```

\*\* Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- NDIM keyword introduced on PLOT subcommand.
- The maximum label length on the PLOT subcommand is increased to 64 for variable names, 255 for variable labels, and 60 for value labels (previous value was 20).

## **Overview**

CATPCA performs principal components analysis on a set of variables. The variables can be given mixed optimal scaling levels, and the relationships among observed variables are not assumed to be linear.

In CATPCA, dimensions correspond to components (that is, an analysis with two dimensions results in two components), and object scores correspond to component scores.

### **Options**

**Optimal Scaling Level.** You can specify the optimal scaling level at which you want to analyze each variable (levels include spline ordinal, spline nominal, ordinal, nominal, multiple nominal, or numerical).

**Discretization.** You can use the DISCRETIZATION subcommand to discretize fractional-value variables or to recode categorical variables.

**Missing Data.** You can use the MISSING subcommand to specify the treatment of missing data on a per-variable basis.

**Supplementary Objects and Variables.** You can specify objects and variables that you want to treat as supplementary to the analysis and then fit them into the solution.

**Read Configuration.** CATPCA can read a configuration from a file through the CONFIGURATION subcommand. This information can be used as the starting point for your analysis or as a fixed solution in which to fit variables.

**Number of Dimensions.** You can specify how many dimensions (components) CATPCA should compute.

**Normalization.** You can specify one of five different options for normalizing the objects and variables.

**Algorithm Tuning.** You can use the MAXITER and CRITITER subcommands to control the values of algorithm-tuning parameters.

**Optional Output.** You can request optional output through the `PRINT` subcommand.

**Optional Plots.** You can request a plot of object points, transformation plots per variable, and plots of category points per variable or a joint plot of category points for specified variables. Other plot options include residuals plots, a biplot, a triplot, component loadings plot, and a plot of projected centroids.

**Writing Discretized Data, Transformed Data, Object (Component) Scores, and Approximations.** You can write the discretized data, transformed data, object scores, and approximations to external files for use in further analyses.

**Saving Transformed Data, Object (Component) Scores, and Approximations.** You can save the transformed variables, object scores, and approximations to the working data file.

### ***Basic Specification***

The basic specification is the `CATPCA` command with the `VARIABLES` and `ANALYSIS` subcommands.

### ***Syntax Rules***

- The `VARIABLES` and `ANALYSIS` subcommands must always appear.
- All subcommands can be specified in any order.
- Variables that are specified in the `ANALYSIS` subcommand must be found in the `VARIABLES` subcommand.
- Variables that are specified in the `SUPPLEMENTARY` subcommand must be found in the `ANALYSIS` subcommand.

### ***Operations***

- If a subcommand is repeated, it causes a syntax error, and the procedure terminates.

### ***Limitations***

- `CATPCA` operates on category indicator variables. The category indicators should be positive integers. You can use the `DISCRETIZATION` subcommand to convert fractional-value variables and string variables into positive integers.
- In addition to system-missing values and user-defined missing values, category indicator values that are less than 1 are treated by `CATPCA` as missing. If one of the values of a categorical variable has been coded 0 or a negative value and you want to treat it as a valid category, use the `COMPUTE` command to add a constant to the values of that variable such that the lowest value will be 1. You can also use the `RANKING` option of the `DISCRETIZATION` subcommand for this purpose, except for variables that you want to treat as numeric, because the characteristic of equal intervals in the data will not be maintained.
- There must be at least three valid cases.
- Split-file has no implications for `CATPCA`.

## Example

```

CATPCA VARIABLES = TEST1 TEST2 TEST3 TO TEST6 TEST7 TEST8
  /ANALYSIS = TEST1 TO TEST2 (WEIGHT=2 LEVEL=ORDI)
              TEST3 TO TEST5 (LEVEL=SPORD INKNOT=3)
              TEST6 TEST7 (LEVEL=SPORD DEGREE=3)
              TEST8 (LEVEL=NUME)
  /DISCRETIZATION = TEST1 (GROUPING NCAT=5 DISTR=UNIFORM)
                  TEST6 (GROUPING) TEST8 (MULTIPLYING)
  /MISSING = TEST5 (ACTIVE) TEST6 (ACTIVE EXTRACAT) TEST8 (LISTWISE)
  /SUPPLEMENTARY = OBJECT(1 3) VARIABLE(TEST1)
  /CONFIGURATION = ('iniconf.sav')
  /DIMENSION = 2
  /NORMALIZATION = VPRINCIPAL
  /MAXITER = 150
  /CRITITER = .000001
  /PRINT = DESCRIP LOADING CORR QUANT(TEST1 TO TEST3) OBJECT
  /PLOT = TRANS(TEST2 TO TEST5) OBJECT(TEST2 TEST3)
  /SAVE = TRDATA OBJECT
  /OUTFILE = TRDATA('/data/trans.sav') OBJECT('/data/obs.sav') .

```

- `VARIABLES` defines variables. The keyword `TO` refers to the order of the variables in the working data file.
- The `ANALYSIS` subcommand defines variables that are used in the analysis. *TEST1* and *TEST2* have a weight of 2. For the other variables, `WEIGHT` is not specified; thus, they have the default weight value of 1. The optimal scaling level for *TEST1* and *TEST2* is ordinal. The optimal scaling level for *TEST3* to *TEST7* is spline ordinal. The optimal scaling level for *TEST8* is numerical. The keyword `TO` refers to the order of the variables in the `VARIABLES` subcommand. The splines for *TEST3* to *TEST5* have degree 2 (default because unspecified) and 3 interior knots. The splines for *TEST6* and *TEST7* have degree 3 and 2 interior knots (default because unspecified).
- `DISCRETIZATION` specifies that *TEST6* and *TEST8*, which are fractional-value variables, are discretized: *TEST6* by recoding into 7 categories with a normal distribution (default because unspecified) and *TEST8* by “multiplying.” *TEST1*, which is a categorical variable, is recoded into 5 categories with a close-to-uniform distribution.
- `MISSING` specifies that objects with missing values on *TEST5* and *TEST6* are included in the analysis; missing values on *TEST5* are replaced with the mode (default if not specified), and missing values on *TEST6* are treated as an extra category. Objects with a missing value on *TEST8* are excluded from the analysis. For all other variables, the default is in effect; that is, missing values (not objects) are excluded from the analysis.
- `CONFIGURATION` specifies *iniconf.sav* as the file containing the coordinates of a configuration that is to be used as the initial configuration (default because unspecified).
- `DIMENSION` specifies 2 as the number of dimensions; that is, 2 components are computed. This setting is the default, so this subcommand could be omitted here.
- The `NORMALIZATION` subcommand specifies optimization of the association between variables. This setting is the default, so this subcommand could be omitted here.
- `MAXITER` specifies 150 as the maximum number of iterations (instead of the default value of 100).
- `CRITITER` sets the convergence criterion to a value that is smaller than the default value.
- `PRINT` specifies descriptives, component loadings and correlations (all default), quantifications for *TEST1* to *TEST3*, and the object (component) scores.

- PLOT requests transformation plots for the variables *TEST2* to *TEST5*, an object points plot labeled with the categories of *TEST2*, and an object points plot labeled with the categories of *TEST3*.
- The SAVE subcommand adds the transformed variables and the component scores to the working data file.
- The OUTFILE subcommand writes the transformed data to a data file called *trans.sav* and writes the component scores to a data file called *obs.sav*, both in the directory */data*.

## **VARIABLES Subcommand**

VARIABLES specifies the variables that may be analyzed in the current CATPCA procedure.

- The VARIABLES subcommand is required.
- At least two variables must be specified, except when the CONFIGURATION subcommand is used with the FIXED keyword.
- The keyword TO on the VARIABLES subcommand refers to the order of variables in the working data file. This behavior of TO is different from the behavior in the variable list in the ANALYSIS subcommand.

## **ANALYSIS Subcommand**

ANALYSIS specifies the variables to be used in the computations, the optimal scaling level, and the variable weight for each variable or variable list. ANALYSIS also specifies supplementary variables and their optimal scaling level. No weight can be specified for supplementary variables.

- At least two variables must be specified, except when the CONFIGURATION subcommand is used with the FIXED keyword.
- All variables on ANALYSIS must be specified on the VARIABLES subcommand.
- The ANALYSIS subcommand is required.
- The keyword TO in the variable list honors the order of variables in the VARIABLES subcommand.
- Optimal scaling levels and variable weights are indicated by the keywords LEVEL and WEIGHT in parentheses following the variable or variable list.

**WEIGHT**            *Specifies the variable weight with a positive integer.* The default value is 1. If WEIGHT is specified for supplementary variables, it is ignored, and a syntax warning is issued.

**LEVEL**            *Specifies the optimal scaling level.*

## Level Keyword

The following keywords are used to indicate the optimal scaling level:

<b>SPORD</b>	<i>Spline ordinal (monotonic)</i> . This setting is the default. The order of the categories of the observed variable is preserved in the optimally scaled variable. Category points will lie on a straight line (vector) through the origin. The resulting transformation is a smooth monotonic piecewise polynomial of the chosen degree. The pieces are specified by the user-specified number and procedure-determined placement of the interior knots.
<b>SPNOM</b>	<i>Spline nominal (nonmonotonic)</i> . The only information in the observed variable that is preserved in the optimally scaled variable is the grouping of objects in categories. The order of the categories of the observed variable is not preserved. Category points will lie on a straight line (vector) through the origin. The resulting transformation is a smooth, possibly nonmonotonic, piecewise polynomial of the chosen degree. The pieces are specified by the user-specified number and procedure-determined placement of the interior knots.
<b>MNOM</b>	<i>Multiple nominal</i> . The only information in the observed variable that is preserved in the optimally scaled variable is the grouping of objects in categories. The order of the categories of the observed variable is not preserved. Category points will be in the centroid of the objects in the particular categories. Multiple indicates that different sets of quantifications are obtained for each dimension.
<b>ORDI</b>	<i>Ordinal</i> . The order of the categories on the observed variable is preserved in the optimally scaled variable. Category points will lie on a straight line (vector) through the origin. The resulting transformation fits better than SPORD transformation but is less smooth.
<b>NOMI</b>	<i>Nominal</i> . The only information in the observed variable that is preserved in the optimally scaled variable is the grouping of objects in categories. The order of the categories of the observed variable is not preserved. Category points will lie on a straight line (vector) through the origin. The resulting transformation fits better than SPNOM transformation but is less smooth.
<b>NUME</b>	<i>Numerical</i> . Categories are treated as equally spaced (interval level). The order of the categories and the equal distances between category numbers of the observed variables are preserved in the optimally scaled variable. Category points will lie on a straight line (vector) through the origin. When all variables are scaled at the numerical level, the CATPCA analysis is analogous to standard principal components analysis.

## SPORD and SPNOM Keywords

The following keywords are used with SPORD and SPNOM:

<b>DEGREE</b>	<i>The degree of the polynomial</i> . It can be any positive integer. The default degree is 2.
<b>INKNOT</b>	<i>The number of interior knots</i> . The minimum is 0, and the maximum is the number of categories of the variable minus 2. If the specified value is too large, the procedure adjusts the number of interior knots to the maximum. The default number of interior knots is 2.

## DISCRETIZATION Subcommand

DISCRETIZATION specifies fractional-value variables that you want to discretize. Also, you can use DISCRETIZATION for ranking or for two ways of recoding categorical variables.

- A string variable's values are always converted into positive integers, according to the internal numeric representations. DISCRETIZATION for string variables applies to these integers.



- When the `DISCRETIZATION` subcommand is omitted or used without a variable list, fractional-value variables are converted into positive integers by grouping them into seven categories with a distribution of close to “normal.”
- When no specification is given for variables in a variable list following `DISCRETIZATION`, these variables are grouped into seven categories with a distribution of close to “normal.”
- In `CATPCA`, values that are less than 1 are considered to be missing (see `MISSING` subcommand). However, when discretizing a variable, values that are less than 1 are considered to be valid and are thus included in the discretization process.

<b>GROUPING</b>	<i>Recode into the specified number of categories or recode intervals of equal size into categories.</i>
<b>RANKING</b>	<i>Rank cases. Rank 1 is assigned to the case with the smallest value on the variable.</i>
<b>MULTIPLYING</b>	<i>Multiply the standardized values of a fractional-value variable by 10, round, and add a value such that the lowest value is 1.</i>

### ***GROUPING Keyword***

`GROUPING` has the following keywords:

<b>NCAT</b>	<i>Number of categories. When <code>NCAT</code> is not specified, the number of categories is set to 7.</i>
<b>EQINTV</b>	<i>Recode intervals of equal size. The size of the intervals must be specified (no default). The resulting number of categories depends on the interval size.</i>

### ***NCAT Keyword***

`NCAT` has the keyword `DISTR`, which has the following keywords:

<b>NORMAL</b>	<i>Normal distribution. This setting is the default when <code>DISTR</code> is not specified.</i>
<b>UNIFORM</b>	<i>Uniform distribution.</i>

### ***MISSING Subcommand***

In `CATPCA`, we consider a system-missing value, user-defined missing values, and values that are less than 1 as missing values. The `MISSING` subcommand allows you to indicate how to handle missing values for each variable.

<b>PASSIVE</b>	<i>Exclude missing values on a variable from analysis. This setting is the default when <code>MISSING</code> is not specified. Passive treatment of missing values means that in optimizing the quantification of a variable, only objects with nonmissing values on the variable are involved and that only the nonmissing values of variables contribute to the solution. Thus, when <code>PASSIVE</code> is specified, missing values do not affect the</i>
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analysis. Further, if all variables are given passive treatment of missing values, objects with missing values on every variable are treated as supplementary.

**ACTIVE**

*Impute missing values.* You can choose to use mode imputation. You can also consider objects with missing values on a variable as belonging to the same category and impute missing values with an extra category indicator.

**LISTWISE**

*Exclude cases with missing values on a variable.* The cases that are used in the analysis are cases without missing values on the specified variables. Also, any variable that is not included in the subcommand receives this specification.

- The ALL keyword may be used to indicate all variables. If ALL is used, it must be the only variable specification.
- A mode or extracat imputation is done before listwise deletion.

**PASSIVE Keyword**

If correlations are requested on the PRINT subcommand, and passive treatment of missing values is specified for a variable, the missing values must be imputed. For the correlations of the quantified variables, you can specify the imputation with one of the following keywords:

**MODEIMPU**      *Impute missing values on a variable with the mode of the quantified variable.*  
MODEIMPU is the default.

**EXTRACAT**      *Impute missing values on a variable with the quantification of an extra category.*  
This treatment implies that objects with a missing value are considered to belong to the same (extra) category.

Note that with passive treatment of missing values, imputation applies only to correlations and is done afterward. Thus, the imputation has no effect on the quantification or the solution.

**ACTIVE Keyword**

The ACTIVE keyword has the following keywords:

**MODEIMPU**      *Impute missing values on a variable with the most frequent category (mode).* When there are multiple modes, the smallest category indicator is used. MODEIMPU is the default.

**EXTRACAT**      *Impute missing values on a variable with an extra category indicator.* This implies that objects with a missing value are considered to belong to the same (extra) category.

Note that with active treatment of missing values, imputation is done before the analysis starts and thus will affect the quantification and the solution.

**SUPPLEMENTARY Subcommand**

The SUPPLEMENTARY subcommand specifies the objects and/or variables that you want to treat as supplementary. Supplementary variables must be found in the ANALYSIS subcommand. You cannot weight supplementary objects and variables (specified weights are ignored). For

supplementary variables, all options on the `MISSING` subcommand can be specified except `LISTWISE`.

<b>OBJECT</b>	<i>Objects that you want to treat as supplementary are indicated with an object number list in parentheses following <code>OBJECT</code>. The keyword <code>TO</code> is allowed. The <code>OBJECT</code> specification is not allowed when <code>CONFIGURATION = FIXED</code>.</i>
<b>VARIABLE</b>	<i>Variables that you want to treat as supplementary are indicated with a variable list in parentheses following <code>VARIABLE</code>. The keyword <code>TO</code> is allowed and honors the order of variables in the <code>VARIABLES</code> subcommand. The <code>VARIABLE</code> specification is ignored when <code>CONFIGURATION = FIXED</code>, because in that case all variables in the <code>ANALYSIS</code> subcommand are automatically treated as supplementary variables.</i>

## ***CONFIGURATION Subcommand***

The `CONFIGURATION` subcommand allows you to read data from a file containing the coordinates of a configuration. The first variable in this file should contain the coordinates for the first dimension, the second variable should contain the coordinates for the second dimension, and so forth.

<b>INITIAL(file)</b>	<i>Use the configuration in the external file as the starting point of the analysis.</i>
<b>FIXED(file)</b>	<i>Fit variables in the fixed configuration that is found in the external file. The variables to fit in should be specified on the <code>ANALYSIS</code> subcommand but will be treated as supplementary. The <code>SUPPLEMENTARY</code> subcommand and variable weights are ignored.</i>

## ***DIMENSION Subcommand***

`DIMENSION` specifies the number of dimensions (components) that you want CATPCA to compute.

- The default number of dimensions is 2.
- `DIMENSION` is followed by an integer indicating the number of dimensions.
- If there are no variables specified as `MNOM` (multiple nominal), the maximum number of dimensions that you can specify is the smaller of the number of observations minus 1 and the total number of variables.
- If some or all of the variables are specified as `MNOM` (multiple nominal), the maximum number of dimensions is the smaller of a) the number of observations minus 1 and b) the total number of valid `MNOM` variable levels (categories) plus the number of `SPORD`, `SPNOM`, `ORDI`, `NOMI`, and `NUME` variables minus the number of `MNOM` variables (if the `MNOM` variables do not have missing values to be treated as passive). If there are `MNOM` variables with missing values to be treated as passive, the maximum number of dimensions is the smaller of a) the number of observations minus 1 and b) the total number of valid `MNOM` variable levels (categories) plus the number of `SPORD`, `SPNOM`, `ORDI`, `NOMI`, and `NUME` variables, minus the larger of c) 1 and d) the number of `MNOM` variables without missing values to be treated as passive.
- If the specified value is too large, CATPCA adjusts the number of dimensions to the maximum.
- The minimum number of dimensions is 1.

## ***NORMALIZATION Subcommand***

The `NORMALIZATION` subcommand specifies one of five options for normalizing the object scores and the variables. Only one normalization method can be used in a given analysis.

<b>VPRINCIPAL</b>	<i>This option optimizes the association between variables. With <code>VPRINCIPAL</code>, the coordinates of the variables in the object space are the component loadings (correlations with object scores) for <code>SPORD</code>, <code>SPNOM</code>, <code>ORDI</code>, <code>NOMI</code>, and <code>NUME</code> variables, and the centroids for <code>MNOM</code> variables. This setting is the default if the <code>NORMALIZATION</code> subcommand is not specified. This setting is useful when you are primarily interested in the correlations between the variables.</i>
<b>OPRINCIPAL</b>	<i>This option optimizes distances between objects. This setting is useful when you are primarily interested in differences or similarities between the objects.</i>
<b>SYMMETRICAL</b>	<i>Use this normalization option if you are primarily interested in the relation between objects and variables.</i>
<b>INDEPENDENT</b>	<i>Use this normalization option if you want to examine distances between objects and correlations between variables separately.</i>

The fifth method allows the user to specify any real value in the closed interval  $[-1, 1]$ . A value of 1 is equal to the `OPRINCIPAL` method, a value of 0 is equal to the `SYMMETRICAL` method, and a value of  $-1$  is equal to the `VPRINCIPAL` method. By specifying a value that is greater than  $-1$  and less than 1, the user can spread the eigenvalue over both objects and variables. This method is useful for making a tailor-made biplot or triplot. If the user specifies a value outside of this interval, the procedure issues a syntax error message and terminates.

## ***MAXITER Subcommand***

`MAXITER` specifies the maximum number of iterations that the procedure can go through in its computations. If not all variables are specified as `NUME` and/or `MNOM`, the output starts from iteration 0, which is the last iteration of the initial phase, in which all variables except `MNOM` variables are treated as `NUME`.

- If `MAXITER` is not specified, the maximum number of iterations is 100.
- The specification on `MAXITER` is a positive integer indicating the maximum number of iterations. There is no uniquely predetermined (that is, hard-coded) maximum for the value that can be used.

## ***CRITITER Subcommand***

`CRITITER` specifies a convergence criterion value. `CATPCA` stops iterating if the difference in fit between the last two iterations is less than the `CRITITER` value.

- If `CRITITER` is not specified, the convergence value is 0.00001.
- The specification on `CRITITER` is any positive value.

## **PRINT Subcommand**

The Model Summary (Cronbach's alpha and Variance Accounted For) and the HISTORY statistics (the variance accounted for, the loss, and the increase in variance accounted for) for the initial solution (if applicable) and last iteration are always displayed. That is, they cannot be controlled by the PRINT subcommand. The PRINT subcommand controls the display of additional optional output. The output of the procedure is based on the transformed variables. However, the keyword OCORR can be used to request the correlations of the original variables, as well.

The default keywords are DESCRIP, LOADING, and CORR. However, when some keywords are specified, the default is nullified and only what was specified comes into effect. If a keyword is duplicated or if a contradicting keyword is encountered, the last specified keyword silently becomes effective (in case of contradicting use of NONE, only the keywords following NONE are effective). An example is as follows:

```
/PRINT <=> /PRINT = DESCRIP LOADING CORR
```

```
/PRINT = VAF VAF <=> /PRINT = VAF
```

```
/PRINT = VAF NONE CORR <=> /PRINT = CORR
```

If a keyword that can be followed by a variable list is duplicated, a syntax error occurs, and the procedure will terminate.

The following keywords can be specified:

<b>DESCRIP(varlist)</b>	<i>Descriptive statistics (frequencies, missing values, and mode).</i> The variables in the varlist must be specified on the VARIABLES subcommand but need not appear on the ANALYSIS subcommand. If DESCRIP is not followed by a varlist, descriptives tables are displayed for all variables in the varlist on the ANALYSIS subcommand.
<b>VAF</b>	<i>Variance accounted for (centroid coordinates, vector coordinates, and total) per variable and per dimension.</i>
<b>LOADING</b>	<i>Component loadings for variables with optimal scaling level that result in vector quantification (that is, SPORD, SPNOM, ORDI, NOMI, and NUME).</i>
<b>QUANT(varlist)</b>	<i>Category quantifications and category coordinates for each dimension.</i> Any variable in the ANALYSIS subcommand may be specified in parentheses after QUANT. (For MNOM variables, the coordinates are the quantifications.) If QUANT is not followed by a variable list, quantification tables are displayed for all variables in the varlist on the ANALYSIS subcommand.
<b>HISTORY</b>	<i>History of iterations.</i> For each iteration (including 0, if applicable), the variance accounted for, the loss (variance not accounted for), and the increase in variance accounted for are shown.
<b>CORR</b>	<i>Correlations of the transformed variables and the eigenvalues of this correlation matrix.</i> If the analysis includes variables with optimal scaling level MNOM, <i>ndim</i> (the number of dimensions in the analysis) correlation matrices are computed; in the <i>i</i> th matrix, the quantifications of dimension <i>i</i> , <i>i</i> = 1, ... <i>ndim</i> , of MNOM variables are used to compute the correlations. For variables with missing values specified to be treated as PASSIVE on the MISSING subcommand, the missing values are imputed according to the specification on the PASSIVE keyword (if no specification is made, mode imputation is used).

<b>OCORR</b>	<i>Correlations of the original variables and the eigenvalues of this correlation matrix.</i> For variables with missing values specified to be treated as <code>PASSIVE</code> on the <code>MISSING</code> subcommand, the missing values are imputed with the variable mode.
<b>OBJECT((varname)varlist)</b>	<i>Object scores (component scores).</i> Following the keyword, a varlist can be given in parentheses to display variables (category indicators), along with object scores. If you want to use a variable to label the objects, this variable must occur in parentheses as the first variable in the varlist. If no labeling variable is specified, the objects are labeled with case numbers. The variables to display, along with the object scores and the variable to label the objects, must be specified on the <code>VARIABLES</code> subcommand but need not appear on the <code>ANALYSIS</code> subcommand. If no variable list is given, only the object scores are displayed.
<b>NONE</b>	<i>No optional output is displayed.</i> The only output that is shown is the model summary and the <code>HISTORY</code> statistics for the initial iteration (if applicable) and last iteration.

The keyword `TO` in a variable list can only be used with variables that are in the `ANALYSIS` subcommand, and `TO` applies only to the order of the variables in the `ANALYSIS` subcommand. For variables that are in the `VARIABLES` subcommand but not in the `ANALYSIS` subcommand, the keyword `TO` cannot be used. For example, if `/VARIABLES = v1 TO v5` and `/ANALYSIS = v2 v1 v4`, then `/PLOT OBJECT(v1 TO v4)` will give two object plots (one plot labeled with `v1` and one plot labeled with `v4`).

## ***PLOT Subcommand***

The `PLOT` subcommand controls the display of plots. The default keywords are `OBJECT` and `LOADING`. That is, the two keywords are in effect when the `PLOT` subcommand is omitted or when the `PLOT` subcommand is given without any keyword. If a keyword is duplicated (for example, `/PLOT = RESID RESID`), only the last keyword is effective. If the keyword `NONE` is used with other keywords (for example, `/PLOT = RESID NONE LOADING`), only the keywords following `NONE` are effective. When keywords contradict, the later keyword overwrites the earlier keywords.

- All the variables to be plotted must be specified on the `ANALYSIS` subcommand.
- If the variable list following the keywords `CATEGORIES`, `TRANS`, `RESID`, and `PROJCENTR` is empty, it will cause a syntax error, and the procedure will terminate.
- The variables in the variable list for labeling the object point following `OBJECT`, `BIPLOT`, and `TRIPLOT` must be specified on the `VARIABLES` subcommand but need not appear on the `ANALYSIS` subcommand. This flexibility means that variables that are not included in the analysis can still be used to label plots.
- The keyword `TO` in a variable list can only be used with variables that are in the `ANALYSIS` subcommand, and `TO` applies only to the order of the variables in the `ANALYSIS` subcommand. For variables that are in the `VARIABLES` subcommand but not in the `ANALYSIS` subcommand, the keyword `TO` cannot be used. For example, if `/VARIABLES = v1 TO v5` and

/ANALYSIS = v2 v1 v4, then /PLOT OBJECT(v1 TO v4) will give two object plots, one plot labeled with v1 and one plot labeled with v4.

- For multidimensional plots, all of the dimensions in the solution are produced in a matrix scatterplot if the number of dimensions in the solution is greater than 2 and the NDIM plot keyword is not specified; if the number of dimensions in the solution is 2, a scatterplot is produced.

The following keywords can be specified:

<b>OBJECT(varlist)(n)</b>	<i>Plots of the object points.</i> Following the keyword, a list of variables in parentheses can be given to indicate that plots of object points labeled with the categories of the variables should be produced (one plot for each variable). The variables to label the objects must be specified on the VARIABLES subcommand but need not appear on the ANALYSIS subcommand. If the variable list is omitted, a plot that is labeled with case numbers is produced.
<b>CATEGORY(varlist)(n)</b>	<i>Plots of the category points.</i> Both the centroid coordinates and the vector coordinates are plotted. A list of variables must be given in parentheses following the keyword. For variables with optimal scaling level MNOM, categories are in the centroids of the objects in the particular categories. For all other optimal scaling levels, categories are on a vector through the origin.
<b>LOADING(varlist (CENTR(varlist)))(l)</b>	<i>Plot of the component loadings optionally with centroids.</i> By default, all variables with an optimal scaling level that results in vector quantification (that is, SPORD, SPNOM, ORDI, NOMI, and NUME) are included in this plot. LOADING can be followed by a varlist to select the loadings to include in the plot. When "LOADING (" or the varlist following "LOADING (" is followed by the keyword CENTR in parentheses, centroids are included in the plot for all variables with optimal scaling level MNOM. CENTR can be followed by a varlist in parentheses to select MNOM variables whose centroids are to be included in the plot. When all variables have the MNOM scaling level, this plot cannot be produced.
<b>TRANS(varlist(n))(n)</b>	<i>Transformation plots per variable (optimal category quantifications against category indicators).</i> Following the keyword, a list of variables in parentheses must be given. MNOM variables in the varlist can be followed by a number of dimensions in parentheses to indicate that you want to display $p$ transformation plots, one plot for each of the first $p$ dimensions. If the number of dimensions is not specified, a plot for the first dimension is produced.
<b>RESID(varlist(n))(n)</b>	<i>Plot of residuals per variable (approximation against optimal category quantifications).</i> Following the keyword, a list of variables in parentheses must be given. MNOM variables in the varlist can be followed by a number of dimensions in parentheses to indicate that you want to display $p$ residual plots, one plot for each of the first $p$ dimensions. If the number of dimensions is not specified, a plot for the first dimension is produced.

<b>BIPLOT(keyword(varlist)) (varlist)(n)</b>	<i>Plot of objects and variables.</i> The coordinates for the variables can be chosen to be component loading or centroids, using the <code>LOADING</code> or <code>CENTR</code> keyword in parentheses following <code>BIPLOT</code> . When no keyword is given, component loadings are plotted. When <code>NORMALIZATION = INDEPENDENT</code> , this plot is incorrect and therefore not available. Following <code>LOADING</code> or <code>CENTR</code> , a list of variables in parentheses can be given to indicate the variables to be included in the plot. If the variable list is omitted, a plot including all variables is produced. Following <code>BIPLOT</code> , a list of variables in parentheses can be given to indicate that plots with objects that are labeled with the categories of the variables should be produced (one plot for each variable). The variables to label the objects must be specified on the <code>VARIABLES</code> subcommand but need not appear on the <code>ANALYSIS</code> subcommand. If the variable list is omitted, a plot with objects labeled with case numbers is produced.
<b>TRIPLLOT(varlist(varlist))(n)</b>	<i>A plot of object points, component loadings for variables with an optimal scaling level that results in vector quantification (that is, <code>SPORD</code>, <code>SPNOM</code>, <code>ORDI</code>, <code>NOMI</code>, and <code>NUME</code>), and centroids for variables with optimal scaling level <code>MNOM</code>.</i> Following the keyword, a list of variables in parentheses can be given to indicate the variables to include in the plot. If the variable list is omitted, all variables are included. The varlist can contain a second varlist in parentheses to indicate that triplots with objects labeled with the categories of the variables in this variable list should be produced (one plot for each variable). The variables to label the objects must be specified on the <code>VARIABLES</code> subcommand but need not appear on the <code>ANALYSIS</code> subcommand. If this second variable list is omitted, a plot with objects labeled with case numbers is produced. When <code>NORMALIZATION = INDEPENDENT</code> , this plot is incorrect and therefore not available.
<b>JOINTCAT(varlist)(n)</b>	<i>Joint plot of the category points for the variables in the varlist.</i> If no varlist is given, the category points for all variables are displayed.
<b>PROJCENTR(varname, varlist)(n)</b>	<i>Plot of the centroids of a variable projected on each of the variables in the varlist.</i> You cannot project centroids of a variable on variables with <code>MNOM</code> optimal scaling level; thus, a variable that has <code>MNOM</code> optimal scaling level can be specified as the variable to be projected but not in the list of variables to be projected on. When this plot is requested, a table with the coordinates of the projected centroids is also displayed.
<b>NONE</b>	<i>No plots.</i>

- For all keywords except `NONE`, the user can specify an optional parameter *l* in parentheses after the variable list in order to control the global upper boundary of variable name/label and value label lengths in the plot. Note that this boundary is applied uniformly to all variables in the list. The label length parameter *l* can take any non-negative integer that is less than or equal to the applicable maximum length (64 for variable names, 255 for variable labels, and 60 for value labels). If *l* = 0, names/values instead of variable/value labels are displayed to indicate variables/categories. If *l* is not specified, `CATPCA` assumes that each variable name/label and value label is displayed at its full length. If *l* is an integer that is larger than the applicable maximum, we reset it to the applicable maximum but do not issue a warning. If a positive value of *l* is given but some or all variables/category values do not have labels, then, for those variables/values, the names/values themselves are used as the labels.

In addition to the plot keywords, the following keyword can be specified:

<b>NDIM(value,value)</b>	<i>Dimension pairs to be plotted.</i> <code>NDIM</code> is followed by a pair of values in parentheses. If <code>NDIM</code> is not specified or is specified without parameter values, a matrix scatterplot including all dimensions is produced.
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- The first value (an integer that can range from 1 to the number of dimensions in the solution minus 1) indicates the dimension that is plotted against higher dimensions.
- The second value (an integer that can range from 2 to the number of dimensions in the solution) indicates the highest dimension to be used in plotting the dimension pairs.
- The `NDIM` specification applies to all requested multidimensional plots.

### **BIPLOT Keyword**

`BIPLOT` takes the following keywords:

<b>LOADING</b> (varlist)	<i>Object points and component loadings.</i>
<b>CENTR</b> (varlist)	<i>Object points and centroids.</i>

### **SAVE Subcommand**

The `SAVE` subcommand is used to add the transformed variables (category indicators that are replaced with optimal quantifications), the object scores, and the approximation to the working data file. Excluded cases are represented by a dot (the system-missing symbol) on every saved variable.

<b>TRDATA</b>	<i>Transformed variables. Missing values that are specified to be treated as passive are represented by a dot.</i>
<b>OBJECT</b>	<i>Object (component) scores.</i>
<b>APPROX</b>	<i>Approximation for variables that do not have optimal scaling level <code>MNOM</code>. For variables with <code>MNOM</code> scaling level, the approximations in dimension <math>s</math> are the object scores in dimension <math>s</math>.</i>

- Following `TRDATA`, a rootname and the number of dimensions to be saved for variables that are specified as `MNOM` can be specified in parentheses.
- For variables that are not specified as `MNOM`, `CATPCA` adds two numbers separated by the symbol `_`. For variables that are specified as `MNOM`, `CATPCA` adds three numbers. The first number uniquely identifies the source variable names, and the last number uniquely identifies the `CATPCA` procedures with the successfully executed `SAVE` subcommands. For variables that are specified as `MNOM`, the middle number corresponds to the dimension number (see the next bullet for more details). Only one rootname can be specified, and it can contain up to five characters for variables that are not specified as `MNOM` and three characters for variables that are specified as `MNOM`. If more than one rootname is specified, the first rootname is used. If a rootname contains more than five characters (`MNOM` variables), the first five characters are used at most. If a rootname contains more than three characters (`MNOM` variables), the first three characters are used at most.
- If a rootname is not specified for `TRDATA`, rootname `TRA` is used to automatically generate unique variable names. The formulas are `ROOTNAME $k$ _ $n$`  and `ROOTNAME $k$ _ $m$ _ $n$` . In this formula,  $k$  increments from 1 to identify the source variable names by using the source variables' position numbers in the `ANALYSIS` subcommand,  $m$  increments from 1 to identify the dimension number, and  $n$  increments from 1 to identify the `CATPCA` procedures with the successfully executed `SAVE` subcommands for a given data file in a continuous session.

For example, with three variables specified on `ANALYSIS`, `LEVEL = MNOM` for the second variable, and with two dimensions to save, the first set of default names—if they do not exist in the data file—would be `TRA1_1`, `TRA2_1_1`, `TRA2_2_1`, and `TRA3_1`. The next set of default names—if they do not exist in the data file—would be `TRA1_2`, `TRA2_1_2`, `TRA2_2_2`, and `TRA3_2`. However, if, for example, `TRA1_2` already exists in the data file, the default names should be attempted as `TRA1_3`, `TRA2_1_3`, `TRA2_2_3`, and `TRA3_3`. That is, the last number increments to the next available integer.

- Following `OBJECT`, a rootname and the number of dimensions can be specified in parentheses, to which CATPCA adds two numbers separated by the symbol `_`. The first number corresponds to the dimension number. The second number uniquely identifies the CATPCA procedures with the successfully executed `SAVE` subcommands (see the next bullet for more details). Only one rootname can be specified, and it can contain up to five characters. If more than one rootname is specified, the first rootname is used; if a rootname contains more than five characters, the first five characters are used at most.
- If a rootname is not specified for `OBJECT`, rootname `OBSCO` is used to automatically generate unique variable names. The formula is `ROOTNAME $m$ _ $n$` . In this formula,  $m$  increments from 1 to identify the dimension number, and  $n$  increments from 1 to identify the CATPCA procedures with the successfully executed `SAVE` subcommands for a given data file in a continuous session. For example, if two dimensions are specified following `OBJECT`, the first set of default names—if they do not exist in the data file—would be `OBSCO1_1` and `OBSCO2_1`. The next set of default names—if they do not exist in the data file—would be `OBSCO1_2` and `OBSCO2_2`. However, if, for example, `OBSCO2_2` already exists in the data file, the default names should be attempted as `OBSCO1_3` and `OBSCO2_3`. That is, the second number increments to the next available integer.
- Following `APPROX`, a rootname can be specified in parentheses, to which CATPCA adds two numbers separated by the symbol `_`. The first number uniquely identifies the source variable names, and the last number uniquely identifies the CATPCA procedures with the successfully executed `SAVE` subcommands (see the next bullet for more details). Only one rootname can be specified, and it can contain up to five characters. If more than one rootname is specified, the first rootname is used; if a rootname contains more than five characters, the first five characters are used at most.
- If a rootname is not specified for `APPROX`, rootname `APP` is used to automatically generate unique variable names. The formula is `ROOTNAME $k$ _ $n$` . In this formula,  $k$  increments from 1 to identify the source variable names by using the source variables' position numbers in the `ANALYSIS` subcommand. Additionally,  $n$  increments from 1 to identify the CATPCA procedures with the successfully executed `SAVE` subcommands for a given data file in a continuous session. For example, with three variables specified on `ANALYSIS` and `LEVEL = MNOM` for the second variable, the first set of default names—if they do not exist in the data file—would be `APP1_1`, `APP2_1`, and `APP3_1`. The next set of default names—if they do not exist in the data file—would be `APP1_2`, `APP2_2`, and `APP3_2`. However, if, for example, `APP1_2` already exists in the data file, the default names should be attempted as `APP1_3`, `APP2_3`, and `APP3_3`. That is, the last number increments to the next available integer.
- Variable labels are created automatically. (They are shown in the Notes table and can also be displayed in the Data Editor window.)
- If the number of dimensions is not specified, the `SAVE` subcommand saves all dimensions.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand is used to write the discretized data, transformed data (category indicators replaced with optimal quantifications), the object scores, and the approximation to a data file or previously declared data set. Excluded cases are represented by a dot (the system-missing symbol) on every saved variable.

<b>DISCRDATA</b> ( <code>'savfile'</code>   <code>'dataset'</code> )	<i>Discretized data.</i>
<b>TRDATA</b> ( <code>'savfile'</code>   <code>'dataset'</code> )	<i>Transformed variables.</i> This setting is the default if the <code>OUTFILE</code> subcommand is specified with a filename and without a keyword. Missing values that are specified to be treated as passive are represented by a dot.
<b>OBJECT</b> ( <code>'savfile'</code>   <code>'dataset'</code> )	<i>Object (component) scores.</i>
<b>APPROX</b> ( <code>'savfile'</code>   <code>'dataset'</code> )	<i>Approximation for variables that do not have optimal scaling level MNOM.</i>

- Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. Data sets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files. The names should be different for each of the keywords.

In principle, the active data set should not be replaced by this subcommand, and the asterisk (\*) file specification is not supported. This strategy also prevents `OUTFILE` interference with the `SAVE` subcommand.

# CATREG

CATREG is available in the Categories option.

```
CATREG VARIABLES = varlist

/ANALYSIS =
  depvar [([LEVEL={SPORD**}] [DEGREE={2}] [INKNOT={2}])]
          {n} {n}
          {SPNOM } [DEGREE={2}] [INKNOT={2}]
          {n} {n}
          {ORDI }
          {NOMI }
          {NUME }
  WITH indvarlist [([LEVEL={SPORD**}] [DEGREE={2}] [INKNOT={2}])]
                  {n} {n}
                  {SPNOM } [DEGREE={2}] [INKNOT={2}]
                  {n} {n}
                  {ORDI }
                  {NOMI }
                  {NUME }

[/DISCRETIZATION = [varlist [({GROUPING })] [{NCAT*={7}}] [DISTR={NORMAL }]]]
                   {n} {UNIFORM}
                   {EQINTV=n }
                   {RANKING }
                   {MULTIPLYING}

[/MISSING = [{varlist}({LISTWISE**})]
            {ALL** } {MODEIMPU }
            {EXTRACAT }

[/SUPPLEMENTARY = OBJECT(objlist)]

[/INITIAL = [{NUMERICAL**}]
            {RANDOM }
            {MULTISTART } ({50 }) ('savfile'|'dataset')
            {n }
            {ALL}
            {FIXSIGNS } (n) ('filename')

[/MAXITER = [{100**}]
            {n }

[/CRITITER = [{.00001**}]
            {value }

[/REGULARIZATION = [{NONE**}]
                   {RIDGE } [({ 0, 1.0, 0.02})] ('filename')
                   {(value, value, value) }
                   {LASSO } [({ 0, 1.0, 0.02})] ('filename')
                   {(value, value, value) }
                   {ENET } [({ 0, 1.0, 0.1})( 0, 1.0, .02)] ('filename')
                   {(value, value, value)(value, value, value)}

[/RESAMPLE = [{NONE** }]]
            {CROSSVAL }[({10})]
            {n }
            {BOOTSTRAP}[({50})]
            {n }

[/PRINT = [R**] [COEFF**] [DESCRIP**](varlist)] [HISTORY] [ANOVA**]
          [CORR] [OCORR] [QUANT](varlist)] [REGU] [NONE]]

[/PLOT = [TRANS(varlist)](h)] [RESID(varlist)](h)] [REGU({valuelist})]
        {ALL }

[/SAVE = [TRDATA[({TRA }))] [PRED[({PRE }))] [RES[({RES }))]
        {rootname} {rootname} {rootname}
```

```
[/OUTFILE = [TRDATA('savfile'|'dataset')] [DISCRDATA('savfile'|'dataset')]] .
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

#### Release 13.0

- The maximum category label length on the `PLOT` subcommand is increased to 60 (previous value was 20).

#### Release 17.0

- `MULTISTART` and `FIXSIGNS` keywords added to `INITIAL` subcommand.
- `REGULARIZATION` subcommand added.
- `RESAMPLE` subcommand added.
- `REGU` keyword added to `PRINT` subcommand.
- `REGU` keyword added to `PLOT` subcommand.
- `SUPPLEMENTARY` categories not occurring in data used to create the model are now interpolated.

## **Overview**

CATREG (categorical regression with optimal scaling using alternating least squares) quantifies categorical variables using optimal scaling, resulting in an optimal linear regression equation for the transformed variables. The variables can be given mixed optimal scaling levels, and no distributional assumptions about the variables are made.

### **Options**

**Transformation Type.** You can specify the transformation type (spline ordinal, spline nominal, ordinal, nominal, or numerical) at which you want to analyze each variable.

**Discretization.** You can use the `DISCRETIZATION` subcommand to discretize fractional-value variables or to recode categorical variables.

**Initial Configuration.** You can specify the kind of initial configuration through the `INITIAL` subcommand. Also, multiple systematic starts or fixed signs for the regression coefficients can be specified through this subcommand.

**Tuning the Algorithm.** You can control the values of algorithm-tuning parameters with the `MAXITER` and `CRITITER` subcommands.

**Regularized regression.** You can specify one of three methods for regularized regression: Ridge regression, the Lasso, or the Elastic Net.

**Resampling.** You can specify cross validation or the .632 bootstrap for estimation of prediction error.

**Missing Data.** You can specify the treatment of missing data with the `MISSING` subcommand.

**Optional Output.** You can request optional output through the `PRINT` subcommand.

**Transformation Plot per Variable.** You can request a plot per variable of its quantification against the category numbers.

**Residual Plot per Variable.** You can request an overlay plot per variable of the residuals and the weighted quantification against the category numbers.

**Ridge, Lasso, or Elastic Net plot.** You can request a plot of the regularized coefficients paths. For the Elastic Net, the plots for all values of the Ridge penalty can be requested, or plots for specified values of the Ridge penalty.

**Writing External Data.** You can write the transformed data (category numbers replaced with optimal quantifications) to an outfile for use in further analyses. You can also write the discretized data to an outfile.

**Saving Variables.** You can save the transformed variables, the predicted values, and/or the residuals in the working data file.

### ***Basic Specification***

The basic specification is the command `CATREG` with the `VARIABLES` and `ANALYSIS` subcommands.

### ***Syntax Rules***

- The `VARIABLES` and `ANALYSIS` subcommands must always appear, and the `VARIABLES` subcommand must be the first subcommand specified. The other subcommands, if specified, can be in any order.
- Variables specified in the `ANALYSIS` subcommand must be found in the `VARIABLES` subcommand.
- In the `ANALYSIS` subcommand, exactly one variable must be specified as a dependent variable and at least one variable must be specified as an independent variable after the keyword `WITH`.
- The word `WITH` is reserved as a keyword in the `CATREG` procedure. Thus, it may not be a variable name in `CATREG`. Also, the word `TO` is a reserved word.

### ***Operations***

- If a subcommand is specified more than once, the last one is executed but with a syntax warning. Note this is true also for the `VARIABLES` and `ANALYSIS` subcommands.

### ***Limitations***

- If more than one dependent variable is specified in the `ANALYSIS` subcommand, `CATREG` is not executed.

- CATREG operates on category indicator variables. The category indicators should be positive integers. You can use the `DISCRETIZATION` subcommand to convert fractional-value variables and string variables into positive integers. If `DISCRETIZATION` is not specified, fractional-value variables are automatically converted into positive integers by grouping them into seven categories with a close to normal distribution and string variables are automatically converted into positive integers by ranking.
- In addition to system missing values and user defined missing values, CATREG treats category indicator values less than 1 as missing. If one of the values of a categorical variable has been coded 0 or some negative value and you want to treat it as a valid category, use the `COMPUTE` command to add a constant to the values of that variable such that the lowest value will be 1. You can also use the `RANKING` option of the `DISCRETIZATION` subcommand for this purpose, except for variables you want to treat as numerical, since the characteristic of equal intervals in the data will not be maintained.
- There must be at least three valid cases.
- The number of valid cases must be greater than the number of independent variables plus 1.
- The maximum number of independent variables is 200.
- Split-File has no implications for CATREG.

## Examples

```

CATREG VARIABLES = TEST1 TEST3 TEST2 TEST4 TEST5 TEST6
  TEST7 TO TEST9 STATUS01 STATUS02
  /ANALYSIS TEST4 (LEVEL=NUME)
  WITH TEST1 TO TEST2 (LEVEL=SPORD DEGREE=1 INKNOT=3)
  TEST5 TEST7 (LEVEL=SPNOM) TEST8 (LEVEL=ORDI)
  STATUS01 STATUS02 (LEVEL=NOMI)
  /DISCRETIZATION = TEST1 (GROUPING NCAT=5 DISTR=UNIFORM)
  TEST5 (GROUPING) TEST7 (MULTIPLYING)

  /INITIAL = RANDOM
  /MAXITER = 100
  /CRITITER = .000001
  /RESAMPLE BOOTSTRAP (100)
  /MISSING = MODEIMPU
  /PRINT = R COEFF DESCRIP ANOVA QUANT(TEST1 TO TEST2 STATUS01 STATUS02)
  /PLOT = TRANS (TEST2 TO TEST7 TEST4)
  /SAVE
  /OUTFILE = '/data/qdata.sav' .

```

- `VARIABLES` defines variables. The keyword `TO` refers to the order of the variables in the working data file.
- The `ANALYSIS` subcommand defines variables used in the analysis. It is specified that `TEST4` is the dependent variable, with optimal scaling level numerical and that the variables `TEST1`, `TEST2`, `TEST3`, `TEST5`, `TEST7`, `TEST8`, `STATUS01`, and `STATUS02` are the independent variables to be used in the analysis. (The keyword `TO` refers to the order of the variables in the `VARIABLES` subcommand.) The optimal scaling level for `TEST1`, `TEST2`, and `TEST3` is spline ordinal; for `TEST5` and `TEST7`, spline nominal; for `TEST8`, ordinal; and for `STATUS01` and `STATUS02`, nominal. The splines for `TEST1` and `TEST2` have degree 1 and three interior knots, and the splines for `TEST5` and `TEST7` have degree 2 and two interior knots (default because unspecified).

- DISCRETIZATION specifies that *TEST5* and *TEST7*, which are fractional-value variables, are discretized: *TEST5* by recoding into seven categories with a normal distribution (default because unspecified) and *TEST7* by “multiplying.” *TEST1*, which is a categorical variable, is recoded into five categories with a close-to-uniform distribution.
- Because there are nominal variables, a random initial solution is requested by the INITIAL subcommand.
- MAXITER specifies the maximum number of iterations to be 100. This is the default, so this subcommand could be omitted here.
- CRITITER sets the convergence criterion to a value smaller than the default value.
- To include cases with missing values, the MISSING subcommand specifies that for each variable, missing values are replaced with the most frequent category (the mode).
- RESAMPLE specifies the .632 bootstrap for estimation of the prediction error using 100 bootstrap samples (in stead of the default of 50).
- PRINT specifies the correlations, the coefficients, the descriptive statistics for all variables, the ANOVA table, the category quantifications for variables *TEST1*, *TEST2*, *TEST3*, *STATUS01*, and *STATUS02*, and the transformed data list of all cases.
- PLOT is used to request quantification plots for the variables *TEST2*, *TEST5*, *TEST7*, and *TEST4*.
- The SAVE subcommand adds the transformed variables to the working data file. The names of these new variables are *TRANS1\_1*, ..., *TRANS9\_1*.
- The OUTFILE subcommand writes the transformed data to a data file called *qdata.sav* in the directory */data*.

### Example: Multiple Systematic Starts

```
CATREG
...
/INITIAL MULTISTART(ALL) ('c:\data\startsigns.sav')
/PRINT = R COEFF DESCRIP ANOVA QUANT(TEST1 TO TEST2 STATUS01 STATUS02)
/PLOT = TRANS (TEST2 TO TEST7 TEST4)
/SAVE TRDATA PRED RES
/OUTFILE = TRDATA('c:\data\qdata.sav') DISCRDATA('c:\data\discr.sav')
```

- Because the ordinal and spline ordinal scaling levels are specified for some variables, there is chance of obtaining a suboptimal solution when applying the numerical or random initial solution. To ascertain obtaining the optimal solution, all multiple systematic starts are used. Using all systematic starts is feasible here because the number of variables with (spline) ordinal scaling is only 3; then the number of all starts is 2 to the power of 3 is 8. With a larger number of variables with (spline) ordinal scaling level, a reduced number of starts is recommended, which can be requested by specifying `/INITIAL MULTISTART(value)`.
- The specifications at the PRINT, PLOT, SAVE, and OUTFILE subcommands will be applied to the optimal solution.

### Example: Fixing Initial Signs for Regression Coefficients

```
CATREG
...
```



```
/INITIAL FIXSIGNS (63) ('c:\data\startsigns.sav')
```

- The `INITIAL` subcommand specifies using a specific set of fixed signs for the regression coefficients. The signs are in the file *startsigns.sav* in the directory *c:\data*. This file was created by a previous run of `CATREG` with keyword `MULTISTART` at the `INITIAL` subcommand (see previous example). The signs of start number 63 are specified to be used.

### **Example: Elastic Net Regularization**

```
CATREG
```

```
...
/REGULARIZATION ENET (.5 2.5 .25) (.01 3.8 .05) ('c:\data\regu_enet.sav')
/PRINT = REGU R COEFF DESCRIP ANOVA QUANT(TEST1 TO TEST2)
/PLOT = REGU (.75 1.5) TRANS (TEST2 TO TEST7 TEST4)
/SAVE TRDATA PRED RES
/OUTFILE = TRDATA('c:\data\qdata.sav') DISCRDATA('c:\data\discr.sav').
```

- `REGULARIZATION` specifies application of Elastic Net regularization, with start value of the Lasso penalty 0.01, stop value 3.8, and increment 0.05, resulting in 76 regularized models, with Lasso penalty values 0.01, 0.06, ..., 3.76. To each of these 76 Lasso models 10 Ridge penalties are applied (0.5, 0.75, ..., 2.5), resulting in  $76 \times 10 = 760$  Elastic Net models.
- `PRINT` specifies displaying a table with the penalty values, R-squared, and the regression coefficients for each regularized model. The contents of this table is written to a data file called *regu\_enet.sav* in the directory *c:\data*.
- The `PLOT` subcommand requests two Elastic Net plots: a Lasso plot with a fixed Ridge penalty of 0.75 and a Lasso plot with a fixed Ridge penalty of 1.50. Any other keywords than `REGU` at the `PLOT` subcommand are ignored.
- Other specifications than `REGU` at the `PRINT` and `PLOT` subcommands, the `SAVE` subcommand and the `TRDATA` keyword at the `OUTFILE` subcommand are ignored.

### **Example: Elastic Net Regularization with Crossvalidation Resampling**

```
CATREG
```

```
...
/REGULARIZATION ENET (.5 2.5 .25) (.01 3.8 .05) ('c:\data\regu_enet.sav')
/RESAMPLE CROSSVAL (5)
/PRINT = REGU R COEFF DESCRIP ANOVA
/PLOT = REGU (.75 1.5) TRANS (TEST2 TO TEST7 TEST4)
/SAVE TRDATA PRED RES
/OUTFILE = TRDATA('c:\data\qdata.sav') DISCRDATA('c:\data\discr.sav').
```

- `REGULARIZATION` is the same as in the previous example.
- The `RESAMPLE` subcommand specifies 5-fold cross-validation to estimate the prediction error for each of the 760 Elastic Net models.
- `PRINT` specifies displaying a table with the penalty values, R-squared, the regression coefficients, and the estimated prediction error for each regularized model. The contents of this table is written to a data file called *regu\_enet.sav* in the directory *c:\data*.
- The specification at the `PLOT` subcommand result in the same plots as in the previous example.
- The other specifications at the `PRINT` and `PLOT` subcommands, and the `SAVE` and `OUTFILE` specifications will be applied to the model with lowest prediction error.

**Example: Obtaining a Specific Elastic Net Model**

```

CATREG
...
/REGULARIZATION ENET (1.25 1.25 0) (.46 .46 0)
/PRINT = R COEFF DESCRIP ANOVA QUANT(TEST1 TO TEST2 STATUS01 STATUS02)
/PLOT = TRANS (TEST2 TO TEST7 TEST4)
/SAVE TRDATA PRED RES
/OUTFILE = TRDATA('c:\data\qdata.sav') DISCRDATA('c:\data\discr.sav').

```

- REGULARIZATION is specified here (stop value equal to start value, increment zero) to obtain output for a specific Elastic Net model: the model with penalty values 1.25 (Ridge) and .46 (Lasso).

**VARIABLES Subcommand**

VARIABLES specifies the variables that may be analyzed in the current CATREG procedure.

- The VARIABLES subcommand is required and precedes all other subcommands.
- The keyword TO on the VARIABLES subcommand refers to the order of variables in the working data file. (Note that this behavior of TO is different from that in the `indvarlist` on the ANALYSIS subcommand.)

**ANALYSIS Subcommand**

ANALYSIS specifies the dependent variable and the independent variables following the keyword WITH.

- All the variables on ANALYSIS must be specified on the VARIABLES subcommand.
- The ANALYSIS subcommand is required and follows the VARIABLES subcommand.
- The first variable list contains exactly one variable as the dependent variable, while the second variable list following WITH contains at least one variable as an independent variable. Each variable may have at most one keyword in parentheses indicating the transformation type of the variable.
- The keyword TO in the independent variable list honors the order of variables on the VARIABLES subcommand.
- Optimal scaling levels are indicated by the keyword LEVEL in parentheses following the variable or variable list.

**LEVEL**      *Specifies the optimal scaling level.*

## **LEVEL Keyword**

The following keywords are used to indicate the optimal scaling level:

<b>SPORD</b>	<i>Spline ordinal (monotonic)</i> . This is the default for a variable listed without any optimal scaling level, for example, one without <code>LEVEL</code> in the parentheses after it or with <code>LEVEL</code> without a specification. Categories are treated as ordered. The order of the categories of the observed variable is preserved in the optimally scaled variable. Categories will be on a straight line through the origin. The resulting transformation is a smooth nondecreasing piecewise polynomial of the chosen degree. The pieces are specified by the number and the placement of the interior knots.
<b>SPNOM</b>	<i>Spline nominal (non-monotonic)</i> . Categories are treated as unordered. Objects in the same category obtain the same quantification. Categories will be on a straight line through the origin. The resulting transformation is a smooth piecewise polynomial of the chosen degree. The pieces are specified by the number and the placement of the interior knots.
<b>ORDI</b>	<i>Ordinal</i> . Categories are treated as ordered. The order of the categories of the observed variable is preserved in the optimally scaled variable. Categories will be on a straight line through the origin. The resulting transformation fits better than <code>SPORD</code> transformation, but is less smooth.
<b>NOMI</b>	<i>Nominal</i> . Categories are treated as unordered. Objects in the same category obtain the same quantification. Categories will be on a straight line through the origin. The resulting transformation fits better than <code>SPNOM</code> transformation, but is less smooth.
<b>NUME</b>	<i>Numerical</i> . Categories are treated as equally spaced (interval level). The order of the categories and the differences between category numbers of the observed variables are preserved in the optimally scaled variable. Categories will be on a straight line through the origin. When all variables are scaled at the numerical level, the <code>CATREG</code> analysis is analogous to standard multiple regression analysis.

## **SPORD and SPNOM Keywords**

The following keywords are used with `SPORD` and `SPNOM` :

<b>DEGREE</b>	<i>The degree of the polynomial</i> . If <code>DEGREE</code> is not specified the degree is assumed to be 2.
<b>INKNOT</b>	<i>The number of the interior knots</i> . If <code>INKNOT</code> is not specified the number of interior knots is assumed to be 2.

## **DISCRETIZATION Subcommand**

`DISCRETIZATION` specifies fractional-value variables that you want to discretize. Also, you can use `DISCRETIZATION` for ranking or for two ways of recoding categorical variables.

- A string variable's values are always converted into positive integers by assigning category indicators according to the ascending alphanumeric order. `DISCRETIZATION` for string variables applies to these integers.
- When the `DISCRETIZATION` subcommand is omitted, or when the `DISCRETIZATION` subcommand is used without a varlist, fractional-value variables are converted into positive integers by grouping them into seven categories (or into the number of distinct values of the variable if this number is less than 7) with a close to normal distribution.

- When no specification is given for variables in a varlist following `DISCRETIZATION`, these variables are grouped into seven categories with a close-to-normal distribution.
- In `CATREG`, a system-missing value, user-defined missing values, and values less than 1 are considered to be missing values (see next section). However, in discretizing a variable, values less than 1 are considered to be valid values, and are thus included in the discretization process. System-missing values and user-defined missing values are excluded.

<b>GROUPING</b>	<i>Recode into the specified number of categories.</i>
<b>RANKING</b>	<i>Rank cases. Rank 1 is assigned to the case with the smallest value on the variable.</i>
<b>MULTIPLYING</b>	<i>Multiplying the standardized values (z-scores) of a fractional-value variable by 10, rounding, and adding a value such that the lowest value is 1.</i>

### ***GROUPING Keyword***

<b>NCAT</b>	<i>Recode into ncat categories. When <code>NCAT</code> is not specified, the number of categories is set to 7 (or the number of distinct values of the variable if this number is less than 7). The valid range is from 2 to 36. You may either specify a number of categories or use the keyword <code>DISTR</code>.</i>
<b>EQINTV</b>	<i>Recode intervals of equal size into categories. The interval size must be specified (there is no default value). The resulting number of categories depends on the interval size.</i>

### ***DISTR Keyword***

`DISTR` has the following keywords:

<b>NORMAL</b>	<i>Normal distribution. This is the default when <code>DISTR</code> is not specified.</i>
<b>UNIFORM</b>	<i>Uniform distribution.</i>

### ***MISSING Subcommand***

In `CATREG`, we consider a system missing value, user defined missing values, and values less than 1 as missing values. However, in discretizing a variable (see previous section), values less than 1 are considered as valid values. The `MISSING` subcommand allows you to indicate how to handle missing values for each variable.

<b>LISTWISE</b>	<i>Exclude cases with missing values on the specified variable(s). The cases used in the analysis are cases without missing values on the variable(s) specified. This is the default applied to all variables, when the <code>MISSING</code> subcommand is omitted or is specified without variable names or keywords. Also, any variable that is not included in the subcommand gets this specification.</i>
<b>MODEIMPU</b>	<i>Impute missing value with mode. All cases are included and the imputations are treated as valid observations for a given variable. When there are multiple modes, the smallest mode is used.</i>
<b>EXTRACAT</b>	<i>Impute missing values on a variable with an extra category indicator. This implies that objects with a missing value are considered to belong to the same (extra) category. This category is treated as nominal, regardless of the optimal scaling level of the variable.</i>

- The ALL keyword may be used to indicate all variables. If it is used, it must be the only variable specification.
- A mode or extra-category imputation is done before listwise deletion.

## ***SUPPLEMENTARY Subcommand***

The SUPPLEMENTARY subcommand specifies the objects that you want to treat as supplementary. You cannot weight supplementary objects (specified weights are ignored). This subcommand can be used to specify test cases.

**OBJECT**            *Supplementary objects.* Objects that you want to treat as supplementary are indicated with an object number list in parentheses following OBJECT. The keyword TO is allowed—for example, OBJECT(1 TO 1 3 5 TO 9).

- Supplementary object are excluded from the analysis. The quantifications resulting from the analysis for the active objects are applied to the categories of supplementary objects, and predicted and residual values for supplementary objects are provided.
- If a supplementary object has a category that does not occur in the active data, the following strategies are applied:

If the variable on which the non-occurring category occurs has a numeric or spline scaling level, and the non-occurring category lies within the range of categories in the active data, then interpolation is applied.

If the variable has numeric scaling level and the non-occurring category lies outside the range of categories in the active data, then extrapolation is applied.

Otherwise the case is excluded.

- Excluded cases are represented by a dot (the sysmis symbol) on every saved variable.

## ***INITIAL Subcommand***

INITIAL specifies the method used to compute the initial value/configuration.

- The specification on INITIAL is keyword NUMERICAL, RANDOM, MULTISTART or FIXSIGNS. If INITIAL is not specified, NUMERICAL is the default.

### **NUMERICAL**

*Treat all variables as numerical.* This is usually best to use when there are only numerical and/or ordinal variables.

### **RANDOM**

*Provide a random initial value.* This should be used only when there is at least one nominal variable.

**MULTISTART(integer|ALL)('savfile'|'dataset')**

*Multiple Systematic Starts.* Multiple final solutions are computed and the best solution is selected. For each solution the same initial values are used, but with different signs for the regression coefficients of variables with ordinal or spline ordinal scaling level. This option is only applicable when there is at least one variable with ordinal or spline ordinal scaling level. With these scaling levels, the CATREG algorithm can result in a suboptimal solution. The optimal solution is always found when multiple systematic starts (using all possible sign patterns for the regression coefficients) are applied. You can specify ALL in parentheses following the keyword to ascertain obtaining the optimal solution. However, the number of all possible sign patterns for the regression coefficients is 2 to the power of  $q$ , where  $q$  is the number of variables with ordinal or spline ordinal scaling level. So, the number of all possible sign patterns rapidly increases with increasing  $q$ .

When  $q$  is large, a reduced number of multiple systematic starts can be requested by specifying a value in parentheses following the keyword. This option selects a reduced number of sign patterns by applying a hierarchical strategy combined with a percentage of loss criterion. The value to specify is the threshold for the percentage of loss of variance that a variable suffers due to ordinal restriction. Specify a non-negative value less than or equal to 100. A variable for which this percentage is below the specified threshold, is not allowed to have a negative sign. So, specifying a threshold value excludes sign patterns in which variables with a loss of variance percentage below the specified threshold have a negative sign. Thus, the higher the threshold, the more sign patterns will be excluded. With this option, obtaining the optimal solution is not guaranteed, but the chance of obtaining a suboptimal solution is diminished. Also, if with the reduced number of starts the optimal solution is not found, the chance that the suboptimal solution is much different from the optimal solution is diminished. Note that there is a trade-off between the chance of obtaining a suboptimal solution and the number of starts: a higher threshold results in more reduction of the number of starts, but a higher chance of obtaining a suboptimal solution.

When this keyword is used, a dataset name or filename in parentheses must be specified. The signs of the regression coefficients for each start will be written to this file.

To give an impression of computing time: when  $q$  is 15, and all variables have seven categories (in CATREG, CPU time depends upon the number of categories, not upon the number of cases), the number of all starts is 32768, which requires 4 minutes on a 2.2 Ghz computer. When  $q$  is 20, the number of all starts is 1048576, requiring 4.5 hours, and when  $q$  is 21, the number of all starts is 2097152, requiring 11.5 hours.

#### **FIXSIGNS(integer startnumber)('savfile' | 'dataset')**

Use fixed signs for the regression coefficients. The signs (indicated by 1 and -1) need to be in (a row of) the specified dataset or file. The integer-valued startnumber to specify is the case number of the row in this file that contains the signs that are to be used. If in a previous run MULTISTART was specified, a file containing the signs for each start was created and can be used here.

## **MAXITER Subcommand**

MAXITER specifies the maximum number of iterations CATREG can go through in its computations. Note that the output starts from the iteration number 0, which is the initial value before any iteration, when INITIAL = NUMERICAL is in effect.

- If MAXITER is not specified, CATREG will iterate up to 100 times.
- The specification on MAXITER is a positive integer indicating the maximum number of iterations. There is no uniquely predetermined (hard coded) maximum for the value that can be used.

## **CRITITER Subcommand**

CRITITER specifies a convergence criterion value. CATREG stops iterating if the difference in fit between the last two iterations is less than the CRITITER value.

- If CRITITER is not specified, the convergence value is 0.00001.
- The specification on CRITITER is any value less than or equal to 0.1 and greater than or equal to 0.000001. (Values less than the lower bound might seriously affect performance. Therefore, they are not supported.)

## **REGULARIZATION Subcommand**

REGULARIZATION specifies the method for regularized regression.

The specification on REGULARIZATION is keyword NONE, RIDGE, LASSO or ENET. If REGULARIZATION is not specified, NONE is the default. Also, a dataset name or filename must be specified. The statistics and coefficients and, if applicable, the estimated prediction error, for all regularized models will be written to this file.

### **NONE**

*No regularization.*

### **RIDGE(start value, stop value, increment)('savfile'|'dataset')**

*Ridge Regression.* A value list in parentheses following the keyword should be given. The first value specifies the start value of the penalty parameter, the second value the stop value, and the third value specifies the increment.

### **LASSO(start value, stop value, increment)('savfile'|'dataset')**

*LASSO (Least Absolute Shrinkage and Selection Operator).* A value list in parentheses following the keyword should be given. The first value specifies the start value of the penalty parameter, the second value the stop value, and the third value specifies the increment.

### **ENET(start, stop , incr)(start, stop , incr)('savfile'|'dataset')**

*Elastic Net.* Two value lists in parentheses following the keyword should be given. The first list specifies the start, stop, and increment values for the Ridge penalty, the second list specifies the start, stop, and increment values for the Lasso penalty.

- If a REGULARIZATION method is specified without specification of a resample method at the RESAMPLE subcommand or specification of test cases at the SUPPLEMENTARY subcommand, any other keywords than REGU at the PRINT and PLOT subcommands are ignored. Also, the SAVE subcommand, and the TRDATA keyword at the OUTFILE subcommand are ignored.
- If a resample method is specified at the RESAMPLE subcommand, or if test cases are specified at the SUPPLEMENTARY subcommand, specified PRINT, PLOT, SAVE, and OUTFILE output will be given for the model with lowest prediction error or with lowest Test MSE.
- Output of an analysis with a specific value of the penalty parameter(s) is obtained by setting the start value(s) to specific penalty value(s), the stop value(s) equal to the start value(s) and the increment value(s) to 0.

## RESAMPLE Subcommand

RESAMPLE specifies the resampling method used for estimation of the prediction error.

The specification on RESAMPLE is keyword NONE, CROSSVAL or BOOTSTRAP. If RESAMPLE is not specified, NONE is the default.

### NONE

*No resampling.*

### CROSSVAL(integer)

*Cross-validation.* The keyword can be followed by a positive integer in parentheses specifying the number of folds. If this value is not specified, 10-fold cross-validation is used.

### BOOTSTRAP(integer)

*.632 Bootstrap.* The keyword can be followed by a positive integer in parentheses specifying the number of bootstrap samples. If this value is not specified, 50 bootstrap samples are used.

## PRINT Subcommand

The PRINT subcommand controls the display of output. The output of the CATREG procedure is always based on the transformed variables. However, the correlations of the original predictor variables can be requested as well by the keyword Ocorr. The default keywords are R, COEFF, DESCRIP, and ANOVA. That is, the four keywords are in effect when the PRINT subcommand is omitted or when the PRINT subcommand is given without any keyword. If a keyword is duplicated or it encounters a contradicting keyword, such as /PRINT = R R NONE, then the last one silently becomes effective.

- The REGU keyword is only applicable if a REGULARIZATION method is specified.

### R

*Multiple R.* Includes  $R^2$ , adjusted  $R^2$ , and adjusted  $R^2$  taking the optimal scaling into account.

### COEFF

*Standardized regression coefficients (beta).* This option gives three tables: a Coefficients table that includes betas, standard error of the betas,  $t$  values, and significance; a Coefficients-Optimal Scaling table, with the standard error of the betas taking the optimal scaling degrees of freedom into account; and a table with the zero-order, part, and partial correlation, Pratt's relative importance measure for the transformed predictors, and the tolerance before and after transformation. If the tolerance for a transformed predictor is lower than the default tolerance value in the Regression procedure (0.0001) but higher than  $10E-12$ , this is reported in an annotation. If the tolerance is lower than  $10E-12$ , then the COEFF computation for this variable is not done and this is reported in an annotation. Note that the regression model includes the intercept coefficient but that its estimate does not exist because the coefficients are standardized.

### DESCRIP(varlist)

*Descriptive statistics (frequencies, missing values, and mode).* The variables in the varlist must be specified on the VARIABLES subcommand but need not appear on the ANALYSIS subcommand. If DESCRIP is not followed by a varlist, Descriptives tables are displayed for all of the variables in the variable list on the ANALYSIS subcommand.



<b>HISTORY</b>	<i>History of iterations.</i> For each iteration, including the starting values for the algorithm, the multiple $R$ and the regression error (square root of $(1 - \text{multiple } R^2)$ ) are shown. The increase in multiple $R$ is listed from the first iteration.
<b>ANOVA</b>	<i>Analysis-of-variance tables.</i> This option includes regression and residual sums of squares, mean squares, and $F$ . This options gives two ANOVA tables: one with degrees of freedom for the regression equal to the number of predictor variables and one with degrees of freedom for the regression taking the optimal scaling into account.
<b>CORR</b>	<i>Correlations of the transformed predictors.</i>
<b>OCORR</b>	<i>Correlations of the original predictors.</i>
<b>QUANT(varlist)</b>	<i>Category quantifications.</i> Any variable in the ANALYSIS subcommand may be specified in parentheses after QUANT. If QUANT is not followed by a varlist, Quantification tables are displayed for all variables in the variable list on the ANALYSIS subcommand.
<b>REGU</b>	<i>Penalty values, R-squared, and the regression coefficients for each regularized model, and, if a RESAMPLE method is specified or if supplementary objects (tests cases) are specified, the prediction error or Test MSE.</i>
<b>NONE</b>	<i>No PRINT output is shown.</i> This is to suppress the default PRINT output.

- The keyword TO in a variable list can be used only with variables that are in the ANALYSIS subcommand, and TO applies only to the order of the variables in the ANALYSIS subcommand. For variables that are in the VARIABLES subcommand but not in the ANALYSIS subcommand, the keyword TO cannot be used. For example, if /VARIABLES = v1 TO v5 and /ANALYSIS is v2 v1 v4, then /PRINT QUANT(v1 TO v4) will give two quantification plots, one for v1 and one for v4. (/PRINT QUANT(v1 TO v4 v2 v3 v5) will give quantification tables for v1, v2, v3, v4, and v5.)

## ***PLOT Subcommand***

The PLOT subcommand controls the display of plots.

- The REGU keyword is only applicable if a REGULARIZATION method is specified.
- In this subcommand, if no plot keyword is given, then no plot is created. Further, if the variable list following the plot keyword is empty, then no plot is created, either.
- All of the variables to be plotted must be specified in the ANALYSIS subcommand. Further, for the residual plots, the variables must be independent variables.

<b>TRANS(varlist)(l)</b>	<i>Transformation plots (optimal category quantifications against category indicators).</i> A list of variables must come from the ANALYSIS variable list and must be given in parentheses following the keyword. Further, the user can specify an optional parameter l in parentheses after the variable list in order to control the global upper boundary of category label lengths in the plot. Note that this boundary is applied uniformly to all transformation plots.
<b>RESID(varlist)(l)</b>	<i>Residual plots (residuals when the dependent variable is predicted from all predictor variables in the analysis except the predictor variable in varlist, against category indicators, and the optimal category quantifications multiplied with beta against category indicators).</i> A list of variables must come from the ANALYSIS variable list's independent variables and must be given in parentheses following the keyword. Further, the user can specify an optional parameter l in parentheses after

the variable list in order to control the global upper boundary of category label lengths in the plot. Note that this boundary is applied uniformly to all residual plots.

**REGU(valuelist)**

*Ridge, Lasso, or Elastic Net plot(s), depending on the regularization method specified at the REGULARIZATION subcommand. A value or valuelist of Ridge penalties must be given in parentheses following the keyword if the regularization method is Elastic Net. The Elastic Net method results in multiple plots: a Lasso plot for each value of the Ridge penalty. To obtain all Elastic Net plots, the keyword ALL in stead of a valuelist can be used.*

- The category label length parameter (1) can take any non-negative integer less than or equal to 60. If 1 = 0, values instead of value labels are displayed to indicate the categories on the  $x$  axis in the plot. If 1 is not specified, CATREG assumes that each value label at its full length is displayed as a plot's category label. If 1 is an integer larger than 60, then we reset it to 60 but do not issue a warning.
- If a positive value of 1 is given but if some or all of the values do not have value labels, then for those values, the values themselves are used as the category labels.
- The keyword TO in a variable list can be used only with variables that are in the ANALYSIS subcommand, and TO applies only to the order of the variables in the ANALYSIS subcommand. For variables that are in the VARIABLES subcommand but not in the ANALYSIS subcommand, the keyword TO cannot be used. For example, if /VARIABLES = v1 TO v5 and /ANALYSIS is v2 v1 v4, then /PLOT TRANS(v1 TO v4) will give two transformation plots, one for v1 and for v4. (/PLOT TRANS(v1 TO v4 v2 v3 v5) will give transformation plots for v1, v2, v3, v4, and v5.)

## SAVE Subcommand

The SAVE subcommand is used to add the transformed variables (category indicators replaced with optimal quantifications), the predicted values, and the residuals to the working data file.

Excluded cases are represented by a dot (the sysmis symbol) on every saved variable.

<b>TRDATA</b>	<i>Transformed variables.</i>
<b>PRED</b>	<i>Predicted values.</i>
<b>RES</b>	<i>Residuals.</i>

- A variable rootname can be specified with each of the keywords. Only one rootname can be specified with each keyword, and it can contain up to five characters (if more than one rootname is specified with a keyword, the first rootname is used; if a rootname contains more than five characters, the first five characters are used at most). If a rootname is not specified, the default rootnames (TRA, PRE, and RES) are used.
- CATREG adds two numbers separated by an underscore ( ) to the rootname. The formula is *ROOTNAMEk\_n*, where  $k$  increments from 1 to identify the source variable names by using the source variables' position numbers in the ANALYSIS subcommand (that is, the dependent variable has the position number 1, and the independent variables have the position numbers 2, 3, ..., etc., as they are listed), and  $n$  increments from 1 to identify the CATREG procedures with the successfully executed SAVE subcommands for a given data file in a continuous

session. For example, with two predictor variables specified on `ANALYSIS`, the first set of default names for the transformed data, if they do not exist in the data file, would be `TRA1_1` for the dependent variable, and `TRA2_1`, `TRA3_1` for the predictor variables. The next set of default names, if they do not exist in the data file, would be `TRA1_2`, `TRA2_2`, `TRA3_2`. However, if, for example, `TRA1_2` already exists in the data file, then the default names should be attempted as `TRA1_3`, `TRA2_3`, `TRA3_3`—that is, the last number increments to the next available integer.

- Variable labels are created automatically. (They are shown in the Procedure Information Table (the Notes table) and can also be displayed in the Data Editor window.)

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand is used to write the discretized data and/or the transformed data (category indicators replaced with optimal quantifications) to a data file or previously declared data set name. Excluded cases are represented by a dot (the `sysmis` symbol) on every saved variable.

<code>DISCRDATA('savfile' dataset')</code>	<i>Discretized data.</i>
<code>TRDATA('savfile' dataset')</code>	<i>Transformed variables.</i>

- Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. Data sets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files.
- An active data set, in principle, should not be replaced by this subcommand, and the asterisk (\*) file specification is not supported. This strategy also prevents the `OUTFILE` interference with the `SAVE` subcommand.

# CCF

```
CCF VARIABLES= series names [WITH series names]

[/DIFF={1}]
      {n}

[/SDIFF={1}]
      {n}

[/PERIOD=n]

[/{NOLOG**}]
      {LN      }

[/SEASONAL]

[/MXCROSS={7**}]
      {n      }

[/APPLY[='model name']]
```

**\*\*Default if the subcommand is omitted and there is no corresponding specification on the TSET command.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
CCF VARIABLES = VARX VARY.
```

## **Overview**

CCF displays and plots the cross-correlation functions of two or more time series. You can also display and plot the cross-correlations of transformed series by requesting natural log and differencing transformations within the procedure.

### **Options**

**Modifying the Series.** You can request a natural log transformation of the series using the LN subcommand and seasonal and nonseasonal differencing to any degree using the SDIFF and DIFF subcommands. With seasonal differencing, you can also specify the periodicity on the PERIOD subcommand.

**Statistical Display.** You can control which series are paired by using the keyword WITH. You can specify the range of lags for which you want values displayed and plotted with the MXCROSS subcommand, overriding the maximum specified on TSET. You can also display and plot values at periodic lags only using the SEASONAL subcommand.

### **Basic Specification**

The basic specification is two or more series names. By default, CCF automatically displays the cross-correlation coefficient and standard error for the negative lags (second series leading), the positive lags (first series leading), and the 0 lag for all possible pair combinations in the series list. It also plots the cross-correlations and marks the bounds of two standard errors on the plot. By default, CCF displays and plots values up to 7 lags (lags  $-7$  to  $+7$ ), or the range specified on TSET.

### **Subcommand Order**

- Subcommands can be specified in any order.

### **Syntax Rules**

- The VARIABLES subcommand can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each one is executed.

### **Operations**

- Subcommand specifications apply to all series named on the CCF command.
- If the LN subcommand is specified, any differencing requested on that CCF command is done on the log-transformed series.
- Confidence limits are displayed in the plot, marking the bounds of two standard errors at each lag.

### **Limitations**

- A maximum of 1 VARIABLES subcommand. There is no limit on the number of series named on the list.

## **Example**

```
CCF VARIABLES = VARX VARY
  /LN
  /DIFF=1
  /SDIFF=1
  /PERIOD=12
  /MXCROSS=25.
```

- This example produces a plot of the cross-correlation function for *VARX* and *VARY* after a natural log transformation, differencing, and seasonal differencing have been applied to both series. Along with the plot, the cross-correlation coefficients and standard errors are displayed for each lag.
- LN transforms the data using the natural logarithm (base  $e$ ) of each series.
- DIFF differences each series once.
- SDIFF and PERIOD apply one degree of seasonal differencing with a periodicity of 12.
- MXCROSS specifies 25 for the maximum range of positive and negative lags for which output is to be produced (lags  $-25$  to  $+25$ ).

## VARIABLES Subcommand

VARIABLES specifies the series to be plotted and is the only required subcommand.

- The minimum VARIABLES specification is a pair of series names.
- If you do not use the keyword WITH, each series is paired with every other series in the list.
- If you specify the keyword WITH, every series named before WITH is paired with every series named after WITH.

### Example

```
CCF VARIABLES=VARA VARB WITH VARC VARD.
```

- This example displays and plots the cross-correlation functions for the following pairs of series: *VARA* with *VARC*, *VARA* with *VARD*, *VARB* with *VARC*, and *VARB* with *VARD*.
- *VARA* is not paired with *VARB*, and *VARC* is not paired with *VARD*.

## DIFF Subcommand

DIFF specifies the degree of differencing used to convert a nonstationary series to a stationary one with a constant mean and variance before obtaining cross-correlations.

- You can specify 0 or any positive integer on DIFF.
- If DIFF is specified without a value, the default is 1.
- The number of values used in the calculations decreases by 1 for each degree of differencing.

### Example

```
CCF VARIABLES = VARX VARY
/DIFF=1.
```

- This command differences series *VARX* and *VARY* before calculating and plotting the cross-correlation function.

## SDIFF Subcommand

If the series exhibits seasonal or periodic patterns, you can use SDIFF to seasonally difference the series before obtaining cross-correlations.

- The specification on SDIFF indicates the degree of seasonal differencing and can be 0 or any positive integer.
- If SDIFF is specified without a value, the degree of seasonal differencing defaults to 1.
- The number of seasons used in the calculations decreases by 1 for each degree of seasonal differencing.
- The length of the period used by SDIFF is specified on the PERIOD subcommand. If the PERIOD subcommand is not specified, the periodicity established on the TSET or DATE command is used (see the PERIOD subcommand).

**Example**

```
CCF VARIABLES = VAR01 WITH VAR02 VAR03
  /SDIFF=1.
```

- In this example, one degree of seasonal differencing using the periodicity established on the TSET or DATE command is applied to the three series.
- Two cross-correlation functions are then plotted, one for the pair *VAR01* and *VAR02*, and one for the pair *VAR01* and *VAR03*.

**PERIOD Subcommand**

PERIOD indicates the length of the period to be used by the SDIFF or SEASONAL subcommands.

- The specification on PERIOD indicates how many observations are in one period or season and can be any positive integer greater than 1.
- PERIOD is ignored if it is used without the SDIFF or SEASONAL subcommands.
- If PERIOD is not specified, the periodicity established on TSET PERIOD is in effect. If TSET PERIOD is not specified, the periodicity established on the DATE command is used. If periodicity was not established anywhere, the SDIFF and SEASONAL subcommands will not be executed.

**Example**

```
CCF VARIABLES = VARX WITH VARY
  /SDIFF=1
  /PERIOD=6.
```

- This command applies one degree of seasonal differencing with a periodicity of 6 to both series and computes and plots the cross-correlation function.

**LN and NOLOG Subcommands**

LN transforms the data using the natural logarithm (base  $e$ ) of each series and is used to remove varying amplitude over time. NOLOG indicates that the data should not be log transformed. NOLOG is the default.

- There are no additional specifications on LN or NOLOG.
- Only the last LN or NOLOG subcommand on a CCF command is executed.
- LN and NOLOG apply to all series named on the CCF command.
- If a natural log transformation is requested and any values in either series in a pair are less than or equal to 0, the CCF for that pair will not be produced because nonpositive values cannot be log transformed.
- NOLOG is generally used with an APPLY subcommand to turn off a previous LN specification.

**Example**

```
CCF VARIABLES = VAR01 VAR02
```

/LN.

- This command transforms the series *VAR01* and *VAR02* using the natural log before computing cross-correlations.

## **SEASONAL Subcommand**

Use `SEASONAL` to focus attention on the seasonal component by displaying and plotting cross-correlations at periodic lags only.

- There are no additional specifications on `SEASONAL`.
- If `SEASONAL` is specified, values are displayed and plotted at the periodic lags indicated on the `PERIOD` subcommand. If no `PERIOD` subcommand is specified, the periodicity first defaults to the `TSET PERIOD` specification and then to the `DATE` command periodicity. If periodicity is not established anywhere, `SEASONAL` is ignored (see the `PERIOD` subcommand).
- If `SEASONAL` is not used, cross-correlations for all lags up to the maximum are displayed and plotted.

### **Example**

```
CCF VARIABLES = VAR01 VAR02 VAR03
/SEASONAL.
```

- This command plots and displays cross-correlations at periodic lags.
- By default, the periodicity established on `TSET PERIOD` (or the `DATE` command) is used. If no periodicity is established, cross-correlations for all lags are displayed and plotted.

## **MXCROSS Subcommand**

`MXCROSS` specifies the maximum range of lags for a series.

- The specification on `MXCROSS` must be a positive integer.
- If `MXCROSS` is not specified, the default range is the value set on `TSET MXCROSS`. If `TSET MXCROSS` is not specified, the default is 7 (lags -7 to +7).
- The value specified on the `MXCROSS` subcommand overrides the value set on `TSET MXCROSS`.

### **Example**

```
CCF VARIABLES = VARX VARY
/MXCROSS=5.
```

- The maximum number of cross-correlations can range from lag -5 to lag +5.

## **APPLY Subcommand**

`APPLY` allows you to use a previously defined CCF model without having to repeat the specifications.



- The only specification on `APPLY` is the name of a previous model enclosed in single or double quotes. If a model name is not specified, the model specified on the previous `CCF` command is used.
- To change one or more model specifications, specify the subcommands of only those portions you want to change after the `APPLY` subcommand.
- If no series are specified on the command, the series that were originally specified with the model being applied are used.
- To change the series used with the model, enter new series names before or after the `APPLY` subcommand.

### Example

```
CCF VARIABLES = VARX VARY
  /LN
  /DIFF=1
  /MXCROSS=25.
CCF VARIABLES = VARX VARY
  /LN
  /DIFF=1
  /SDIFF=1
  /PERIOD=12
  /MXCROSS=25.
CCF VARIABLES = VARX VAR01
  /APPLY.
CCF VARIABLES = VARX VAR01
  /APPLY='MOD_1'.
```

- The first command displays and plots the cross-correlation function for *VARX* and *VARY* after each series is log transformed and differenced. The maximum range is set to 25 lags. This model is assigned the name *MOD\_1* as soon as the command is executed.
- The second command displays and plots the cross-correlation function for *VARX* and *VARY* after each series is log transformed, differenced, and seasonally differenced with a periodicity of 12. The maximum range is again set to 25 lags. This model is assigned the name *MOD\_2*.
- The third command requests the cross-correlation function for the series *VARX* and *VAR01* using the same model and the same range of lags as used for *MOD\_2*.
- The fourth command applies *MOD\_1* (from the first command) to the series *VARX* and *VAR01*.

## References

Box, G. E. P., and G. M. Jenkins. 1976. *Time series analysis: Forecasting and control*, Rev. ed. San Francisco: Holden-Day.

# CD

```
CD 'directory specification'.
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 13.0

- Command introduced.

## **Example**

```
CD '/main/sales/consumer_division/2004/data'.  
GET FILE='julydata.sav'.  
INSERT FILE='../commands/monthly_report.sps'.
```

## **Overview**

CD changes the working directory location, making it possible to use relative paths for subsequent file specifications in command syntax, including data files specified on commands such as GET and SAVE, command syntax files specified on commands such as INSERT and INCLUDE, and output files specified on commands such as OMS and WRITE.

## **Basic Specification**

The only specification is the command name followed by a quoted directory specification.

- The directory specification can contain a drive specification.
- The directory specification can be a previously defined file handle (see the FILE HANDLE command for more information).
- The directory specification can include paths defined in operating system environment variables.

## **Operations**

The change in the working directory remains in effect until some other condition occurs that changes the working directory during the session, such as explicitly changing the working directory on another CD command or an INSERT command with a CD keyword that specifies a different directory.

- If the directory path is a relative path, it is relative to the current working directory.
- If the directory specification contains a filename, the filename portion is ignored.

- If the last (most-nested) subdirectory in the directory specification does not exist, then it is assumed to be a filename and is ignored.
- If any directory specification prior to the last directory (or file) is invalid, the command will fail, and an error message is issued.

### **Limitations**

The CD command has no effect on the relative directory location for SET TLOOK file specifications. File specifications for the TLOOK subcommand of the SET command should include complete path information.

## **Examples**

### **Working with Absolute Paths**

```
CD '/sales/data/july.sav'.  
CD '/sales/data/july'.  
CD '/sales/data/july'.
```

If */sales/data* is a valid directory:

- The first CD command will ignore the filename *july.sav* and set the working directory to */sales/data*.
- If the subdirectory *july* exists, the second CD command will change the working directory to */sales/data/july*; otherwise, it will change the working directory to */sales/data*.
- The third CD command will fail if the *dqta* subdirectory doesn't exist.

### **Working with Relative Paths**

```
CD '/sales'.  
CD 'data'.  
CD 'july'.
```

If */sales* is a valid directory:

- The first CD command will change the working directory to */sales*.
- The relative path in the second CD command will change the working directory to */sales/data*.
- The relative path in the third CD command will change the working directory to */sales/data/july*.

## **Preserving and Restoring the Working Directory Setting**

The original working directory can be preserved with the PRESERVE command and later restored with the RESTORE command.

### **Example**

```
CD '/sales/data'.  
PRESERVE.  
CD '/commands/examples'.  
RESTORE.
```

- PRESERVE retains the working directory location set on the preceding CD command.
- The second CD command changes the working directory.
- RESTORE resets the working directory back to */sales/data*.

# ***CLEAR TIME PROGRAM***

```
CLEAR TIME PROGRAM.
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Overview***

CLEAR TIME PROGRAM deletes all time-dependent covariates created in the previous TIME PROGRAM command. It is primarily used in interactive mode to remove temporary variables associated with the time program so that you can redefine time-dependent covariates. It is not necessary to use this command if you have run a procedure that executes the TIME PROGRAM transformations, since all temporary variables created by TIME PROGRAM are automatically deleted.

### ***Basic Specification***

The only specification is the command itself. CLEAR TIME PROGRAM has no additional specifications.

## ***Example***

```
TIME PROGRAM.  
COMPUTE Z=AGE + T_.  
  
CLEAR TIME PROGRAM.  
  
TIME PROGRAM.  
COMPUTE Z=AGE + T_ - 18.  
  
COXREG SURVIVAL WITH Z  
  /STATUS SURVSTA EVENT(1).
```

- The first TIME PROGRAM command defines the time-dependent covariate *Z* as the current age.
- The CLEAR TIME PROGRAM command deletes the time-dependent covariate *Z*.
- The second TIME PROGRAM command redefines the time-dependent covariate *Z* as the number of years since turning 18.. *Z* is then specified as a covariate in COXREG.

# CLEAR TRANSFORMATIONS

CLEAR TRANSFORMATIONS

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Overview

CLEAR TRANSFORMATIONS discards previous data transformation commands.

### Basic Specification

The only specification is the command itself. CLEAR TRANSFORMATIONS has no additional specifications.

### Operations

- CLEAR TRANSFORMATIONS discards all data transformation commands that have accumulated since the last procedure.
- CLEAR TRANSFORMATIONS has no effect if a command file is submitted to your operating system for execution. It generates a warning when a command file is present.
- Be sure to delete CLEAR TRANSFORMATIONS and any unwanted transformation commands from the journal file if you plan to submit the file to the operating system for batch mode execution. Otherwise, the unwanted transformations will cause problems.

## Examples

```
GET FILE="/data/query.sav".
FREQUENCIES=ITEM1 ITEM2 ITEM3.
RECODE ITEM1, ITEM2, ITEM3 (0=1) (1=0) (2=-1).
COMPUTE INDEXQ=(ITEM1 + ITEM2 + ITEM3)/3.
VARIABLE LABELS INDEXQ 'SUMMARY INDEX OF QUESTIONS'.
CLEAR TRANSFORMATIONS.
DISPLAY DICTIONARY.
```

- The GET and FREQUENCIES commands are executed.
- The RECODE, COMPUTE, and VARIABLE LABELS commands are transformations. They do not affect the data until the next procedure is executed.
- The CLEAR TRANSFORMATIONS command discards the RECODE, COMPUTE, and VARIABLE LABELS commands.
- The DISPLAY command displays the working file dictionary. Data values and labels are exactly as they were when the FREQUENCIES command was executed. The variable *INDEXQ* does not exist because CLEAR TRANSFORMATIONS discarded the COMPUTE command.

# CLUSTER

CLUSTER is available in the Statistics Base option.

```
CLUSTER varlist [/MISSING={EXCLUDE**} [INCLUDE]]

[/MEASURE={ SEUCLID**          }
           {EUCLID              }
           {COSINE              }
           {CORRELATION         }
           {BLOCK               }
           {CHEBYCHEV          }
           {POWER(p,r)         }
           {MINKOWSKI(p)       }
           {CHISQ              }
           {PH2                }
           {RR[ (p[, np])]     }
           {SM[ (p[, np])]     }
           {JACCARD[ (p[, np])] }
           {DICE[ (p[, np])]   }
           {SS1[ (p[, np])]    }
           {RT[ (p[, np])]     }
           {SS2[ (p[, np])]    }
           {K1[ (p[, np])]     }
           {SS3[ (p[, np])]    }
           {K2[ (p[, np])]     }
           {SS4[ (p[, np])]    }
           {HAMANN[ (p[, np])] }
           {OCHIAI[ (p[, np])] }
           {SS5[ (p[, np])]    }
           {PHI[ (p[, np])]    }
           {LAMBDA[ (p[, np])] }
           {D[ (p[, np])]      }
           {Y[ (p[, np])]      }
           {Q[ (p[, np])]      }
           {BEUCLID[ (p[, np])] }
           {SIZE[ (p[, np])]   }
           {PATTERN[ (p[, np])] }
           {BSEUCLID[ (p[, np])] }
           {BSHAPE[ (p[, np])] }
           {DISPER[ (p[, np])] }
           {VARIANCE[ (p[, np])] }
           {BLWMN[ (p[, np])] }
           }

[/METHOD={ BAVERAGE** [(rootname)] [...]
          {WAVERAGE }
          {SINGLE }
          {COMPLETE }
          {CENTROID }
          {MEDIAN }
          {WARD }
          {DEFAULT** }

[/SAVE=CLUSTER({level })] [/ID=varname]
           {min,max}

[/PRINT=[CLUSTER({level })] [DISTANCE] [SCHEDULE**] [NONE]
           {min,max}

[/PLOT=[VICICLE** [(min[, max[, inc]])]] [DENDROGRAM] [NONE]
           {HICICLE[ (min[, max[, inc]])]}

[/MATRIX=[IN({'savfile'|'dataset'})] [OUT({'savfile'|'dataset'})]]
           {* } {* }
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Example**

```
CLUSTER V1 TO V4  
  /PLOT=DENDROGRAM  
  /PRINT=CLUSTER (2,4) .
```

## **Overview**

CLUSTER produces hierarchical clusters of items based on distance measures of dissimilarity or similarity. The items being clustered are usually cases from the active dataset, and the distance measures are computed from their values for one or more variables. You can also cluster variables if you read in a matrix measuring distances between variables. Cluster analysis is discussed in Anderberg (1973).

### **Options**

**Cluster Measures and Methods.** You can specify one of 37 similarity or distance measures on the MEASURE subcommand and any of the seven methods on the METHOD subcommand.

**New Variables.** You can save cluster membership for specified solutions as new variables in the active dataset using the SAVE subcommand.

**Display and Plots.** You can display cluster membership, the distance or similarity matrix used to cluster variables or cases, and the agglomeration schedule for the cluster solution with the PRINT subcommand. You can request either a horizontal or vertical icicle plot or a dendrogram of the cluster solution and control the cluster levels displayed in the icicle plot with the PLOT subcommand. You can also specify a variable to be used as a case identifier in the display on the ID subcommand.

**Matrix Input and Output.** You can write out the distance matrix and use it in subsequent CLUSTER, PROXIMITIES, or ALSICAL analyses or read in matrices produced by other CLUSTER or PROXIMITIES procedures using the MATRIX subcommand.

### **Basic Specification**

The basic specification is a variable list. CLUSTER assumes that the items being clustered are cases and uses the squared Euclidean distances between cases on the variables in the analysis as the measure of distance.

### **Subcommand Order**

- The variable list must be specified first.
- The remaining subcommands can be specified in any order.



**Syntax Rules**

- The variable list and subcommands can each be specified once.
- More than one clustering method can be specified on the `METHOD` subcommand.

**Operations**

The `CLUSTER` procedure involves four steps:

- First, `CLUSTER` obtains distance measures of similarities between or distances separating initial clusters (individual cases or individual variables if the input is a matrix measuring distances between variables).
- Second, it combines the two nearest clusters to form a new cluster.
- Third, it recomputes similarities or distances of existing clusters to the new cluster.
- It then returns to the second step until all items are combined in one cluster.

This process yields a hierarchy of cluster solutions, ranging from one overall cluster to as many clusters as there are items being clustered. Clusters at a higher level can contain several lower-level clusters. Within each level, the clusters are disjoint (each item belongs to only one cluster).

- `CLUSTER` identifies clusters in solutions by sequential integers (1, 2, 3, and so on).

**Limitations**

- `CLUSTER` stores cases and a lower-triangular matrix of proximities in memory. Storage requirements increase rapidly with the number of cases. You should be able to cluster 100 cases using a small number of variables in an 80K workspace.
- `CLUSTER` does not honor weights.

**Example**

```
CLUSTER V1 TO V4
  /PLOT=DENDROGRAM
  /PRINT=CLUSTER (2 4).
```

- This example clusters cases based on their values for all variables between and including *V1* and *V4* in the active dataset.
- The analysis uses the default measure of distance (squared Euclidean) and the default clustering method (average linkage between groups).
- `PLOT` requests a dendrogram.
- `PRINT` displays a table of the cluster membership of each case for the two-, three-, and four-cluster solutions.

**Used with the *PROXIMITIES* command to create distances**

```
PROXIMITIES price engine_s horsepow wheelbas width length
  curb_wgt fuel_cap mpg
  /MATRIX OUT ('C:/TEMP/spssclus.tmp')
  /VIEW= CASE
  /MEASURE= SEUCLID
  /PRINT NONE
```

*CLUSTER*

```

    /ID= model
    /STANDARDIZE= VARIABLE Z .
CLUSTER
  /MATRIX IN ('C:/TEMP/spssclus.tmp')
  /METHOD SINGLE
  /ID= model
  /PRINT SCHEDULE
  /PLOT DENDROGRAM.
ERASE FILE= 'C:/TEMP/spssclus.tmp'.

```

- The `PROXIMITIES` command creates distances suitable for cluster analysis using squared Euclidean distance as the measure and z score standardization of the variables, and saves them to a temporary data file using the `MATRIX` subcommand.
- The `CLUSTER` command reads the distances using the `MATRIX` subcommand.
- The `METHOD` subcommand specifies a single linkage analysis.
- The `ID` subcommand specifies that values of *model* should be used to identify cases in the output.
- The `PRINT` subcommand requests the agglomeration schedule.
- The `PLOT` subcommand requests a dendrogram.
- The `ERASE` command removes the temporary file.

***Variable List***

The variable list identifies the variables used to compute similarities or distances between cases.

- The variable list is required except when matrix input is used. It must be specified before the optional subcommands.
- If matrix input is used, the variable list can be omitted. The names for the items in the matrix are used to compute similarities or distances.
- You can specify a variable list to override the names for the items in the matrix. This allows you to read in a subset of cases for analysis. Specifying a variable that does not exist in the matrix results in an error.

***MEASURE Subcommand***

`MEASURE` specifies the distance or similarity measure used to cluster cases.

- If the `MEASURE` subcommand is omitted or included without specifications, squared Euclidean distances are used.
- Only one measure can be specified.

## Measures for Interval Data

For interval data, use any one of the following keywords on MEASURE:

<b>SEUCLID</b>	<p><i>Squared Euclidean distance.</i> The distance between two items, <math>x</math> and <math>y</math>, is the sum of the squared differences between the values for the items. SEUCLID is the measure commonly used with centroid, median, and Ward's methods of clustering. SEUCLID is the default and can also be requested with keyword DEFAULT.</p> $SEUCLID(x, y) = \sum_i (x_i - y_i)^2$
<b>EUCLID</b>	<p><i>Euclidean distance.</i> This is the default specification for MEASURE. The distance between two items, <math>x</math> and <math>y</math>, is the square root of the sum of the squared differences between the values for the items.</p> $EUCLID(x, y) = \sqrt{\sum_i (x_i - y_i)^2}$
<b>CORRELATION</b>	<p><i>Correlation between vectors of values.</i> This is a pattern similarity measure.</p> $CORRELATION(x, y) = \frac{\sum_i (Z_{xi} Z_{yi})}{N-1}$ <p>where <math>Z_{xi}</math> is the <math>z</math> score (standardized) value of <math>x</math> for the <math>i</math>th case or variable, and <math>N</math> is the number of cases or variables.</p>
<b>COSINE</b>	<p><i>Cosine of vectors of values.</i> This is a pattern similarity measure.</p> $COSINE(x, y) = \frac{\sum_i (x_i y_i)}{\sqrt{(\sum_i x_i^2)(\sum_i y_i^2)}}$
<b>CHEBYCHEV</b>	<p><i>Chebychev distance metric.</i> The distance between two items is the maximum absolute difference between the values for the items.</p> $CHEBYCHEV(x, y) = \max_i  x_i - y_i $
<b>BLOCK</b>	<p><i>City-block or Manhattan distance.</i> The distance between two items is the sum of the absolute differences between the values for the items.</p> $BLOCK(x, y) = \sum_i  x_i - y_i $
<b>MINKOWSKI(p)</b>	<p><i>Distance in an absolute Minkowski power metric.</i> The distance between two items is the <math>p</math>th root of the sum of the absolute differences to the <math>p</math>th power between the values for the items. Appropriate selection of the integer parameter <math>p</math> yields Euclidean and many other distance metrics.</p> $MINKOWSKI(x, y) = (\sum_i  x_i - y_i ^p)^{1/p}$
<b>POWER(p,r)</b>	<p><i>Distance in an absolute power metric.</i> The distance between two items is the <math>r</math>th root of the sum of the absolute differences to the <math>p</math>th power between the values for the items. Appropriate selection of the integer parameters <math>p</math> and <math>r</math> yields Euclidean, squared Euclidean, Minkowski, city-block, and many other distance metrics.</p> $POWER(x, y) = (\sum_i  x_i - y_i ^p)^{1/r}$

## Measures for Frequency Count Data

For frequency count data, use any one of the following keywords on MEASURE:

**CHISQ** *Based on the chi-square test of equality for two sets of frequencies.* The magnitude of this dissimilarity measure depends on the total frequencies of the two cases or variables whose dissimilarity is computed. Expected values are from the model of independence of cases or variables  $x$  and  $y$ .

$$CHISQ(x, y) = \sqrt{\frac{\sum_i (x_i - E(x_i))^2}{E(x_i)} + \frac{\sum_i (y_i - E(y_i))^2}{E(y_i)}}$$

**PH2** *Phi-square between sets of frequencies.* This is the CHISQ measure normalized by the square root of the combined frequency. Therefore, its value does not depend on the total frequencies of the two cases or variables whose dissimilarity is computed.

$$PH2(x, y) = \sqrt{\frac{\frac{\sum_i (x_i - E(x_i))^2}{E(x_i)} + \frac{\sum_i (y_i - E(y_i))^2}{E(y_i)}}{N}}$$

## Measures for Binary Data

Different binary measures emphasize different aspects of the relationship between sets of binary values. However, all the measures are specified in the same way. Each measure has two optional integer-valued parameters,  $p$  (present) and  $np$  (not present).

- If both parameters are specified, CLUSTER uses the value of the first as an indicator that a characteristic is present and the value of the second as an indicator that a characteristic is absent. CLUSTER skips all other values.
- If only the first parameter is specified, CLUSTER uses that value to indicate presence and all other values to indicate absence.
- If no parameters are specified, CLUSTER assumes that 1 indicates presence and 0 indicates absence.

Using the indicators for presence and absence within each item (case or variable), CLUSTER constructs a  $2 \times 2$  contingency table for each pair of items in turn. It uses this table to compute a proximity measure for the pair.

		Item 2 characteristics	
		Present	Absent
Item 1 characteristics	Present	$a$	$b$
	Absent	$c$	$d$

CLUSTER computes all binary measures from the values of  $a$ ,  $b$ ,  $c$ , and  $d$ . These values are tallied across variables (when the items are cases) or across cases (when the items are variables). For example, if the variables  $V, W, X, Y, Z$  have values 0, 1, 1, 0, 1 for case 1 and values 0, 1, 1, 0, 0 for case 2 (where 1 indicates presence and 0 indicates absence), the contingency table is as follows:

		Case 2 characteristics	
		Present	Absent
Case 1 characteristics			

<b>Present</b>	2	1
<b>Absent</b>	0	2

The contingency table indicates that both cases are present for two variables ( $W$  and  $X$ ), both cases are absent for two variables ( $V$  and  $Y$ ), and case 1 is present and case 2 is absent for one variable ( $Z$ ). There are no variables for which case 1 is absent and case 2 is present.

The available binary measures include matching coefficients, conditional probabilities, predictability measures, and others.

**Matching Coefficients.** The table below shows a classification scheme for matching coefficients. In this scheme, *matches* are joint presences (value  $a$  in the contingency table) or joint absences (value  $d$ ). *Nonmatches* are equal in number to value  $b$  plus value  $c$ . Matches and nonmatches may or may not be weighted equally. The three coefficients JACCARD, DICE, and SS2 are related monotonically, as are SM, SS1, and RT. All coefficients in the table are similarity measures, and all except two (K1 and SS3) range from 0 to 1. K1 and SS3 have a minimum value of 0 and no upper limit.

Table 31-1  
Binary matching coefficients in CLUSTER

	Joint absences excluded from numerator	Joint absences included in numerator
<b>All matches included in denominator</b>		
Equal weight for matches and nonmatches	RR	SM
Double weight for matches		SS1
Double weight for nonmatches		RT
<b>Joint absences excluded from denominator</b>		
Equal weight for matches and nonmatches	JACCARD	
Double weight for matches	DICE	
Double weight for nonmatches	SS2	
<b>All matches excluded from denominator</b>		
Equal weight for matches and nonmatches	K1	SS3

**RR[(p,np)]**

*Russell and Rao similarity measure.* This is the binary dot product.

$$RR(x, y) = \frac{a}{a+b+c+d}$$

**SM[(p,np)]**

*Simple matching similarity measure.* This is the ratio of the number of matches to the total number of characteristics.

$$SM(x, y) = \frac{a+d}{a+b+c+d}$$

**JACCARD**[(p[,np])] *Jaccard similarity measure.* This is also known as the *similarity ratio*.

$$JACCARD(x, y) = \frac{a}{a+b+c}$$

**DICE**[(p[,np])] *Dice (or Czekanowski or Sorenson) similarity measure.*

$$DICE(x, y) = \frac{2a}{2a+b+c}$$

**SS1**[(p[,np])] *Sokal and Sneath similarity measure 1.*

$$SS1(x, y) = \frac{2(a+d)}{2(a+d)+b+c+d}$$

**RT**[(p[,np])] *Rogers and Tanimoto similarity measure.*

$$RT(x, y) = \frac{a+d}{a+d+2(b+c)}$$

**SS2**[(p[,np])] *Sokal and Sneath similarity measure 2.*

$$SS2(x, y) = \frac{a}{a+2(b+c)}$$

**K1**[(p[,np])] *Kulczynski similarity measure 1.* This measure has a minimum value of 0 and no upper limit. It is undefined when there are no nonmatches ( $b=0$  and  $c=0$ ).

$$K1(x, y) = \frac{a}{b+c}$$

**SS3**[(p[,np])] *Sokal and Sneath similarity measure 3.* This measure has a minimum value of 0 and no upper limit. It is undefined when there are no nonmatches ( $b=0$  and  $c=0$ ).

$$SS3(x, y) = \frac{a+d}{b+c}$$

**Conditional Probabilities.** The following binary measures yield values that can be interpreted in terms of conditional probability. All three are similarity measures.

**K2**[(p[,np])] *Kulczynski similarity measure 2.* This yields the average conditional probability that a characteristic is present in one item given that the characteristic is present in the other item. The measure is an average over both items acting as predictors. It has a range of 0 to 1.

$$K2(x, y) = \frac{a/(a+b)+a/(a+c)}{2}$$

**SS4**[(p[,np])] *Sokal and Sneath similarity measure 4.* This yields the conditional probability that a characteristic of one item is in the same state (presence or absence) as the characteristic of the other item. The measure is an average over both items acting as predictors. It has a range of 0 to 1.

$$SS4(x, y) = \frac{a/(a+b)+a/(a+c)+d/(b+d)+d/(c+d)}{4}$$

**HAMANN**[(p[,np])] *Hamann similarity measure.* This measure gives the probability that a characteristic has the same state in both items (present in both or absent from both) minus the probability that a characteristic has different states in the two items (present in one and absent from the other). **HAMANN** has a range of  $-1$  to  $+1$  and is monotonically related to **SM**, **SS1**, and **RT**.

$$HAMANN(x, y) = \frac{(a+d)-(b+c)}{a+b+c+d}$$

**Predictability Measures.** The following four binary measures assess the association between items as the predictability of one given the other. All four measures yield similarities.

**LAMBDA[(p[,np])]**

*Goodman and Kruskal's lambda (similarity).* This coefficient assesses the predictability of the state of a characteristic on one item (present or absent) given the state on the other item. Specifically, LAMBDA measures the proportional reduction in error using one item to predict the other when the directions of prediction are of equal importance. LAMBDA has a range of 0 to 1.

$$LAMBDA(x, y) = \frac{t_1 - t_2}{2(a+b+c+d) - t_2}$$

where

$$t_1 = \max(a, b) + \max(c, d) + \max(a, c) + \max(b, d)$$

$$t_2 = \max(a + c, b + d) + \max(a + d, c + d).$$

**D[(p[,np])]**

*Anderberg's D (similarity).* This coefficient assesses the predictability of the state of a characteristic on one item (present or absent) given the state on the other. D measures the actual reduction in the error probability when one item is used to predict the other. The range of D is 0 to 1.

$$D(x, y) = \frac{t_1 - t_2}{2(a+b+c+d)}$$

where

$$t_1 = \max(a, b) + \max(c, d) + \max(a, c) + \max(b, d)$$

$$t_2 = \max(a + c, b + d) + \max(a + d, c + d).$$

**Y[(p[,np])]**

*Yule's Y coefficient of colligation (similarity).* This is a function of the cross-ratio for a  $2 \times 2$  table. It has a range of  $-1$  to  $+1$ .

$$Y(x, y) = \frac{\sqrt{ad} - \sqrt{bc}}{\sqrt{ad} + \sqrt{bc}}$$

**Q[(p[,np])]**

*Yule's Q (similarity).* This is the  $2 \times 2$  version of Goodman and Kruskal's ordinal measure *gamma*. Like Yule's Y, Q is a function of the cross-ratio for a  $2 \times 2$  table and has a range of  $-1$  to  $+1$ .

$$Q(x, y) = \frac{ad - bc}{ad + bc}$$

**Other Binary Measures.** The remaining binary measures available in CLUSTER are either binary equivalents of association measures for continuous variables or measures of special properties of the relationship between items.

**OCHIAI[(p[,np])]**

*Ochiai similarity measure.* This is the binary form of the cosine. It has a range of 0 to 1.

$$OCHIAI(x, y) = \sqrt{\frac{a}{a+b} \frac{a}{a+c}}$$

**SS5[(p[,np])]**

*Sokal and Sneath similarity measure 5.* The range is 0 to 1.

$$SS5(x, y) = \frac{ad}{\sqrt{(a+b)(a+c)(b+d)(c+d)}}$$

<b>PHI</b> [(p[,np])]	<i>Fourfold point correlation (similarity)</i> . This is the binary form of the Pearson product-moment correlation coefficient.
	$PHI(x, y) = \frac{ad-bc}{\sqrt{(a+b)(a+c)(b+d)(c+d)}}$
<b>BEUCLID</b> [(p[,np])]	<i>Binary Euclidean distance</i> . This is a distance measure. Its minimum value is 0, and it has no upper limit.
	$BEUCLID(x, y) = \sqrt{b+c}$
<b>BSEUCLID</b> [(p[,np])]	<i>Binary squared Euclidean distance</i> . This is a distance measure. Its minimum value is 0, and it has no upper limit.
	$BSEUCLID(x, y) = b+c$
<b>SIZE</b> [(p[,np])]	<i>Size difference</i> . This is a dissimilarity measure with a minimum value of 0 and no upper limit.
	$SIZE(x, y) = \frac{(b-c)^2}{(a+b+c+d)^2}$
<b>PATTERN</b> [(p[,np])]	<i>Pattern difference</i> . This is a dissimilarity measure. The range is 0 to 1.
	$PATTERN(x, y) = \frac{bc}{(a+b+c+d)^2}$
<b>BSHAPE</b> [(p[,np])]	<i>Binary shape difference</i> . This dissimilarity measure has no upper or lower limit.
	$BSHAPE(x, y) = \frac{(a+b+c+d)(b+c)-(b-c)^2}{(a+b+c+d)^2}$
<b>DISPER</b> [(p[,np])]	<i>Dispersion similarity measure</i> . The range is -1 to +1.
	$DISPER(x, y) = \frac{ad-bc}{(a+b+c+d)^2}$
<b>VARIANCE</b> [(p[,np])]	<i>Variance dissimilarity measure</i> . This measure has a minimum value of 0 and no upper limit.
	$VARIANCE(x, y) = \frac{b+c}{4(a+b+c+d)}$
<b>BLWMN</b> [(p[,np])]	<i>Binary Lance-and-Williams nonmetric dissimilarity measure</i> . This measure is also known as the Bray-Curtis nonmetric coefficient. The range is 0 to 1.
	$BLWMN(x, y) = \frac{b+c}{2a+b+c}$

## **METHOD Subcommand**

METHOD specifies one or more clustering methods.

- If the METHOD subcommand is omitted or included without specifications, the method of average linkage between groups is used.
- Only one METHOD subcommand can be used, but more than one method can be specified on it.



- When the number of items is large, **CENTROID** and **MEDIAN** require significantly more CPU time than other methods.

<b>BAVERAGE</b>	<i>Average linkage between groups (UPGMA).</i> <b>BAVERAGE</b> is the default and can also be requested with keyword <b>DEFAULT</b> .
<b>WAVERAGE</b>	<i>Average linkage within groups.</i>
<b>SINGLE</b>	<i>Single linkage or nearest neighbor.</i>
<b>COMPLETE</b>	<i>Complete linkage or furthest neighbor.</i>
<b>CENTROID</b>	<i>Centroid clustering (UPGMC).</i> Squared Euclidean distances are commonly used with this method.
<b>MEDIAN</b>	<i>Median clustering (WPGMC).</i> Squared Euclidean distances are commonly used with this method.
<b>WARD</b>	<i>Ward's method.</i> Squared Euclidean distances are commonly used with this method.

### Example

```
CLUSTER V1 V2 V3
/METHOD=SINGLE COMPLETE WARDS.
```

- This example clusters cases based on their values for the variables *V1*, *V2*, and *V3* and uses three clustering methods: single linkage, complete linkage, and Ward's method.

## SAVE Subcommand

**SAVE** allows you to save cluster membership at specified solution levels as new variables in the active dataset.

- The specification on **SAVE** is the **CLUSTER** keyword, followed by either a single number indicating the level (number of clusters) of the cluster solution or a range separated by a comma indicating the minimum and maximum numbers of clusters when membership of more than one solution is to be saved. The number or range must be enclosed in parentheses and applies to all methods specified on **METHOD**.
- You can specify a rootname in parentheses after each method specification on the **METHOD** subcommand. **CLUSTER** forms new variable names by appending the number of the cluster solution to the rootname.
- If no rootname is specified, **CLUSTER** forms variable names using the formula  $CLU_n_m$ , where  $m$  increments to create a unique rootname for the set of variables saved for one method and  $n$  is the number of the cluster solution.
- The names and descriptive labels of the new variables are displayed in the procedure information notes.
- You cannot use the **SAVE** subcommand if you are replacing the active dataset with matrix materials ([For more information, see the topic Matrix Output on p. 319.](#))

### Example

```
CLUSTER A B C
/METHOD=BAVERAGE SINGLE (SINMEM) WARD
/SAVE=CLUSTERS(3,5).
```

- This command creates nine new variables: *CLU5\_1*, *CLU4\_1*, and *CLU3\_1* for *BAVERAGE*, *SINMEM5*, *SINMEM4*, and *SINMEM3* for *SINGLE*, and *CLU5\_2*, *CLU4\_2*, and *CLU3\_2* for *WARD*. The variables contain the cluster membership for each case at the five-, four-, and three-cluster solutions using the three clustering methods. Ward's method is the third specification on *METHOD* but uses the second set of default names, since it is the second method specified without a rootname.
- The order of the new variables in the active dataset is the same as listed above, since the solutions are obtained in the order from 5 to 3.
- New variables are listed in the procedure information notes.

## ***ID Subcommand***

*ID* names a string variable to be used as the case identifier in cluster membership tables, icicle plots, and dendrograms. If the *ID* subcommand is omitted, cases are identified by case numbers alone.

- When used with the *MATRIX IN* subcommand, the variable specified on the *ID* subcommand identifies the labeling variable in the matrix file.

## ***PRINT Subcommand***

*PRINT* controls the display of cluster output (except plots, which are controlled by the *PLOT* subcommand).

- If the *PRINT* subcommand is omitted or included without specifications, an agglomeration schedule is displayed. If any keywords are specified on *PRINT*, the agglomeration schedule is displayed only if explicitly requested.
- *CLUSTER* automatically displays summary information (the method and measure used, the number of cases) for each method named on the *METHOD* subcommand. This summary is displayed regardless of specifications on *PRINT*.

You can specify any or all of the following on the *PRINT* subcommand:

### **SCHEDULE**

*Agglomeration schedule.* The agglomeration schedule shows the order and distances at which items and clusters combine to form new clusters. It also shows the cluster level at which an item joins a cluster. *SCHEDULE* is the default and can also be requested with the keyword *DEFAULT*.

### **CLUSTER(min,max)**

*Cluster membership.* For each item, the display includes the value of the case identifier (or the variable name if matrix input is used), the case sequence number, and a value (1, 2, 3, and so on) identifying the cluster to which that case belongs in a given cluster solution. Specify either a single integer value in parentheses indicating the level of a single solution or a minimum value and a maximum value indicating a range of solutions for which display is desired. If the number of clusters specified exceeds the number produced, the largest number of clusters is used (the number of items minus 1). If *CLUSTER* is specified more than once, the last specification is used.

<b>DISTANCE</b>	<i>Proximities matrix.</i> The proximities matrix table displays the distances or similarities between items computed by CLUSTER or obtained from an input matrix. DISTANCE produces a large volume of output and uses significant CPU time when the number of cases is large.
<b>NONE</b>	<i>None of the above.</i> NONE overrides any other keywords specified on PRINT.

**Example**

```
CLUSTER V1 V2 V3 /PRINT=CLUSTER(3,5).
```

- This example displays cluster membership for each case for the three-, four-, and five-cluster solutions.

**PLOT Subcommand**

PLOT controls the plots produced for each method specified on the METHOD subcommand. For icicle plots, PLOT allows you to control the cluster solution at which the plot begins and ends and the increment for displaying intermediate cluster solutions.

- If the PLOT subcommand is omitted or included without specifications, a vertical icicle plot is produced.
- If any keywords are specified on PLOT, only those plots requested are produced.
- The icicle plots are generated as pivot tables and the dendrogram is generated as text output.
- If there is not enough memory for a dendrogram or an icicle plot, the plot is skipped and a warning is issued.
- The size of an icicle plot can be controlled by specifying range values or an increment for VICICLE or HICICLE. Smaller plots require significantly less workspace and time.

<b>VICICLE(min,max,inc)</b>	<i>Vertical icicle plot.</i> This is the default. The range specifications are optional. If used, they must be integer and must be enclosed in parentheses. The specification <i>min</i> is the cluster solution at which to start the display (the default is 1), and the specification <i>max</i> is the cluster solution at which to end the display (the default is the number of cases minus 1). If <i>max</i> is greater than the number of cases minus 1, the default is used. The increment to use between cluster solutions is <i>inc</i> (the default is 1). If <i>max</i> is specified, <i>min</i> must be specified, and if <i>inc</i> is specified, both <i>min</i> and <i>max</i> must be specified. If VICICLE is specified more than once, only the last range specification is used.
<b>HICICLE(min,max,inc)</b>	<i>Horizontal icicle plot.</i> The range specifications are the same as for VICICLE. If both VICICLE and HICICLE are specified, the last range specified is used for both. If a range is not specified on the last instance of VICICLE or HICICLE, the defaults are used even if a range is specified earlier.
<b>DENDROGRAM</b>	<i>Tree diagram.</i> The dendrogram is scaled by the joining distances of the clusters.
<b>NONE</b>	<i>No plots.</i>

**Example**

```
CLUSTER V1 V2 V3 /PLOT=VICICLE(1,20).
```

- This example produces a vertical icicle plot for the 1-cluster through the 20-cluster solution.

### **Example**

```
CLUSTER V1 V2 V3 /PLOT=VICICLE(1,151,5).
```

- This example produces a vertical icicle plot for every fifth cluster solution starting with 1 and ending with 151 (1 cluster, 6 clusters, 11 clusters, and so on).

## **MISSING Subcommand**

MISSING controls the treatment of cases with missing values. A case that has a missing value for any variable on the variable list is omitted from the analysis. By default, user-missing values are excluded from the analysis.

<b>EXCLUDE</b>	<i>Exclude cases with user-missing values. This is the default.</i>
<b>INCLUDE</b>	<i>Include cases with user-missing values. Only cases with system-missing values are excluded.</i>

## **MATRIX Subcommand**

MATRIX reads and writes IBM® SPSS® Statistics data files.

- Either IN or OUT and a matrix file in parentheses are required. When both IN and OUT are used on the same CLUSTER procedure, they can be specified on separate MATRIX subcommands or on the same subcommand.
- The input or output matrix information is displayed in the procedure information notes.

**OUT ('savfile' | 'dataset')** *Write a matrix data file.* Specify either a quoted file specification, a previously declared dataset (DATASET DECLARE), or an asterisk in parentheses (\*). If you specify an asterisk (\*), the matrix data file replaces the active dataset.

**IN ('savfile' | 'dataset')** *Read a matrix data file.* Specify either a quoted file specification, a previously declared dataset (DATASET DECLARE), or an asterisk in parentheses (\*). The asterisk specifies the active dataset. A matrix file read from an external file does not replace the active dataset.

When a matrix is produced using the MATRIX OUT subcommand, it corresponds to a unique dataset. All subsequent analyses performed on this matrix would match the corresponding analysis on the original data. However, if the data file is altered in any way, this would no longer be true.

For example, if the original file is edited or rearranged, it would in general no longer correspond to the initially produced matrix. You need to make sure that the data match the matrix whenever inferring the results from the matrix analysis. Specifically, when saving the cluster membership into an active dataset in the CLUSTER procedure, the proximity matrix in the MATRIX IN statement must match the current active dataset.

## Matrix Output

- CLUSTER writes proximity-type matrices with *ROWTYPE\_* values of PROX. CLUSTER neither reads nor writes additional statistics with its matrix materials. For more information, see the topic [Format of the Matrix Data File on p. 319](#).
- The matrices produced by CLUSTER can be used by subsequent CLUSTER procedures or by the PROXIMITIES and ALSICAL procedures.
- Any documents contained in the active dataset are not transferred to the matrix file.

## Matrix Input

- CLUSTER can read matrices written by a previous CLUSTER command or by PROXIMITIES, or created by MATRIX DATA. When the input matrix contains distances between variables, CLUSTER clusters all or a subset of the variables.
- Values for split-file variables should precede values for *ROWTYPE\_*, *CASENO\_* and the labeling variable (if present) should come after *ROWTYPE\_* and before *VARNAME\_*.
- If *CASENO\_* is of type string rather than numeric, it will be considered unavailable and a warning is issued.
- If *CASENO\_* appears on a variable list, a syntax error results.
- CLUSTER ignores unrecognized *ROWTYPE\_* values.
- When you are reading a matrix created with MATRIX DATA, you should supply a value label for PROX of either *SIMILARITY* or *DISSIMILARITY* so that the matrix is correctly identified. If you do not supply a label, CLUSTER assumes *DISSIMILARITY*. (See “Format of the Matrix Data File” below.)
- The program reads variable names, variable and value labels, and print and write formats from the dictionary of the matrix data file.
- MATRIX=IN cannot be specified unless an active dataset has already been defined. To read an existing matrix data file at the beginning of a session, use GET to retrieve the matrix file and then specify IN(\*) on MATRIX.
- The variable list on CLUSTER can be omitted when a matrix data file is used as input. By default, all cases or variables in the matrix data file are used in the analysis. Specify a variable list when you want to read in a subset of items for analysis.

## Format of the Matrix Data File

- The matrix data file can include three special variables created by the program: *ROWTYPE\_*, *ID*, and *VARNAME\_*.
- The variable *ROWTYPE\_* is a string variable with the value PROX (for proximity measure). PROX is assigned value labels containing the distance measure used to create the matrix and either *SIMILARITY* or *DISSIMILARITY* as an identifier. The variable *VARNAME\_* is a short string variable whose values are the names of the new variables. The variable *CASENO\_* is a numeric variable with values equal to the original case numbers.
- *ID* is included only when an identifying variable is not specified on the ID subcommand. *ID* is a short string and takes the value CASE *m*, where *m* is the actual number of each case. Note that *m* may not be consecutive if cases have been selected.

- If an identifying variable is specified on the `ID` subcommand, it takes the place of `ID` between `ROWTYPE_` and `VARNAME_`. Up to 20 characters can be displayed for the identifying variable.
- `VARNAME_` is a string variable that takes the values `VAR1`, `VAR2`, ..., `VARn` to correspond to the names of the distance variables in the matrix (`VAR1`, `VAR2`, ..., `VARn`, where `n` is the number of cases in the largest split file). The numeric suffix for the variable names is consecutive and may not be the same as the actual case number.
- The remaining variables in the matrix file are the distance variables used to form the matrix. The distance variables are assigned variable labels in the form of `CASE m` to identify the actual number of each case.

### Split Files

- When split-file processing is in effect, the first variables in the matrix data file are the split variables, followed by `ROWTYPE_`, the case-identifier variable or `ID`, `VARNAME_`, and the distance variables.
- A full set of matrix materials is written for each split-file group defined by the split variables.
- A split variable cannot have the same name as any other variable written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split file must be in effect when that matrix is read by any procedure.

### Missing Values

Missing-value treatment affects the values written to a matrix data file. When reading a matrix data file, be sure to specify a missing-value treatment on `CLUSTER` that is compatible with the treatment that was in effect when the matrix materials were generated.

### Example: Output to External File

```
DATA LIST FILE=ALMANAC1 RECORDS=3
  /1 CITY 6-18(A) POP80 53-60
  /2 CHURCHES 10-13 PARKS 14-17 PHONES 18-25 TVS 26-32
    RADIOST 33-35 TVST 36-38 TAXRATE 52-57(2) .
N OF CASES 8.

CLUSTER CHURCHES TO TAXRATE
  /ID=CITY
  /MEASURE=EUCLID
  /MATRIX=OUT (CLUSMTX) .
```

- `CLUSTER` reads raw data from file `ALMANAC1` and writes one set of matrix materials to file `CLUSMTX`.
- The active dataset is still the `ALMANAC1` file defined on `DATA LIST`. Subsequent commands are executed on `ALMANAC1`.

### Example: Output Replacing Active Dataset

```
DATA LIST FILE=ALMANAC1 RECORDS=3
```

```

/1 CITY 6-18(A) POP80 53-60
/2 CHURCHES 10-13 PARKS 14-17 PHONES 18-25 TVS 26-32
   RADIOST 33-35 TVST 36-38 TAXRATE 52-57(2) .
N OF CASES 8 .

CLUSTER CHURCHES TO TAXRATE
/ID=CITY
/MEASURE=EUCLID
/MATRIX=OUT(*) .
LIST .

```

- CLUSTER writes the same matrix as in the previous example. However, the matrix data file replaces the active dataset. The LIST command is executed on the matrix file, not on *ALMANACI*.

### Example: Input from Active Dataset

```

GET FILE=CLUSMTX .
CLUSTER
/ID=CITY
/MATRIX=IN(*) .

```

- This example starts a new session and reads an existing matrix data file. GET retrieves the matrix data file *CLUSMTX*.
- MATRIX=IN specifies an asterisk because the matrix data file is the active dataset. If MATRIX=IN(CLUSMTX) is specified, the program issues an error message.
- If the GET command is omitted, the program issues an error message.

### Example: Input from External File

```

GET FILE=PRSNL .
FREQUENCIES VARIABLE=AGE .

CLUSTER
/ID=CITY
/MATRIX=IN(CLUSMTX) .

```

- This example performs a frequencies analysis on the file *PRSNL* and then uses a different file for CLUSTER. The file is an existing matrix data file.
- The variable list is omitted on the CLUSTER command. By default, all cases in the matrix file are used in the analysis.
- MATRIX=IN specifies the matrix data file *CLUSMTX*.
- *CLUSMTX* does not replace *PRSNL* as the active dataset.

### Example: Input from Active Dataset

```

GET FILE='data/crime.sav' .
PROXIMITIES MURDER TO MOTOR
/VIEW=VARIABLE
/MEASURE=PH2
/MATRIX=OUT(*) .
CLUSTER
/MATRIX=IN(*) .

```

- PROXIMITIES uses the data from *crime.sav*, which is now the active dataset. The VIEW subcommand specifies computation of proximity values between variables. The MATRIX subcommand writes the matrix to the active dataset.
- MATRIX=IN(\*) on the CLUSTER command reads the matrix materials from the active dataset. Since the matrix contains distances between variables, CLUSTER clusters variables based on distance measures in the input. The variable list is omitted on the CLUSTER command, so all variables are used in the analysis. The slash preceding the MATRIX subcommand is required because there is an implied variable list. Without the slash, CLUSTER would attempt to interpret MATRIX as a variable name rather than a subcommand name.



# CODEBOOK

CODEBOOK is available in the Statistics Base option.

*Note:* Square brackets used in the CODEBOOK syntax chart are required parts of the syntax and are not used to indicate optional elements. Equals signs (=) used in the syntax chart are required elements. All subcommands are optional.

```
CODEBOOK variable [level] variable [level] variable [level]...

/VARINFO
  POSITION LABEL TYPE FORMAT MEASURE ROLE ATTRIBUTES VALUELABELS MISSING
  RESERVEDATTRIBUTES

/FILEINFO NAME LOCATION CASECOUNT LABEL ATTRIBUTES DOCUMENTS WEIGHT
  RESERVEDATTRIBUTES

/STATISTICS NONE COUNT PERCENT MEAN STDDEV QUANTILES

/OPTIONS MAXCATS=200*
          VARORDER={FILE**      }
                {ALPHA         }
                {VARLIST†      }
                {MEASURE       }
                {ATTRIBUTE(name)}
          SORT={ASCENDING** }
                {DESCENDING  }
```

\*Default if subcommand or keyword omitted.

\*\*Default if subcommand omitted and there is no variable list.

†Default if subcommand omitted and the command includes a variable list.

- If the VARINFO subcommand is omitted, all variable information except RESERVEDATTRIBUTES is included.
- If the STATISTICS subcommand is omitted, all statistics are included.

## **Release History**

Release 17.0

- Command introduced.

Release 18

- ROLE keyword added to VARINFO subcommand.

## **Example**

```
CODEBOOK Age Income $MultCars.
```

## Overview

CODEBOOK reports the dictionary information — such as variable names, variable labels, value labels, missing values — and summary statistics for all or specified variables and multiple response sets in the active dataset. For nominal and ordinal variables and multiple response sets, summary statistics include counts and percents. For scale variables, summary statistics include mean, standard deviation, and quartiles.

### Options

Optionally, you can:

- Specify the variables and/or multiple response sets to include in the report.
- Choose the types of dictionary information to display.
- Suppress the display of summary statistics for any nominal and ordinal variables and multiple response set with more than a specified number of unique values.
- Override the defined measurement level for a variable, thereby changing the available summary statistics for the variable.
- Include a table of file information, such as file name and location and number of cases.
- Sort the variables in the report by variable name or label or other dictionary attributes such as type, format, or measurement level.

### Basic Specification

The basic specification is the command name `CODEBOOK` with no additional specifications.

### Subcommand Order

The command name is followed by the optional variable list, followed by the optional subcommands in any order.

### Syntax Rules

- Variables and multiple response sets listed on the optional variable list must already exist in the active dataset.
- Each variable can only be specified or implied once.
- Multiple response set names must include the leading dollar sign (\$).
- The keyword `TO` can be used to specify consecutive variables in file order. It cannot be used to specify a list of multiple response sets.
- Each subcommand can only be specified once.
- Equals signs and square brackets shown in examples are required syntax elements.

### Operations

- By default, `CODEBOOK` reads the active dataset and causes the execution of any pending transformations.

- If `/STATISTICS NONE` is specified, CODEBOOK does not read the active dataset or execute pending transformations.
- `SPLIT FILE` status is ignored. This includes split-file groups created by the `MULTIPLE IMPUTATION` command (available in the Missing Values add-on option).
- `FILTER` status is honored for computing summary statistics.

## Examples

### **CODEBOOK with No Additional Specifications**

```
CODEBOOK.
```

The default output includes:

- Variable information for all variables in the dataset, except for reserved system attributes.
- Counts and percents for all categories of nominal and ordinal variables, labeled categories of scale variables, and multiple response sets.
- Mean, standard deviation, and quartiles for scale variables.

### **Specifying Specific Variables, Variable Information, and Statistics**

```
CODEBOOK Var1 Var3 [N] $Multvar
  /VARINFO LABEL MEASURE VALUELABELS MISSING
  /STATISTICS COUNT MEAN
  /OPTIONS MAXCATS=10.
```

- The results will only include information for the two specified variables and the one multiple response set.
- `Var3 [N]` indicates that *Var3* should be treated as nominal for summary statistics. This has no effect on the defined measurement level for the variable or the measurement level displayed in the results.
- Dictionary information will be limited to variable label, measurement level, value labels, and missing values.
- Only counts will be included for nominal/ordinal variables, multiple response sets and labeled categories of scale variables.
- Only the mean will be included for scale variables.
- For nominal/ordinal variables, multiple response sets, and labeled values of scale variables, `MAXCATS=10` will suppress the display of value labels and counts if there are more than 10 unique, valid values.

## Variable List

The optional variable list specification allows you to limit the results to specified variables and/or multiple response sets and override the defined measurement level for specified numeric variables.

- Each variable or multiple response set can be specified or implied only once.
- Multiple response set names must include the leading dollar sign (\$).

- Keyword `TO` can be used to specify consecutive variables in file order. It cannot be used to specify a list of multiple response sets.
- Keyword `ALL` can be used to specify all variables (does not include multiple response sets). Note: `ALL` cannot be used in combination with a list of one or more specific variable names since a variable can be specified only once.
- Keyword `$ALL` can be used to specify all multiple response sets (does not include variables). Note: `$ALL` cannot be used in combination with a list of one or more specific multiple response sets, since a set can be specified only once.

### ***Overriding Defined Measurement Level***

Available summary statistics are determined by the measurement level of the variable. You can override the defined measurement level by including a measurement level specification in square brackets after the variable name: `[N]` for nominal, `[O]` for ordinal, and `[S]` for scale. This does not change the defined measurement level for the variable, and if results include measurement level, the defined measurement level is displayed in the results.

- For string variables and multiple response sets, measurement level can only be nominal or ordinal.
- Measurement level specification cannot be used with keywords `ALL` or `$ALL`.

### ***Example***

```
CODEBOOK Var1 Var3 [N] Var5 TO Var8 [N] $Multvar.
```

- *Var3* and *Var8* will be treated as nominal for summary statistics, so the available summary statistics for those two variables are counts and percents.
- The defined measurement level will be used to determine available summary statistics for all other variables, including all the variables preceding *Var8* in the set of variables defined by `Var5 TO Var8`.

## ***VARINFO Subcommand***

The optional `VARINFO` subcommand allows you to control the variable information included in the results.

- By default, all available variable information, with the exception of reserved system attributes, is included.
- If you include the `VARINFO` subcommand, it should be followed by one or more of the available keywords that indicate the dictionary information to include.

The available options are:

**POSITION.** *File position.* An integer that represents the position of the variable in file order. This is not available for multiple response sets.

**LABEL.** *Defined variable label.* For more information, see the topic [VARIABLE LABELS](#) on p. 2130.

**TYPE.** *Fundamental data type.* This is either *Numeric*, *String*, or *Multiple Response Set*.

**FORMAT.** *Print format.* The display format for the variable, such as *A4*, *F8.2*, or *DATE11*. For more information, see the topic [Variable Types and Formats](#) on p. 54. This is not available for multiple response sets.

**MEASURE.** *Measurement level.* The possible values are *Nominal*, *Ordinal*, *Scale*, and *Unknown*. The value displayed is the measurement level stored in the dictionary and is not affected by any temporary measurement level override specified by the `CODEBOOK` command. For more information, see the topic [VARIABLE LEVEL](#) on p. 2132. This is not available for multiple response sets. (Note: The measurement level for numeric variables may be “unknown” prior to the first data pass when the measurement level has not been explicitly set, such as data read from an external source or newly created variables. The measurement level for string variables is always known.)

**ROLE.** *Role when using predefined roles in dialogs.* Some dialogs support predefined roles that can be used to pre-select variables for analysis. For more information, see the topic [Overview](#) on p. 2133.

**ATTRIBUTES.** *User-defined custom variable attributes.* Output includes both the names and values for any custom variable attributes associated with each variable. For more information, see the topic [VARIABLE ATTRIBUTE](#) on p. 2127. This is not available for multiple response sets.

**VALUELABELS.** *Defined value labels.* If the `STATISTICS` subcommand includes `COUNT` or `PERCENT`, defined value labels are included in the output even if this keyword is omitted from the `VARINFO` subcommand. For information on defining value labels, see [VALUE LABELS](#).

**MISSING.** *User-defined missing values.* If the `STATISTICS` subcommand includes `COUNT` or `PERCENT`, defined value labels are included in the output even if this keyword is omitted from the `VARINFO` subcommand. For information on defining missing values, see [MISSING VALUES](#). This is not available for multiple response sets.

**RESERVEDATTRIBUTES.** *Reserved system variable attributes.* You can display system attributes, but you should not alter them. System attribute names start with a dollar sign (\$) . Non-display attributes, with names that begin with either “@” or “\$@”, are not included. Output includes both the names and values for any system attributes associated with each variable. This is not available for multiple response sets.

### **Example**

```
CODEBOOK  
  /VARINFO LABEL MEASURE VALUELABELS MISSING.
```

## **FILEINFO Subcommand**

The optional `FILEINFO` subcommand allows you to control the file information included in the results.

- By default, no file information is included.
- If you include the `FILEINFO` subcommand, it should be followed by one or more of the available keywords that indicate the file information to include.

The available options are:

**NAME.** *Name of the SPSS Statistics data file.* If the dataset has never been saved in IBM® SPSS® Statistics format, then there is no data file name.

**LOCATION.** *Directory (folder) location of the SPSS Statistics data file.* If the dataset has never been saved in SPSS Statistics format, then there is no location.

**CASECOUNT.** *Number of cases in the active dataset.* This is the total number of cases, including any cases that may be excluded from summary statistics due to filter conditions.

**LABEL.** *File label.* For more information, see the topic [FILE LABEL](#) on p. 725.

**ATTRIBUTES.** *User-defined custom data file attributes.* For more information, see the topic [DATAFILE ATTRIBUTE](#) on p. 569.

**DOCUMENTS.** *Data file document text.* Document text created with the `DOCUMENT` or `ADD DOCUMENT` commands. For more information, see the topic [ADD DOCUMENT](#) on p. 135.

**WEIGHT.** *Weight status* If weighting is on, the name of the weight variable is displayed.

**RESERVEDATTRIBUTES.** *Reserved system data file attributes.* You can display system attributes, but you should not alter them. System attribute names start with a dollar sign (\$) . Non-display attributes, with names that begin with either “@” or “\$@”, are not included. Output includes both the names and values for any system data file attributes.

### **Example**

```
CODEBOOK
  /FILEINFO NAME LOCATION CASECOUNT.
```

- The file information table will only include the name and location of the SPSS Statistics data file and the total number of cases in the file.
- The output will also include a table of variable information and summary statistics for each variable and multiple response set, because all variable information is included by default.

## **STATISTICS Subcommand**

The optional `STATISTICS` subcommand allows you to control the summary statistics that are included in the output, or suppress the display of summary statistics entirely.

- By default, all summary statistics are included.
- If you include the `STATISTICS` subcommand, it should be followed by one or more of the available keywords.

The available options are:

**COUNT.** *Number of cases in each category.* This applies to nominal and ordinal variables, multiple response sets, and labeled values of scale variables.

**PERCENT.** *Percent of cases in each category.* This applies to nominal and ordinal variables, multiple response sets, and labeled values of scale variables. The denominator is the total number of cases, including cases with missing values for the variable. If filtering is in effect, it is the total number of unfiltered cases. For multiple response sets, percentages can sum to more than 100%.

**MEAN.** *Mean.* This applies to scale variables only.

**STDDEV.** *Standard deviation.* This applies to scale variables only.

**QUANTILES.** *25th, 50th (median), and 75th percentiles.* This applies to scale variables only.

**NONE.** *Do not include any summary statistics.* If specified, this can be the only keyword included on the `STATISTICS` subcommand.

### **Example**

```
CODEBOOK  
  /STATISTICS COUNT MEAN.
```

## **OPTIONS Subcommand**

The `OPTIONS` subcommand allows you to suppress the display of value labels, counts, and percents for variables with more than a specified number of values or value labels and control the order in which variables are displayed in the output.

- If the `CODEBOOK` command does not include a variable list, the default display order is ascending file order.
- If the `CODEBOOK` command includes a variable list, the default display order is the order in which the variables are listed on the command.

### **MAXCATS Keyword**

**MAXCATS=N.** *Suppress counts and percents for variables with more than the specified number of valid values.* The default is 200. This applies to nominal and ordinal variables, multiple response sets, and labeled values of scale variables. For scale variables, the number of categories is the number of labeled values.

### **VARORDER Keyword**

The optional `VARORDER` keyword is followed by an equals sign (=) and one of the following alternatives:

**FILE.** *File order.*

**ALPHA.** *Alphabetic order by variable name.*

**VARLIST.** *Order in which variables and multiple response sets are listed on the command.* If there is no variable list, this setting is ignored, and the default file order is used.

**MEASURE.** *Sort by measurement level.* This creates four sorting groups: nominal, ordinal, scale, and unknown. (Note: The measurement level for numeric variables may be “unknown” prior to the first data pass when the measurement level has not been explicitly set, such as data read from

an external source or newly created variables. The measurement level for string variables is always known.)

**ATTRIBUTE (name).** *Alphabetic order by user-defined custom attribute name and value.* In ascending order, variables that don't have the attribute sort to the top, followed by variables that have the attribute but no defined value for the attribute, followed by variables with defined values for the attribute in alphabetic order of the values. [For more information, see the topic VARIABLE ATTRIBUTE on p. 2127.](#)

### ***SORT Keyword***

The optional SORT keyword is followed by an equals sign (=) and one of the following alternatives:

**ASCENDING.** *Ascending order.*

**DESCENDING.** *Descending order.*

### ***Example***

```
CODEBOOK  
  /OPTIONS MAXCATS=50 VARORDER=ALPHA SORT=DESCENDING.
```



# COMMENT

```
{COMMENT} text  
{ * }
```

## Overview

COMMENT inserts explanatory text within the command sequence. Comments are included among the commands printed back in the output; they do not become part of the information saved in a data file. To include commentary in the dictionary of a data file, use the DOCUMENT command.

### Syntax Rules

- The first line of a comment can begin with the keyword COMMENT or with an asterisk (\*). Comment text can extend for multiple lines and can contain any characters. A period is required at the end of the last line to terminate the comment.
- Use /\* and \*/ to set off a comment within a command. The comment can be placed wherever a blank is valid (except within strings) and should be preceded by a blank. Comments within a command cannot be continued onto the next line.
- The closing \*/ is optional when the comment is at the end of the line. The command can continue onto the next line just as if the inserted comment was a blank.
- Comments cannot be inserted within data lines.
- A comment on a separate line by itself within a command will cause an error. The comment line will be interpreted as a blank line, which is interpreted as a command terminator.

## Examples

### Comment As a Separate Command

```
* Create a new variable as a combination of two old variables;  
the new variable is a scratch variable used later in the  
session; it will not be saved with the data file.
```

```
COMPUTE #XYVAR=0.  
IF (XVAR EQ 1 AND YVAR EQ 1) #XYVAR=1.
```

The three-line comment will be included in the display file but will not be part of the data file if the active dataset is saved.

### Comments within Commands

```
IF (RACE EQ 1 AND GENDER EQ 1) GENDERRACE = 1. /*White males.
```

The comment is entered on a command line. The closing \*/ is not needed because the comment is at the end of the line.

***Comment on Separate Line within a Command***

```
FREQUENCIES VARIABLES=Var1 to Var5
/*this will cause an error*/
/FORMAT=NOTABLE
/BARCHART.
```

A comment on a separate line within a command will cause an error. The comment is interpreted as a blank line, and a blank line is interpreted as a command terminator. So `/FORMAT=NOTABLE` will be interpreted as the start of a different command, resulting in an error.

# COMPUTE

```
COMPUTE target variable=expression
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
COMPUTE newvar1=var1+var2.  
COMPUTE newvar2=RND(MEAN(var1 to var4)).  
COMPUTE logicalVar=(var1>5).  
STRING newString (A10).  
COMPUTE newString=CONCAT((RTRIM(stringVar1), stringVar2)).
```

Functions and operators available for COMPUTE are described in Transformation Expressions on p. 68.

## **Overview**

COMPUTE creates new numeric variables or modifies the values of existing string or numeric variables. The variable named on the left of the equals sign is the **target variable**. The variables, constants, and functions on the right side of the equals sign form an **assignment expression**. For a complete discussion of functions, see Transformation Expressions on p. 68.

### **Numeric Transformations**

Numeric variables can be created or modified with COMPUTE. The assignment expression for numeric transformations can include combinations of constants, variables, numeric operators, and functions.

### **String transformations**

String variables can be modified but cannot be created with COMPUTE. However, a new string variable can be declared and assigned a width with the STRING command and then assigned values by COMPUTE. The assignment expression can include string constants, string variables, and any of the string functions. All other functions are available for numeric transformations only.

### **Basic specification**

The basic specification is a target variable, an equals sign (required), and an assignment expression.

## **Syntax rules**

- The target variable must be named first, and the equals sign is required. Only one target variable is allowed per COMPUTE command.

- If the target variable is numeric, the expression must yield a numeric value; if the target variable is a string, the expression must yield a string value.
- Each function must specify at least one argument enclosed in parentheses. If a function has two or more arguments, the arguments must be separated by commas. For a complete discussion of functions and their arguments, see Transformation Expressions on p. 68.
- You can use the `TO` keyword to refer to a set of variables where the argument is a list of variables.

### ***Numeric variables***

- Parentheses are used to indicate the order of execution and to set off the arguments to a function.
- Numeric functions use simple or complex expressions as arguments. Expressions must be enclosed in parentheses.

### ***String variables***

- String values and constants must be enclosed in single or double quotes.
- When strings of different lengths are compared using the `ANY` or `RANGE` functions, the shorter string is right-padded with blanks so that its length equals that of the longer string.

### ***Operations***

- If the target variable already exists, its values are replaced.
- If the target variable does not exist and the assignment expression is numeric, the program creates a new variable.
- If the target variable does not exist and the assignment expression is a string, the program displays an error message and does not execute the command. Use the `STRING` command to declare new string variables before using them as target variables.

### ***Numeric variables***

- New numeric variables created with `COMPUTE` are assigned a dictionary format of `F8.2` and are initialized to the system-missing value for each case (unless the `LEAVE` command is used). Existing numeric variables transformed with `COMPUTE` retain their original dictionary formats. The format of a numeric variable can be changed with the `FORMATS` command.
- All expressions are evaluated in the following order: first functions, then exponentiation, and then arithmetic operations. The order of operations can be changed with parentheses.
- `COMPUTE` returns the system-missing value when it doesn't have enough information to evaluate a function properly. Arithmetic functions that take only one argument cannot be evaluated if that argument is missing. The date and time functions cannot be evaluated if any argument is missing. Statistical functions are evaluated if a sufficient number of arguments is valid. For example, in the command

```
COMPUTE FACTOR = SCORE1 + SCORE2 + SCORE3.
```

*FACTOR* is assigned the system-missing value for a case if any of the three score values is missing. It is assigned a valid value only when all score values are valid. In the command

```
COMPUTE FACTOR = SUM(SCORE1 TO SCORE3) .
```

*FACTOR* is assigned a valid value if at least one score value is valid. It is system-missing only when all three score values are missing. See [Missing values in numeric expressions](#) for information on how to control the minimum number of non-missing arguments required to return a non-missing result.

## String variables

- String variables can be modified but not created on `COMPUTE`. However, a new string variable can be created and assigned a width with the `STRING` command and then assigned new values with `COMPUTE`.
- Existing string variables transformed with `COMPUTE` retain their original dictionary formats. String variables declared on `STRING` and transformed with `COMPUTE` retain the formats assigned to them on `STRING`.
- The format of string variables cannot be changed with `FORMATS`. Instead, use `STRING` to create a new variable with the desired width and then use `COMPUTE` to set the values of the new string equal to the values of the original.
- The string returned by a string expression does not have to be the same width as the target variable. If the target variable is shorter, the result is right-trimmed. If the target variable is longer, the result is right-padded. The program displays no warning messages when trimming or padding.
- To control the width of strings, use the functions that are available for padding (`LPAD`, `RPAD`), trimming (`LTRIM`, `RTRIM`), and selecting a portion of strings (`SUBSTR`).
- To determine whether a character in a string is single-byte or double-byte, use the `MBLEN.BYTE` function. Specify the string and, optionally, its beginning byte position. If the position is not specified, it defaults to 1.

For more information, see the topic [String functions](#) on p. 106.

## Examples

A number of examples are provided to illustrate the use of `COMPUTE`. For a complete list of available functions and detailed function descriptions, see [Transformation Expressions](#).

### Arithmetic operations

```
COMPUTE V1=25-V2 .
COMPUTE V3=(V2/V4)*100 .

DO IF Tenure GT 5 .
COMPUTE Raise=Salary*.12 .
ELSE IF Tenure GT 1 .
COMPUTE Raise=Salary*.1 .
ELSE .
COMPUTE Raise=0 .
END IF .
```

## COMPUTE

- *V1* is 25 minus *V2* for all cases. *V3* is *V2* expressed as a percentage of *V4*.
- *Raise* is 12% of *Salary* if *Tenure* is greater than 5. For remaining cases, *Raise* is 10% of *Salary* if *Tenure* is greater than 1. For all other cases, *Raise* is 0.

**Arithmetic functions**

```
COMPUTE WtChange=ABS(Weight1-Weight2).
COMPUTE NewVar=RND((V1/V2)*100).
COMPUTE Income=TRUNC(Income).
COMPUTE MinSqrt=SQRT(MIN(V1,V2,V3,V4)).
```

```
COMPUTE Test = TRUNC(SQRT(X/Y)) * .5.
COMPUTE Parens = TRUNC(SQRT(X/Y) * .5).
```

- *WtChange* is the absolute value of *Weight1* minus *Weight2*.
- *NewVar* is the percentage *V1* is of *V2*, rounded to an integer.
- *Income* is truncated to an integer.
- *MinSqrt* is the square root of the minimum value of the four variables *V1* to *V4*. `MIN` determines the minimum value of the four variables, and `SQRT` computes the square root.
- The last two examples above illustrate the use of parentheses to control the order of execution. For a case with value 2 for *X* and *Y*, *Test* equals 0.5, since 2 divided by 2 (*X/Y*) is 1, the square root of 1 is 1, truncating 1 returns 1, and 1 times 0.5 is 0.5. However, *Parens* equals 0 for the same case, since `SQRT(X/Y)` is 1, 1 times 0.5 is 0.5, and truncating 0.5 returns 0.

**Statistical functions**

```
COMPUTE NewSalary = SUM(Salary,Raise).
COMPUTE MinValue = MIN(V1,V2,V3,V4).
COMPUTE MeanValue = MEAN(V1,V2,V3,V4).
COMPUTE NewMean = MEAN.3(V1,V2,V3,V4).
```

- *NewSalary* is the sum of *Salary* plus *Raise*.
- *MinValue* is the minimum of the values for *V1* to *V4*.
- *MeanValue* is the mean of the values for *V1* to *V4*. Since the mean can be computed for one, two, three, or four values, *MeanValue* is assigned a valid value as long as any one of the four variables has a valid value for that case.
- In the last example above, the `.3` suffix specifies the minimum number of valid arguments required. *NewMean* is the mean of variables *V1* to *V4* *only* if at least three of these variables have valid values. Otherwise, *NewMean* is system-missing for that case.

**Missing-Value functions**

```
MISSING VALUE V1 V2 V3 (0).
COMPUTE AllValid=V1 + V2 + V3.
COMPUTE UM=VALUE(V1) + VALUE(V2) + VALUE(V3).
COMPUTE SM=SYSMIS(V1) + SYSMIS(V2) + SYSMIS(V3).
COMPUTE M=MISSING(V1) + MISSING(V2) + MISSING(V3).
```

- The `MISSING VALUE` command declares the value 0 as missing for *V1*, *V2*, and *V3*.

- *AllValid* is the sum of three variables only for cases with valid values for all three variables. *AllValid* is assigned the system-missing value for a case if any variable in the assignment expression has a system- or user-missing value.
- The VALUE function overrides user-missing value status. Thus, *UM* is the sum of *V1*, *V2*, and *V3* for each case, including cases with the value 0 (the user-missing value) for any of the three variables. Cases with the system-missing value for *V1*, *V2*, and *V3* are system-missing.
- The SYSMIS function on the third COMPUTE returns the value 1 if the variable is system-missing. Thus, *SM* ranges from 0 to 3 for each case, depending on whether the variables *V1*, *V2*, and *V3* are system-missing for that case.
- The MISSING function on the fourth COMPUTE returns the value 1 if the variable named is system- or user-missing. Thus, *M* ranges from 0 to 3 for each case, depending on whether the variables *V1*, *V2*, and *V3* are user- or system-missing for that case.
- Alternatively, you could use the COUNT command to create the variables *SM* and *M*.

\* Test for listwise deletion of missing values.

```
DATA LIST /V1 TO V6 1-6.
BEGIN DATA
213 56
123457
123457
9234 6
END DATA.
MISSING VALUES V1 TO V6(6,9).

COMPUTE NotValid=NMISS(V1 TO V6).
FREQUENCIES VAR=NotValid.
```

- COMPUTE determines the number of missing values for each case. For each case without missing values, the value of *NotValid* is 0. For each case with one missing value, the value of *NotValid* is 1, and so on. Both system- and user-missing values are counted.
- FREQUENCIES generates a frequency table for *NotValid*. The table gives a count of how many cases have all valid values, how many cases have one missing value, how many cases have two missing values, and so on, for variables *V1* to *V6*. This table can be used to determine how many cases would be dropped in an analysis that uses listwise deletion of missing values. For other ways to check listwise deletion, see the examples for the ELSE command (in the DO IF command) and those for the IF command.

For more information, see the topic [Missing value functions](#) on p. 122.

## String functions

```
DATA LIST FREE / FullName (A20).
BEGIN DATA
"Fred Smith"
END DATA.
STRING FirstName LastName LastFirstName (A20).
COMPUTE #spaceLoc=INDEX(FullName, " ").
COMPUTE FirstName=SUBSTR(FullName, 1, (#spaceLoc-1)).
COMPUTE LastName=SUBSTR(FullName, (#spaceLoc+1)).
COMPUTE LastFirstName=CONCAT(RTRIM(LastName), ", ", FirstName).
COMPUTE LastFirstName=REPLACE(LastFirstName, "Fred", "Ted").
```

---

*COMPUTE*

- The INDEX function returns a number that represents the location of the first blank space in the value of the string variable *FullName*.
- The first SUBSTR function sets *FirstName* to the portion of *FullName* prior to the first space in the value. So, in this example, the value of *FirstName* is “Fred”.
- The second SUBSTR function sets *LastName* to the portion of *FullName* after the first blank space in the value. So, in this example, the value of *LastName* is “Smith”.
- The CONCAT function combines the values of *LastName* and *FirstName*, with a comma and a space between the two values. So, in this example, the value of *LastFirstName* is “Smith, Fred”. Since all string values are right-padded with blank spaces to the defined width of the string variable, the RTRIM function is needed to remove all the extra blank spaces from *LastName*.
- The REPLACE function changes any instances of the string “Fred” in *LastFirstName* to “Ted”. So, in this example, the value of *LastFirstName* is changed to “Smith, Ted”.

For more information, see the topic [String functions](#) on p. 106.

### ***Scoring functions***

```
STRING SPECIES(A20) .  
COMPUTE SCOREPROB=ApplyModel(CREDITMOD1, 'PROBABILIT') .  
COMPUTE SPECIES=StrApplyModel(QUESTMOD1, 'PREDICT') .
```

- *SCOREPROB* is the probability that the value predicted from the model specified by CREDITMOD1 is correct.
- *SPECIES* is the predicted result from the model specified by QUESTMOD1 as applied to the active dataset. The prediction is returned as a string value.



# CONJOINT

CONJOINT is available in the Conjoint option.

```
CONJOINT  [PLAN={*                }]
           {'savfile'|'dataset'}

[/DATA={*                }]
         {'savfile'|'dataset'}

/{SEQUENCE}=varlist
  {RANK      }
  {SCORE     }

[/SUBJECT=variable]

[/FACTORS=varlist['labels'] ({{DISCRETE[{MORE}]}}
                             {      {LESS} }
                             {LINEAR[{MORE}] }
                             {      {LESS} }
                             {IDEAL      }
                             {ANTIIDEAL  }
                             [values['labels']])]

varlist...

[/PRINT={ALL**          } [SUMMARYONLY]]
        {ANALYSIS      }
        {SIMULATION    }
        {NONE          }

[/UTILITY=file]

[/PLOT={ [SUMMARY] [SUBJECT] [ALL]]}
       { [NONE**]          }]
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example:

```
CONJOINT PLAN=' /DATA/CARPLAN.SAV'
  /FACTORS=SPEED (LINEAR MORE) WARRANTY (DISCRETE MORE)
  PRICE (LINEAR LESS) SEATS
  /SUBJECT=SUBJ /RANK=RANK1 TO RANK15 /UTILITY='UTIL.SAV'.
```

## Overview

CONJOINT analyzes score or rank data from full-concept conjoint studies. A plan file that is generated by ORTHOPLAN or entered by the user describes the set of full concepts that are scored or ranked in terms of preference. A variety of continuous and discrete models is available to estimate utilities for each individual subject and for the group. Simulation estimates for concepts that are not rated can also be computed.

### **Options**

**Data Input.** You can analyze data recorded as rankings of an ordered set of profiles (or cards) as the profile numbers arranged in rank order, or as preference scores of an ordered set of profiles.

**Model Specification.** You can specify how each factor is expected to be related to the scores or ranks.

**Display Output.** The output can include the analysis of the experimental data, results of simulation data, or both.

**Writing an External File.** An data file containing utility estimates and associated statistics for each subject can be written for use in further analyses or graphs.

### **Basic Specification**

- The basic specification is `CONJOINT`, a `PLAN` or `DATA` subcommand, and a `SEQUENCE`, `RANK`, or `SCORE` subcommand to describe the type of data.
- `CONJOINT` requires two files: a plan file and a data file. If only the `PLAN` subcommand or the `DATA` subcommand—but not both—is specified, `CONJOINT` will read the file that is specified on the `PLAN` or `DATA` subcommand and use the active dataset as the other file.
- By default, estimates are computed by using the `DISCRETE` model for all variables in the plan file (except those named `STATUS_` and `CARD_`). Output includes Kendall's tau and Pearson's product-moment correlation coefficients measuring the relationship between predicted scores and actual scores. Significance levels for one-tailed tests are displayed.

### **Subcommand Order**

- Subcommands can appear in any order.

### **Syntax Rules**

- Multiple `FACTORS` subcommands are all executed. For all other subcommands, only the last occurrence is executed.

### **Operations**

- Both the plan and data files can be external IBM® SPSS® Statistics data files. In this case, `CONJOINT` can be used before an active dataset is defined.
- The variable `STATUS_` in the plan file must equal 0 for experimental profiles, 1 for holdout profiles, and 2 for simulation profiles. Holdout profiles are judged by the subjects but are not used when `CONJOINT` estimates utilities. Instead, these profiles are used as a check on the validity of the estimated utilities. Simulation profiles are factor-level combinations that are not rated by the subjects but are estimated by `CONJOINT` based on the ratings of the experimental profiles. If there is no `STATUS_` variable, all profiles in the plan file are assumed to be experimental profiles.
- All variables in the plan file except `STATUS_` and `CARD_` are used by `CONJOINT` as factors.
- In addition to the estimates for each individual subject, average estimates for each split-file group that is identified in the data file are computed. The plan file cannot have a split-file structure.

- Factors are tested for orthogonality by CONJOINT. If all of the factors are not orthogonal, a matrix of Cramér's  $V$  statistics is displayed to describe the non-orthogonality.
- When SEQUENCE or RANK data are used, CONJOINT internally reverses the ranking scale so that the computed coefficients are positive.
- The plan file cannot be sorted or modified in any way after the data are collected, because the sequence of profiles in the plan file must match the sequence of values in the data file in a one-to-one correspondence. (CONJOINT uses the order of profiles as they appear in the plan file, not the value of *CARD\_*, to determine profile order.) If RANK or SCORE is the data-recording method, the first response from the first subject in the data file is the rank or score of the first profile in the plan file. If SEQUENCE is the data-recording method, the first response from the first subject in the data file is the profile number (determined by the order of profiles in the plan file) of the most preferred profile.

### Limitations

- Factors must be numeric.
- The plan file cannot contain missing values or case weights. In the active dataset, profiles with missing values on the SUBJECT variable are grouped together and averaged at the end. If any preference data (the ranks, scores, or profile numbers) are missing, that subject is skipped.
- Factors must have at least two levels. The maximum number of levels for each factor is 99. Note that ORTHOPLAN will only produce plans with factors with 9 or fewer levels for each factor.

## Examples

```
CONJOINT PLAN=' /DATA/CARPLAN.SAV'
  /FACTORS=SPEED (LINEAR MORE) WARRANTY (DISCRETE MORE)
  PRICE (LINEAR LESS) SEATS
  /SUBJECT=SUBJ /RANK=RANK1 TO RANK15 /UTILITY='UTIL.SAV'.
```

- The PLAN subcommand specifies the IBM® SPSS® Statistics data file *CARPLAN.SAV* as the plan file containing the full-concept profiles. Because there is no DATA subcommand, the active dataset is assumed to contain the subjects' rankings of these profiles.
- The FACTORS subcommand specifies the ways in which the factors are expected to be related to the rankings. For example, speed is expected to be linearly related to the rankings, so that cars with higher speeds will receive lower (more-preferred) rankings.
- The SUBJECT subcommand specifies the variable *SUBJ* in the active dataset as an identification variable. All consecutive cases with the same value on this variable are combined to estimate utilities.
- The RANK subcommand specifies that each data point is a ranking of a specific profile and identifies the variables in the active dataset that contain these rankings.
- UTILITY writes out an external data file named *UTIL.SAV* containing the utility estimates and associated statistics for each subject.

## PLAN Subcommand

PLAN identifies the file containing the full-concept profiles.

- `PLAN` is followed by quoted file specification for an external IBM® SPSS® Statistics data file or currently open dataset containing the plan. An asterisk instead of a file specification indicates the active dataset.
- If the `PLAN` subcommand is omitted, the active dataset is assumed by default. However, you must specify at least one SPSS Statistics data file or dataset on a `PLAN` or `DATA` subcommand. The active dataset cannot be specified as both the plan file and data file.
- The plan file is a specially prepared file that is generated by `ORTHOPLAN` or entered by the user. The plan file can contain the variables `CARD_` and `STATUS_`, and it must contain the factors of the conjoint study. The value of `CARD_` is a profile identification number. The value of `STATUS_` is 0, 1, or 2, depending on whether the profile is an experimental profile (0), a holdout profile (1), or a simulation profile (2).
- The sequence of the profiles in the plan file must match the sequence of values in the data file.
- Any simulation profiles (`STATUS_=2`) must follow experimental and holdout profiles in the plan file.
- All variables in the plan file except `CARD_` and `STATUS_` are used as factors by `CONJOINT`.

### Example

```
DATA LIST FREE /CARD_ WARRANTY SEATS PRICE SPEED STATUS_ .
BEGIN DATA
1 1 4 14000 130 2
2 1 4 14000 100 2
3 3 4 14000 130 2
4 3 4 14000 100 2
END DATA.
ADD FILES FILE=' /DATA/CARPLAN.SAV' /FILE=*.
CONJOINT PLAN=* /DATA=' /DATA/CARDATA.SAV'
  /FACTORS=PRICE (ANTIIDEAL) SPEED (LINEAR) WARRANTY (DISCRETE MORE)
  /SUBJECT=SUBJ /RANK=RANK1 TO RANK15 /PRINT=SIMULATION.
```

- `DATA LIST` defines six variables—a `CARD_` identification variable, four factors, and a `STATUS_` variable.
- The data between `BEGIN DATA` and `END DATA` are four simulation profiles. Each profile contains a `CARD_` identification number and the specific combination of factor levels of interest.
- The variable `STATUS_` is equal to 2 for all cases (profiles). `CONJOINT` interprets profiles with `STATUS_` equal to 2 as simulation profiles.
- The `ADD FILES` command joins an old plan file, `CARPLAN.SAV`, with the active dataset. Note that the active dataset is indicated last on the `ADD FILES` command so that the simulation profiles are appended to the end of `CARPLAN.SAV`.
- The `PLAN` subcommand on `CONJOINT` defines the new active dataset as the plan file. The `DATA` subcommand specifies a data file from a previous `CONJOINT` analysis.

## DATA Subcommand

`DATA` identifies the file containing the subjects' preference scores or rankings.

- DATA is followed by a quoted file specification for an external IBM® SPSS® Statistics data file or a currently open dataset containing the data. An asterisk instead of a file specification indicates the active dataset.
- If the DATA subcommand is omitted, the active dataset is assumed by default. However, you must specify at least one SPSS Statistics data file on a DATA or PLAN subcommand. The active dataset cannot be specified as both the plan file and data file.
- One variable in the data file can be a subject identification variable. All other variables are the subject responses and are equal in number to the number of experimental and holdout profiles in the plan file.
- The subject responses can be in the form of ranks assigned to an ordered sequence of profiles, scores assigned to an ordered sequence of profiles, or profile numbers in preference order from most liked to least liked.
- Tied ranks or scores are allowed. If tied ranks are present, CONJOINT issues a warning and then proceeds with the analysis. Data recorded in SEQUENCE format, however, cannot have ties, because each profile number must be unique.

### Example

```

DATA LIST FREE /SUBJ RANK1 TO RANK15.
BEGIN DATA
01 3 7 6 1 2 4 9 12 15 13 14 5 8 10 11
02 7 3 4 9 6 15 10 13 5 11 1 8 4 2 12
03 12 13 5 1 14 8 11 2 7 6 3 4 15 9 10
04 3 6 7 4 2 1 9 12 15 11 14 5 8 10 13
05 9 3 4 7 6 10 15 13 5 12 1 8 4 2 11
50 12 13 8 1 14 5 11 6 7 2 3 4 15 10 9
END DATA.
SAVE OUTFILE=' /DATA/RANKINGS.SAV' .
DATA LIST FREE /CARD_ WARRANTY SEATS PRICE SPEED.
BEGIN DATA
 1 1 4 14000 130
 2 1 4 14000 100
 3 3 4 14000 130
 4 3 4 14000 100
 5 5 2 10000 130
 6 1 4 10000 070
 7 3 4 10000 070
 8 5 2 10000 100
 9 1 4 07000 130
10 1 4 07000 100
11 5 2 07000 070
12 5 4 07000 070
13 1 4 07000 070
14 5 2 10000 070
15 5 2 14000 130
END DATA.
CONJOINT PLAN=* /DATA=' /DATA/RANKINGS.SAV'
 /FACTORS=PRICE (ANTIIDEAL) SPEED (LINEAR)
 WARRANTY (DISCRETE MORE)
 /SUBJECT=SUBJ /RANK=RANK1 TO RANK15.

```

- The first set of DATA LIST and BEGIN-END DATA commands creates a data file containing the rankings. This file is saved in the external file *RANKINGS.SAV*.

- The second set of `DATA LIST` and `BEGIN-END DATA` commands defines the plan file as the active dataset.
- The `CONJOINT` command uses the active dataset as the plan file and uses `RANKINGS.SAV` as the data file.

## ***SEQUENCE, RANK, or SCORE Subcommand***

The `SEQUENCE`, `RANK`, or `SCORE` subcommand is specified to indicate the way in which the preference data were recorded.

<b>SEQUENCE</b>	<i>Each data point in the data file is a profile number, starting with the most-preferred profile and ending with the least-preferred profile. This is how the data are recorded if the subject is asked to order the deck of profiles from most preferred to least preferred. The researcher records which profile number was first, which profile number was second, and so on.</i>
<b>RANK</b>	<i>Each data point is a ranking, starting with the ranking of profile 1, then the ranking of profile 2, and so on. This is how the data are recorded if the subject is asked to assign a rank to each profile, ranging from 1 to <math>n</math>, where <math>n</math> is the number of profiles. A lower rank implies greater preference.</i>
<b>SCORE</b>	<i>Each data point is a preference score assigned to the profiles, starting with the score of profile 1, then the score of profile 2, and so on. These types of data might be generated, for example, by asking subjects to use a Likert scale to assign a score to each profile or by asking subjects to assign a number from 1 to 100 to show how much they like the profile. A higher score implies greater preference.</i>

- You must specify one, and only one, of these three subcommands.
- After each subcommand, the names of the variables containing the preference data (the profile numbers, ranks, or scores) are listed. There must be as many variable names listed as there are experimental and holdout profiles in the plan file.

### ***Example***

```
CONJOINT PLAN=* /DATA='DATA.SAV'
  /FACTORS=PRICE (ANTIIDEAL) SPEED (LINEAR) WARRANTY (DISCRETE MORE)
  /SUBJECT=SUBJ
  /RANK=RANK1 TO RANK15.
```

- The `RANK` subcommand indicates that the data are rankings of an ordered sequence of profiles. The first data point after `SUBJ` is variable `RANK1`, which is the ranking that is given by subject 1 to profile 1.
- There are 15 profiles in the plan file, so there must be 15 variables listed on the `RANK` subcommand.
- The example uses the `TO` keyword to refer to the 15 rank variables.

## ***SUBJECT Subcommand***

`SUBJECT` specifies an identification variable. All consecutive cases having the same value on this variable are combined to estimate the utilities.

- If SUBJECT is not specified, all data are assumed to come from one subject, and only a group summary is displayed.
- SUBJECT is followed by the name of a variable in the active dataset.
- If the same SUBJECT value appears later in the data file, it is treated as a different subject.

## ***FACTORS Subcommand***

FACTORS specifies the way in which each factor is expected to be related to the rankings or scores.

- If FACTORS is not specified, the DISCRETE model is assumed for all factors.
- All variables in the plan file except *CARD\_* and *STATUS\_* are used as factors, even if they are not specified on FACTORS.
- FACTORS is followed by a variable list and a model specification in parentheses that describes the expected relationship between scores or ranks and factor levels for that variable list.
- The model specification consists of a model name and, for the DISCRETE and LINEAR models, an optional MORE or LESS keyword to indicate the direction of the expected relationship. Values and value labels can also be specified.
- MORE and LESS keywords will *not* affect estimates of utilities. They are used simply to identify subjects whose estimates do not match the expected direction.

The four available models are as follows:

<b>DISCRETE</b>	<i>No assumption.</i> The factor levels are categorical, and no assumption is made about the relationship between the factor and the scores or ranks. This setting is the default. Specify keyword MORE after DISCRETE to indicate that higher levels of a factor are expected to be more preferred. Specify keyword LESS after DISCRETE to indicate that lower levels of a factor are expected to be more preferred.
<b>LINEAR</b>	<i>Linear relationship.</i> The scores or ranks are expected to be linearly related to the factor. Specify keyword MORE after LINEAR to indicate that higher levels of a factor are expected to be more preferred. Specify keyword LESS after LINEAR to indicate that lower levels of a factor are expected to be more preferred.
<b>IDEAL</b>	<i>Quadratic relationship, decreasing preference.</i> A quadratic relationship is expected between the scores or ranks and the factor. It is assumed that there is an ideal level for the factor, and distance from this ideal point, in either direction, is associated with decreasing preference. Factors that are described with this model should have at least three levels.
<b>ANTIIDEAL</b>	<i>Quadratic relationship, increasing preference.</i> A quadratic relationship is expected between the scores or ranks and the factor. It is assumed that there is a worst level for the factor, and distance from this point, in either direction, is associated with increasing preference. Factors that are described with this model should have at least three levels.

- The DISCRETE model is assumed for those variables that are not listed on the FACTORS subcommand.
- When a MORE or LESS keyword is used with DISCRETE or LINEAR, a reversal is noted when the expected direction does not occur.

- Both `IDEAL` and `ANTIIDEAL` create a quadratic function for the factor. The only difference is whether preference increases or decreases with distance from the point. The estimated utilities are the same for these two models. A reversal is noted when the expected model (`IDEAL` or `ANTIIDEAL`) does not occur.
- The optional value and value label lists allow you to recode data and/or replace value labels. The new values, in the order in which they appear on the value list, replace existing values, starting with the smallest existing value. If a new value is not specified for an existing value, the value remains unchanged.
- New value labels are specified in apostrophes or quotation marks. New values without new labels retain existing labels; new value labels without new values are assigned to values in the order in which they appear, starting with the smallest existing value.
- For each factor that is recoded, a table is displayed, showing the original and recoded values and the value labels.
- If the factor levels are coded in discrete categories (for example, 1, 2, 3), these values are the values used by `CONJOINT` in computations, even if the value labels contain the actual values (for example, 80, 100, 130). Value labels are never used in computations. You can recode the values as described above to change the coded values to the real values. Recoding does not affect `DISCRETE` factors but does change the coefficients of `LINEAR`, `IDEAL`, and `ANTIIDEAL` factors.
- In the output, variables are described in the following order:
  1. All `DISCRETE` variables in the order in which they appear on the `FACTORS` subcommand.
  2. All `LINEAR` variables in the order in which they appear on the `FACTORS` subcommand.
  3. All `IDEAL` and `ANTIIDEAL` factors in the order in which they appear on the `FACTORS` subcommand.

### **Example**

```
CONJOINT DATA='DATA.SAV'
  /FACTORS=PRICE (LINEAR LESS) SPEED (IDEAL 70 100 130)
  WARRANTY (DISCRETE MORE)
  /RANK=RANK1 TO RANK15.
```

- The `FACTORS` subcommand specifies the expected relationships. A linear relationship is expected between price and rankings, so that the higher the price, the lower the preference (higher ranks). A quadratic relationship is expected between speed levels and rankings, and longer warranties are expected to be associated with greater preference (lower ranks).
- The `SPEED` factor has a new value list. If the existing values were 1, 2, and 3, 70 replaces 1, 100 replaces 2, and 130 replaces 3.
- Any variable in the plan file (except `CARD_` and `STATUS_`) that is not listed on the `FACTORS` subcommand uses the `DISCRETE` model.

## **PRINT Subcommand**

`PRINT` controls whether your output includes the analysis of the experimental data, the results of the simulation data, both, or none.



The following keywords are available:

<b>ANALYSIS</b>	<i>Only the results of the experimental data analysis are included.</i>
<b>SIMULATION</b>	<i>Only the results of the simulation data analysis are included. The results of three simulation models—maximum utility, Bradley-Terry-Luce (BTL), and logit—are displayed.</i>
<b>SUMMARYONLY</b>	<i>Only the summaries in the output are included, not the individual subjects. Thus, if you have a large number of subjects, you can see the summary results without having to generate output for each subject.</i>
<b>ALL</b>	<i>The results of both the experimental data and simulation data analyses are included. ALL is the default.</i>
<b>NONE</b>	<i>No results are written to the display file. This keyword is useful if you are interested only in writing the utility file (see “UTILITY Subcommand” below).</i>

## ***UTILITY Subcommand***

UTILITY writes a utility file to the specified IBM® SPSS® Statistics file.

- If UTILITY is not specified, no utility file is written.
- UTILITY is followed by the name of the file to be written.
- The file is specified in the usual manner for your operating system.
- The utility file contains one case for each subject. If SUBJECT is not specified, the utility file contains a single case with statistics for the group as a whole.

The variables that are written to the utility file are in the following order:

- Any SPLIT FILE variables in the active dataset.
- Any SUBJECT variable.
- The constant for the regression equation for the subject. The regression equation constant is named CONSTANT.
- For DISCRETE factors, all of the utilities that are estimated for the subject. The names of the utilities that are estimated with DISCRETE factors are formed by appending a digit after the factor name. The first utility gets a 1, the second utility gets a 2, and so on.
- For LINEAR factors, a single coefficient. The name of the coefficient for LINEAR factors is formed by appending *\_L* to the factor name. (To calculate the predicted score, multiply the factor value by the coefficient.)
- For IDEAL or ANTIIDEAL factors, two coefficients. The name of the two coefficients for IDEAL or ANTIIDEAL factors are formed by appending *\_L* and *\_Q*, respectively, to the factor name. (To use these coefficients in calculating the predicted score, multiply the factor value by the first coefficient and add that to the product of the second coefficient and the square of the factor value.)
- The estimated ranks or scores for all profiles in the plan file. The names of the estimated ranks or scores are of the form *SCORE<sub>n</sub>* for experimental and holdout profiles, or *SIMUL<sub>n</sub>* for simulation profiles, where *n* is the position in the plan file. The name is *SCORE* for experimental and holdout profiles even if the data are ranks.

If the variable names that are created are too long, letters are truncated from the end of the original variable name before new suffixes are appended.

## ***PLOT Subcommand***

The PLOT subcommand produces plots in addition to the output that is usually produced by CONJOINT.

The following keywords are available for this subcommand:

<b>SUMMARY</b>	<i>Produces a bar chart of the importance values for all variables, plus a utility bar chart for each variable. This setting is the default if the PLOT subcommand is specified with no keywords.</i>
<b>SUBJECT</b>	<i>Plots a clustered bar chart of the importance values for each factor, clustered by subjects, and one clustered bar chart for each factor, showing the utilities for each factor level, clustered by subjects. If no SUBJECT subcommand was specified naming the variables, no plots are produced and a warning is displayed.</i>
<b>ALL</b>	<i>Plots both summary and subject charts.</i>
<b>NONE</b>	<i>Does not produce any charts. This setting is the default if the subcommand is omitted.</i>

# CORRELATIONS

CORRELATIONS is available in the Statistics Base option.

```
CORRELATIONS VARIABLES= varlist [WITH varlist] [/varlist...]  
  [/MISSING={PAIRWISE**}  [{INCLUDE}]]  
             {LISTWISE }   {EXCLUDE}  
  [/PRINT={TWOTAIL**}  {SIG**}]  
          {ONETAIL  }   {NOSIG}  
  [/MATRIX=OUT({*          })]  
             {'savfile'|'dataset'}  
  [/STATISTICS=[DESCRIPTIVES] [XPROD] [ALL]]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 16.0

- Added support for SET THREADS and SET MCACHE.

## Example

```
CORRELATIONS VARIABLES=FOOD RENT PUBTRANS TEACHER COOK ENGINEER  
  /MISSING=INCLUDE.
```

## Overview

CORRELATIONS (alias PEARSON CORR) produces Pearson product-moment correlations with significance levels and, optionally, univariate statistics, covariances, and cross-product deviations. Other procedures that produce correlation matrices are PARTIAL CORR, REGRESSION, DISCRIMINANT, and FACTOR.

## Options

**Types of Matrices.** A simple variable list on the VARIABLES subcommand produces a square matrix. You can also request a rectangular matrix of correlations between specific pairs of variables or between variable lists using the keyword WITH on VARIABLES.

**Significance Levels.** By default, CORRELATIONS displays the number of cases and significance levels for each coefficient. Significance levels are based on a two-tailed test. You can request a one-tailed test, and you can display the significance level for each coefficient as an annotation using the PRINT subcommand.

**Additional Statistics.** You can obtain the mean, standard deviation, and number of nonmissing cases for each variable, and the cross-product deviations and covariance for each pair of variables using the `STATISTICS` subcommand.

**Matrix Output.** You can write matrix materials to a data file using the `MATRIX` subcommand. The matrix materials include the mean, standard deviation, number of cases used to compute each coefficient, and Pearson correlation coefficient for each variable. The matrix data file can be read by several other procedures.

### **Basic Specification**

- The basic specification is the `VARIABLES` subcommand, which specifies the variables to be analyzed.
- By default, `CORRELATIONS` produces a matrix of correlation coefficients. The number of cases and the significance level are displayed for each coefficient. The significance level is based on a two-tailed test.

### **Subcommand Order**

- The `VARIABLES` subcommand must be first.
- The remaining subcommands can be specified in any order.

### **Operations**

- The correlation of a variable with itself is displayed as 1.0000.
- A correlation that cannot be computed is displayed as a period (.).
- `CORRELATIONS` does not execute if string variables are specified on the variable list.
- This procedure uses the multithreaded options specified by `SET THREADS` and `SET MCACHE`.

### **Limitations**

- A maximum of 40 variable lists.
- A maximum of 500 variables total per command.
- A maximum of 250 syntax elements. Each individual occurrence of a variable name, keyword, or special delimiter counts as 1 toward this total. Variables implied by the `TO` keyword do not count toward this total.

## **Example**

```
CORRELATIONS
  /VARIABLES=sales mpg
  /PRINT=TWOTAIL NOSIG
  /MISSING=PAIRWISE .
```

- The procedure produces a correlation matrix for *sales* and *mpg*.
- All other options are set to their default values.

## VARIABLES Subcommand

VARIABLES specifies the variable list.

- A simple variable list produces a square matrix of correlations of each variable with every other variable.
- Variable lists joined by the keyword WITH produce a rectangular correlation matrix. Variables before WITH define the rows of the matrix and variables after WITH define the columns.
- The keyword ALL can be used on the variable list to refer to all user-defined variables.
- You can specify multiple VARIABLES subcommands on a single CORRELATIONS command. The slash between the subcommands is required; the keyword VARIABLES is not.

### Example

```
CORRELATIONS VARIABLES=FOOD RENT PUBTRANS TEACHER COOK ENGINEER
/VARIABLES=FOOD RENT WITH COOK TEACHER MANAGER ENGINEER
/MISSING=INCLUDE.
```

- The first VARIABLES subcommand requests a square matrix of correlation coefficients among the variables *FOOD*, *RENT*, *PUBTRANS*, *TEACHER*, *COOK*, and *ENGINEER*.
- The second VARIABLES subcommand requests a rectangular correlation matrix in which *FOOD* and *RENT* are the row variables and *COOK*, *TEACHER*, *MANAGER*, and *ENGINEER* are the column variables.

## PRINT Subcommand

PRINT controls whether the significance level is based on a one- or two-tailed test and whether the number of cases and the significance level for each correlation coefficient are displayed.

<b>TWOTAIL</b>	<i>Two-tailed test of significance.</i> This test is appropriate when the direction of the relationship cannot be determined in advance, as is often the case in exploratory data analysis. This is the default.
<b>ONETAIL</b>	<i>One-tailed test of significance.</i> This test is appropriate when the direction of the relationship between a pair of variables can be specified in advance of the analysis.
<b>SIG</b>	<i>Do not flag significant values.</i> SIG is the default.
<b>NOSIG</b>	<i>Flag significant values.</i> Values significant at the 0.05 level are flagged with a single asterisk; those that are significant at the 0.01 level are flagged with two asterisks.

## STATISTICS Subcommand

The correlation coefficients are automatically displayed in the Correlations table for an analysis specified by a VARIABLES list. STATISTICS requests additional statistics.

<b>DESCRIPTIVES</b>	<i>Display mean, standard deviation, and number of nonmissing cases for each variable on the Variables list in the Descriptive Statistics table. This table precedes all Correlations tables. Variables specified on more than one VARIABLES list are displayed only once. Missing values are handled on a variable-by-variable basis regardless of the missing-value option in effect for the correlations.</i>
<b>XPROD</b>	<i>Display cross-product deviations and covariance for each pair of variables in the Correlations table(s).</i>
<b>ALL</b>	<i>All additional statistics. This produces the same statistics as DESCRIPTIVES and XPROD together.</i>

## MISSING Subcommand

MISSING controls the treatment of missing values.

- The PAIRWISE and LISTWISE keywords are alternatives; however, each can be specified with INCLUDE or EXCLUDE.
- The default is PAIRWISE and EXCLUDE.

<b>PAIRWISE</b>	<i>Exclude missing values pairwise. Cases that have missing values for one or both of a pair of variables for a specific correlation coefficient are excluded from the computation of that coefficient. Since each coefficient is based on all cases that have valid values for that particular pair of variables, this can result in a set of coefficients based on a varying number of cases. The valid number of cases is displayed in the Correlations table. This is the default.</i>
<b>LISTWISE</b>	<i>Exclude missing values listwise. Cases that have missing values for any variable named on any VARIABLES list are excluded from the computation of all coefficients across lists. The valid number of cases is the same for all analyses and is displayed in a single annotation.</i>
<b>INCLUDE</b>	<i>Include user-missing values. User-missing values are included in the analysis.</i>
<b>EXCLUDE</b>	<i>Exclude all missing values. Both user- and system-missing values are excluded from the analysis.</i>

## MATRIX Subcommand

MATRIX writes matrix materials to a data file or previously declared dataset (DATASET DECLARE command). The matrix materials include the mean and standard deviation for each variable, the number of cases used to compute each coefficient, and the Pearson correlation coefficients. Several procedures can read matrix materials produced by CORRELATIONS, including PARTIAL CORR, REGRESSION, FACTOR, and CLUSTER.

- CORRELATIONS cannot write rectangular matrices (those specified with the keyword WITH) to a file.
- If you specify more than one variable list on CORRELATIONS, only the last list that does not use the keyword WITH is written to the matrix data file.

- The keyword `OUT` specifies the file to which the matrix is written. Specify an asterisk to replace the active dataset or a quoted file specification or dataset name, enclosed in parentheses.
- Documents from the original file will not be included in the matrix file and will not be present if the matrix file becomes the working data file.

### **Format of the Matrix Data File**

- The matrix data file has two special variables created by the program: `ROWTYPE_` and `VARNAME_`. The variable `ROWTYPE_` is a short string variable with values `MEAN`, `STDDEV`, `N`, and `CORR` (for Pearson correlation coefficient). The next variable, `VARNAME_`, is a short string variable whose values are the names of the variables used to form the correlation matrix. When `ROWTYPE_` is `CORR`, `VARNAME_` gives the variable associated with that row of the correlation matrix.
- The remaining variables in the file are the variables used to form the correlation matrix.

### **Split Files**

- When split-file processing is in effect, the first variables in the matrix file will be split variables, followed by `ROWTYPE_`, `VARNAME_`, and the variables used to form the correlation matrix.
- A full set of matrix materials is written for each subgroup defined by the split variables.
- A split variable cannot have the same name as any other variable written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split-file specifications must be in effect when that matrix is read by another procedure.

### **Missing Values**

- With pairwise treatment of missing values (the default), a matrix of the number of cases used to compute each coefficient is included with the matrix materials.
- With listwise treatment, a single number indicating the number of cases used to calculate all coefficients is included.

### **Example**

```
GET FILE=CITY /KEEP FOOD RENT PUBTRANS TEACHER COOK ENGINEER.
CORRELATIONS VARIABLES=FOOD TO ENGINEER
/MATRIX OUT(CORRMAT) .
```

- `CORRELATIONS` reads data from the file `CITY` and writes one set of matrix materials to the file `CORRMAT`. The working file is still `CITY`. Subsequent commands are executed on `CITY`.

### **Example**

```
GET FILE=CITY /KEEP FOOD RENT PUBTRANS TEACHER COOK ENGINEER.
CORRELATIONS VARIABLES=FOOD TO ENGINEER
/MATRIX OUT(*) .
LIST.
DISPLAY DICTIONARY.
```

---

*CORRELATIONS*

- *CORRELATIONS* writes the same matrix as in the example above. However, the matrix data file replaces the working file. The `LIST` and `DISPLAY` commands are executed on the matrix file, not on the *CITY* file.

**Example**

```
CORRELATIONS VARIABLES=FOOD RENT COOK TEACHER MANAGER ENGINEER  
/FOOD TO TEACHER /PUBTRANS WITH MECHANIC  
/MATRIX OUT(*).
```

- Only the matrix for *FOOD TO TEACHER* is written to the matrix data file because it is the last variable list that does not use the keyword `WITH`.



# CORRESPONDENCE

CORRESPONDENCE is available in the Categories option.

```
CORRESPONDENCE

/TABLE = {rowvar (min, max) BY colvar (min, max)}
        {ALL (# of rows, # of columns)      }

[/SUPPLEMENTARY = [{rowvar (valuelist)}] [{colvar (valuelist)}]
                  {ROW (valuelist)      } {COLUMN (valuelist)}

[/EQUAL = [{rowvar (valuelist)}] [{colvar (valuelist)}]
          {ROW (valuelist)      } {COLUMN (valuelist)}

[/MEASURE = {CHISQ**}
            {EUCLID  }

[/STANDARDIZE = {RMEAN  }
                {CMEAN  }
                {RCMEAN**}
                {RSUM   }
                {CSUM   }

[/DIMENSION = {2**  }
              {value}

[/NORMALIZATION = {SYMMETRICAL**}
                  {PRINCIPAL  }
                  {RPRINCIPAL }
                  {CPRINCIPAL }
                  {value      }

[/PRINT = [TABLE**] [RPROF] [CPROF] [RPOINTS**] [CPOINTS**]
          [RCONF] [CCONF] [PERMUTATION[(n)]] [DEFAULT] [NONE]]

[/PLOT = [NDIM({value,value})]
        {value,MAX  }
        [RPOINTS[(n)]] [CPOINTS[(n)]] [TRROWS[(n)]]
        [TRCOLUMNS[(n)]] [BIPLOT**[(n)]] [NONE]]

[/OUTFILE = [SCORE('savfile'|'dataset')] [VARIANCE('savfile'|'dataset')]
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 13.0

- For the NDIM keyword on the PLOT subcommand, the default is changed to all dimensions.
- The maximum label length on the PLOT subcommand is increased to 60 (previous value was 20).

## Overview

CORRESPONDENCE displays the relationships between rows and columns of a two-way table graphically by a biplot. It computes the row and column scores and statistics and produces plots based on the scores. Also, confidence statistics are computed.

### Options

**Number of Dimensions.** You can specify how many dimensions CORRESPONDENCE should compute.

**Supplementary Points.** You can specify supplementary rows and columns.

**Equality Restrictions.** You can restrict rows and columns to have equal scores.

**Measure.** You can specify the distance measure to be the chi-square or Euclidean.

**Standardization.** You can specify one of five different standardization methods.

**Method of Normalization.** You can specify one of five different methods for normalizing the row and column scores.

**Confidence Statistics.** You can request computation of confidence statistics (standard deviations and correlations) for row and column scores. For singular values, confidence statistics are always computed.

**Data Input.** You can analyze individual casewise data, aggregated data, or table data.

**Display Output.** You can control which statistics are displayed and plotted.

**Writing Matrices.** You can write the row and column scores and the confidence statistics (variances and covariances) for the singular values to external files.

### Basic Specification

- The basic specification is CORRESPONDENCE and the TABLE subcommand. By default, CORRESPONDENCE computes a two-dimensional solution and displays the correspondence table, the summary table, an overview of the row and column scores, and a biplot of the row and column points.

### Subcommand Order

- The TABLE subcommand must appear first.
- All other subcommands can appear in any order.

### Syntax Rules

- Only one keyword can be specified on the MEASURE subcommand.
- Only one keyword can be specified on the STANDARDIZE subcommand.
- Only one keyword can be specified on the NORMALIZATION subcommand.
- Only one parameter can be specified on the DIMENSION subcommand.

**Operations**

- If a subcommand is specified more than once, only the last occurrence is executed.

**Limitations**

- The table input data and the aggregated input data cannot contain negative values. `CORRESPONDENCE` will treat such values as 0.
- Rows and columns that are specified as supplementary cannot be equalized.
- The maximum number of supplementary points for a variable is 200.
- The maximum number of equalities for a variable is 200.

**Example**

```
CORRESPONDENCE TABLE=MENTAL(1,4) BY SES(1,6)
/PRINT=RPOINTS CPOINTS
/PLOT=RPOINTS CPOINTS.
```

- Two variables, *MENTAL* and *SES*, are specified on the `TABLE` subcommand. *MENTAL* has values ranging from 1 to 4, and *SES* has values ranging from 1 to 6.
- The summary table and overview tables of the row and column scores are displayed.
- The row points plot and the column points plot are produced.

**TABLE Subcommand**

`TABLE` specifies the row and column variables along with their integer value ranges. The two variables are separated by the keyword `BY`.

- The `TABLE` subcommand is required.

**Casewise Data**

- Each variable is followed by an integer value range in parentheses. The value range consists of the variable's minimum value and its maximum value.
- Values outside of the specified range are not included in the analysis.
- Values do not have to be sequential. Empty categories yield a zero in the input table and do not affect the statistics for other categories.

**Example**

```
DATA LIST FREE/VAR1 VAR2.
BEGIN DATA
3 1
6 1
3 1
4 2
4 2
6 3
6 3
6 3
3 2
```

## CORRESPONDENCE

```

4 2
6 3
END DATA.
CORRESPONDENCE TABLE=VAR1 (3, 6) BY VAR2 (1, 3) .

```

- DATA LIST defines two variables, *VAR1* and *VAR2*.
- *VAR1* has three levels, coded 3, 4, and 6. *VAR2* also has three levels, coded 1, 2, and 3.
- Since a range of (3,6) is specified for *VAR1*, CORRESPONDENCE defines four categories, coded 3, 4, 5, and 6. The empty category, 5, for which there is no data, receives system-missing values for all statistics and does not affect the analysis.

## Aggregated Data

To analyze aggregated data, such as data from a crosstabulation where cell counts are available but the original raw data are not, you can use the WEIGHT command before CORRESPONDENCE.

### Example

To analyze a 3×3 table, such as the one shown below, you could use these commands:

```

DATA LIST FREE/ BIRTHORD ANXIETY COUNT.
BEGIN DATA
1 1 48
1 2 27
1 3 22
2 1 33
2 2 20
2 3 39
3 1 29
3 2 42
3 3 47
END DATA.
WEIGHT BY COUNT.
CORRESPONDENCE TABLE=BIRTHORD (1,3) BY ANXIETY (1,3) .

```

- The WEIGHT command weights each case by the value of *COUNT*, as if there are 48 subjects with *BIRTHORD*=1 and *ANXIETY*=1, 27 subjects with *BIRTHORD*=1 and *ANXIETY*=2, and so on.
- CORRESPONDENCE can then be used to analyze the data.
- If any of the table cell values (the values of the WEIGHT variable) equals 0, the WEIGHT command issues a warning, but the CORRESPONDENCE analysis is done correctly.
- The table cell values (the values of the WEIGHT variable) cannot be negative.

Table 37-1  
3 × 3 table

		Anxiety		
		High	Med	Low
Birth order	First	48	27	22
	Second	33	20	39
	Other	29	42	47

## Table Data

- The cells of a table can be read and analyzed directly by using the keyword `ALL` after `TABLE`.
- The columns of the input table must be specified as variables on the `DATA LIST` command. Only columns are defined, not rows.
- `ALL` is followed by the number of rows in the table, a comma, and the number of columns in the table, all in parentheses.
- The row variable is named `ROW`, and the column variable is named `COLUMN`.
- The number of rows and columns specified can be smaller than the actual number of rows and columns if you want to analyze only a subset of the table.
- The variables (columns of the table) are treated as the column categories, and the cases (rows of the table) are treated as the row categories.
- Row categories can be assigned values (category codes) when you specify `TABLE=ALL` by the optional variable `ROWCAT_`. This variable must be defined as a numeric variable with unique values corresponding to the row categories. If `ROWCAT_` is not present, the row index (case) numbers are used as row category values.

### Example

```
DATA LIST /ROWCAT_ 1 COL1 3-4 COL2 6-7 COL3 9-10.
BEGIN DATA
1 50 19 26
2 16 40 34
3 12 35 65
4 11 20 58
END DATA.
VALUE LABELS ROWCAT_ 1 'ROW1' 2 'ROW2' 3 'ROW3' 4 'ROW4'.
CORRESPONDENCE TABLE=ALL(4,3).
```

- `DATA LIST` defines the row category naming variable `ROWCAT_` and the three columns of the table as the variables.
- The `TABLE=ALL` specification indicates that the data are the cells of a table. The (4,3) specification indicates that there are four rows and three columns.
- The column variable is named `COLUMN` with categories labeled `COL1`, `COL2`, and `COL3`.
- The row variable is named `ROW` with categories labeled `ROW1`, `ROW2`, `ROW3`, and `ROW4`.

## DIMENSION Subcommand

`DIMENSION` specifies the number of dimensions you want `CORRESPONDENCE` to compute.

- If you do not specify the `DIMENSION` subcommand, `CORRESPONDENCE` computes two dimensions.
- `DIMENSION` is followed by a positive integer indicating the number of dimensions. If this parameter is omitted, a value of 2 is assumed.
- In general, you should choose as few dimensions as needed to explain most of the variation. The minimum number of dimensions that can be specified is 1. The maximum number of dimensions that can be specified equals the minimum of the number of active rows and the number of active columns minus 1. An active row or column is a nonsupplementary row or

column that is used in the analysis. For example, in a table where the number of rows is 5 (2 of which are supplementary) and the number of columns is 4, the number of active rows (3) is smaller than the number of active columns (4). Thus, the maximum number of dimensions that can be specified is  $(5-2)-1$ , or 2. Rows and columns that are restricted to have equal scores count as 1 toward the number of active rows or columns. For example, in a table with five rows and four columns, where two columns are restricted to have equal scores, the number of active rows is 5 and the number of active columns is  $(4-1)$ , or 3. The maximum number of dimensions that can be specified is  $(3-1)$ , or 2. Empty rows and columns (rows or columns with no data, all zeros, or all missing data) are not counted toward the number of rows and columns.

- If more than the maximum allowed number of dimensions is specified, CORRESPONDENCE reduces the number of dimensions to the maximum.

## **SUPPLEMENTARY Subcommand**

The SUPPLEMENTARY subcommand specifies the rows and/or columns that you want to treat as supplementary (also called passive or illustrative).

- For casewise data, the specification on SUPPLEMENTARY is the row and/or column variable name, followed by a value list in parentheses. The values must be in the value range specified on the TABLE subcommand for the row or column variable.
- For table data, the specification on SUPPLEMENTARY is *ROW* and/or *COLUMN*, followed by a value list in parentheses. The values represent the row or column indices of the table input data.
- The maximum number of supplementary rows or columns is the number of rows or columns minus 2. Rows and columns that are restricted to have equal scores count as 1 toward the number of rows or columns.
- Supplementary rows and columns cannot be equalized.

### **Example**

```
CORRESPONDENCE TABLE=MENTAL(1,8) BY SES(1,6)
/SUPPLEMENTARY MENTAL(3) SES(2,6) .
```

- SUPPLEMENTARY specifies the third level of *MENTAL* and the second and sixth levels of *SES* to be supplementary.

### **Example**

```
CORRESPONDENCE TABLE=ALL(8,6)
/SUPPLEMENTARY ROW(3) COLUMN(2,6) .
```

- SUPPLEMENTARY specifies the third level of the row variable and the second and sixth levels of the column variable to be supplementary.

## ***EQUAL Subcommand***

The `EQUAL` subcommand specifies the rows and/or columns that you want to restrict to have equal scores.

- For casewise data, the specification on `EQUAL` is the row and/or column variable name, followed by a list of at least two values in parentheses. The values must be in the value range specified on the `TABLE` subcommand for the row or column variable.
- For table data, the specification on `EQUAL` is `ROW` and/or `COLUMN`, followed by a value list in parentheses. The values represent the row or column indices of the table input data.
- Rows or columns that are restricted to have equal scores cannot be supplementary.
- The maximum number of equal rows or columns is the number of active rows or columns minus 1.

### ***Example***

```
CORRESPONDENCE TABLE=MENTAL(1,8) BY SES(1,6)
/EQUAL MENTAL(1,2) (6,7) SES(1,2,3).
```

- `EQUAL` specifies the first and second level of `MENTAL`, the sixth and seventh level of `MENTAL`, and the first, second, and third levels of `SES` to have equal scores.

## ***MEASURE Subcommand***

The `MEASURE` subcommand specifies the measure of distance between the row and column profiles.

- Only one keyword can be used.

The following keywords are available:

<b>CHISQ</b>	<i>Chi-square distance.</i> This is the weighted distance, where the weight is the mass of the rows or columns. This is the default specification for <code>MEASURE</code> and is the necessary specification for standard correspondence analysis.
<b>EUCLID</b>	<i>Euclidean distance.</i> The distance is the square root of the sum of squared differences between the values for two rows or columns.

## ***STANDARDIZE Subcommand***

When `MEASURE=EUCLID`, the `STANDARDIZE` subcommand specifies the method of standardization.

- Only one keyword can be used.
- If `MEASURE` is `CHISQ`, only `RCMEAN` standardization can be used, resulting in standard correspondence analysis.

The following keywords are available:

<b>RMEAN</b>	<i>The row means are removed.</i>
<b>CMEAN</b>	<i>The column means are removed.</i>
<b>RCMEAN</b>	<i>Both the row and column means are removed. This is the default specification.</i>
<b>RSUM</b>	<i>First the row totals are equalized and then the row means are removed.</i>
<b>CSUM</b>	<i>First the column totals are equalized and then the column means are removed.</i>

## ***NORMALIZATION Subcommand***

The **NORMALIZATION** subcommand specifies one of five methods for normalizing the row and column scores. Only the scores and confidence statistics are affected; contributions and profiles are not changed.

The following keywords are available:

<b>SYMMETRICAL</b>	<i>For each dimension, rows are the weighted average of columns divided by the matching singular value, and columns are the weighted average of rows divided by the matching singular value. This is the default if the <b>NORMALIZATION</b> subcommand is not specified. Use this normalization method if you are primarily interested in differences or similarities between rows and columns.</i>
<b>PRINCIPAL</b>	<i>Distances between row points and distances between column points are approximations of chi-square distances or of Euclidean distances (depending on <b>MEASURE</b>). The distances represent the distance between the row or column and its corresponding average row or column profile. Use this normalization method if you want to examine both differences between categories of the row variable and differences between categories of the column variable (but not differences between variables).</i>
<b>RPRINCIPAL</b>	<i>Distances between row points are approximations of chi-square distances or of Euclidean distances (depending on <b>MEASURE</b>). This method maximizes distances between row points, resulting in row points that are weighted averages of the column points. This is useful when you are primarily interested in differences or similarities between categories of the row variable.</i>
<b>CPRINCIPAL</b>	<i>Distances between column points are approximations of chi-square distances or of Euclidean distances (depending on <b>MEASURE</b>). This method maximizes distances between column points, resulting in column points that are weighted averages of the row points. This is useful when you are primarily interested in differences or similarities between categories of the column variable.</i>

The fifth method allows the user to specify any value in the range  $-1$  to  $+1$ , inclusive. A value of  $1$  is equal to the **RPRINCIPAL** method, a value of  $0$  is equal to the **SYMMETRICAL** method, and a value of  $-1$  is equal to the **CPRINCIPAL** method. By specifying a value between  $-1$  and  $1$ , the user can spread the inertia over both row and column scores to varying degrees. This method is useful for making tailor-made biplots.

## ***PRINT Subcommand***

Use **PRINT** to control which of several correspondence statistics are displayed. The summary table (singular values, inertia, proportion of inertia accounted for, cumulative proportion of inertia accounted for, and confidence statistics for the maximum number of dimensions) is always



produced. If PRINT is not specified, the input table, the summary table, the overview of row points table, and the overview of column points table are displayed.

The following keywords are available:

<b>TABLE</b>	<i>A crosstabulation of the input variables showing row and column marginals.</i>
<b>RPROFILES</b>	<i>The row profiles. PRINT=RPROFILES is analogous to the CELLS=ROW subcommand in CROSSTABS.</i>
<b>CPROFILES</b>	<i>The column profiles. PRINT=CPROFILES is analogous to the CELLS=COLUMN subcommand in CROSSTABS.</i>
<b>RPOINTS</b>	<i>Overview of row points (mass, scores, inertia, contribution of the points to the inertia of the dimension, and the contribution of the dimensions to the inertia of the points).</i>
<b>CPOINTS</b>	<i>Overview of column points (mass, scores, inertia, contribution of the points to the inertia of the dimension, and the contribution of the dimensions to the inertia of the points).</i>
<b>RCONF</b>	<i>Confidence statistics (standard deviations and correlations) for the active row points.</i>
<b>CCONF</b>	<i>Confidence statistics (standard deviations and correlations) for the active column points.</i>
<b>PERMUTATION(n)</b>	<i>The original table permuted according to the scores of the rows and columns. PERMUTATION can be followed by a number in parentheses indicating the maximum number of dimensions for which you want permuted tables. The default number of dimensions is 1.</i>
<b>NONE</b>	<i>No output other than the SUMMARY table.</i>
<b>DEFAULT</b>	<i>TABLE, RPOINTS, CPOINTS, and the SUMMARY tables. These statistics are displayed if you omit the PRINT subcommand.</i>

## ***PLOT Subcommand***

Use PLOT to produce a biplot of row and column points, plus plots of the row points, column points, transformations of the categories of the row variable, and transformations of the categories of the column variable. If PLOT is not specified or is specified without keywords, a biplot is produced.

The following keywords are available:

<b>TRROWS(n)</b>	<i>Transformation plots for the rows (row category scores against row category indicator values).</i>
<b>TRCOLUMNS(n)</b>	<i>Transformation plots for the columns (column category scores against column category indicator values).</i>
<b>RPOINTS(n)</b>	<i>Plot of the row points.</i>
<b>CPOINTS(n)</b>	<i>Plot of the column points.</i>

**BIPLOT(n)** *Biplot of the row and column points.* This is the default plot. This plot is not available when NORMALIZATION=PRINCIPAL.

**NONE** *No plots.*

- For all of the keywords except NONE the user can specify an optional parameter *l* in parentheses in order to control the global upper boundary of value label lengths in the plot. The label length parameter *l* can take any nonnegative integer less than or equal to the applicable maximum length of 60. If *l* is not specified, CORRESPONDENCE assumes that each value label at its full length is displayed. If *l* is an integer larger than the applicable maximum, then we reset it to the applicable maximum, but do not issue a warning. If a positive value of *l* is given but if some or all of the category values do not have labels, then for those values the values themselves are used as the labels.

In addition to the plot keywords, the following can be specified:

**NDIM(value,value)** *Dimension pairs to be plotted.* NDIM is followed by a pair of values in parentheses. If NDIM is not specified or if NDIM is specified without parameter values, a matrix scatterplot including all dimensions is produced.

- The first value must be any integer from 1 to the number of dimensions in the solution minus 1.
- The second value must be an integer from 2 to the number of dimensions in the solution. The second value must exceed the first. Alternatively, the keyword MAX can be used instead of a value to indicate the highest dimension of the solution.
- For TRROWS and TRCOLUMNS, the first and second values indicate the range of dimensions for which the plots are created.
- For RPOINTS, CPOINTS, and BILOT, the first and second values indicate plotting pairs of dimensions. The first value indicates the dimension that is plotted against higher dimensions. The second value indicates the highest dimension to be used in plotting the dimension pairs.

### Example

```
CORRESPONDENCE TABLE=MENTAL(1,4) BY SES(1,6)
/PLOT NDIM(1,3) BILOT(5) .
```

- BILOT and NDIM(1,3) requests that a scatterplot for dimensions 1 and 2, and a scatterplot for dimensions 1 and 3 should be produced.
- The 5 following BILOT indicates that only the first five characters of each label are to be shown in the biplot matrix.

### Example

```
CORRESPONDENCE TABLE=MENTAL(1,4) BY SES(1,6)
/DIMENSION = 3
/PLOT NDIM(1,MAX) TRROWS .
```

- Three transformation plots for the row categories are produced, one for each dimension from 1 to the highest dimension of the analysis (in this case, 3). The label parameter is not specified, and so the category labels in the plot are shown up their full lengths.

## OUTFILE Subcommand

Use `OUTFILE` to write row and column scores and/or confidence statistics (variances and covariances) for the singular values and row and column scores to an external IBM® SPSS® Statistics data file or previously declared dataset.

`OUTFILE` must be followed by one or both of the following keywords:

**SCORE** ('file'|'dataset')      *Write row and column scores.*  
**VARIANCE** ('file'|'dataset')      *Write variances and covariances.*

- Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. Datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files. The names should be different for the each of the keywords.
- For `VARIANCE`, supplementary and equality constrained rows and columns are not produced in the external file.

The variables in the `SCORE` matrix data file and their values are:

**ROWTYPE\_**      *String variable containing the value ROW for all of the rows and COLUMN for all of the columns.*  
**LEVEL\_**      *String variable containing the values (or value labels, if present) of each original variable.*  
**VARNAME\_**      *String variable containing the original variable names.*  
**DIM1...DIMn**      *Numerical variables containing the row and column scores for each dimension. Each variable is named DIMn, where n represents the dimension number.*

The variables in the `VARIANCE` matrix data file and their values are:

**ROWTYPE\_**      *String variable containing the value COV for all of the cases in the file.*  
**VARNAME\_**      *String variable containing the value SINGULAR, the row variable's name, and the column variable's name.*  
**LEVEL\_**      *String variable containing the row variable's values (or labels), the column variable's values (or labels), and a blank value for VARNAME\_ = SINGULAR.*  
**DIMNMBR\_**      *String variable containing the dimension number.*  
**DIM1...DIMn**      *Numerical variables containing the variances and covariances for each dimension. Each variable is named DIMn, where n represents the dimension number.*

# COUNT

```
COUNT varname=varlist(value list) [/varname=...]
```

*Keywords for numeric value lists:*

LOWEST, LO, HIGHEST, HI, THRU, MISSING, SYSMIS

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
COUNT TARGET=V1 V2 V3 (2).
```

## **Overview**

COUNT creates a numeric variable that, for each case, counts the occurrences of the same value (or list of values) across a list of variables. The new variable is called the *target* variable. The variables and values that are counted are the *criterion* variables and values. Criterion variables can be either numeric or string.

### **Basic Specification**

The basic specification is the target variable, an equals sign, the criterion variable(s), and the criterion value(s) enclosed in parentheses.

### **Syntax Rules**

- Use a slash to separate the specifications for each target variable.
- The criterion variables specified for a single target variable must be either all numeric or all string.
- Each value on a list of criterion values must be separated by a comma or space. String values must be enclosed in quotes.
- The keywords THRU, LOWEST (LO), HIGHEST (HI), SYSMIS, and MISSING can be used only with numeric criterion variables.
- A variable can be specified on more than one criterion variable list.
- You can use the keyword TO to specify consecutive criterion variables that have the same criterion value or values.
- You can specify multiple variable lists for a single target variable to count different values for different variables.

**Operations**

- Target variables are always numeric and are initialized to 0 for each case. They are assigned a dictionary format of F8.2.
- If the target variable already exists, its previous values are replaced.
- COUNT ignores the missing-value status of user-missing values. It counts a value even if that value has been previously declared as missing.
- The target variable is never system-missing. To define user-missing values for target variables, use the RECODE or MISSING VALUES command.
- SYSMIS counts system-missing values for numeric variables.
- MISSING counts both user- and system-missing values for numeric variables.

**Examples****Counting Occurrences of a Single Value**

```
COUNT TARGET=V1 V2 V3 (2).
```

- The value of *TARGET* for each case will be either 0, 1, 2, or 3, depending on the number of times the value 2 occurs across the three variables for each case.
- *TARGET* is a numeric variable with an F8.2 format.

**Counting Occurrences of a Range of Values and System-Missing Values**

```
COUNT QLOW=Q1 TO Q10 (LO THRU 0)
/QSYSMIS=Q1 TO Q10 (SYSMIS).
```

- Assuming that there are 10 variables between and including *Q1* and *Q10* in the active dataset, *QLOW* ranges from 0 to 10, depending on the number of times a case has a negative or 0 value across the variables *Q1* to *Q10*.
- *QSYSMIS* ranges from 0 to 10, depending on how many system-missing values are encountered for *Q1* to *Q10* for each case. User-missing values are not counted.
- Both *QLOW* and *QSYSMIS* are numeric variables and have F8.2 formats.

**Counting Occurrences of String Values**

```
COUNT SVAR=V1 V2 ('male ') V3 V4 V5 ('female').
```

- *SVAR* ranges from 0 to 5, depending on the number of times a case has a value of male for *V1* and *V2* and a value of female for *V3*, *V4*, and *V5*.
- *SVAR* is a numeric variable with an F8.2 format.

# COXREG

COXREG is available in the Advanced Statistics option.

```
COXREG VARIABLES = survival varname [WITH varlist]
/ STATUS = varname [EVENT] (vallist) [LOST (vallist)]
[/STRATA = varname]
[/CATEGORICAL = varname]
[/CONTRAST (varname) = {DEVIATION (refcat)}]
                        {SIMPLE (refcat) }
                        {DIFFERENCE }
                        {HELMERT }
                        {REPEATED }
                        {POLYNOMIAL(metric)}
                        {SPECIAL (matrix) }
                        {INDICATOR (refcat)}

[/METHOD = {ENTER**} ] [{varlist}]
      {BSTEP [{COND}]} {ALL }
      {LR }
      {WALD}
      {FSTEP [{COND}]}
      {LR }
      {WALD}

[/MISSING = {EXCLUDE**}]
      {INCLUDE }

[/PRINT = [{DEFAULT**}] [CI ({95})]]
      {SUMMARY } {n }
      {BASELINE }
      {CORR }
      {ALL }

[/CRITERIA = [{BCON}({1E-4**})] [LCON({1E-5**})]
      {PCON} { n } { n }
      [ITERATE({20**})]
      { n }
      [PIN({0.05**})] [POUT({0.1**})]
      { n } { n }

[/PLOT = [NONE**] [SURVIVAL] [HAZARD] [LML] [OMS]]
[/PATTERN = [varname(value)...] [BY varname]]
[/OUTFILE = [COEFF('savfile' | 'dataset')] [TABLE('savfile' | 'dataset')]
      [PARAMETER('file')]]
[/SAVE = tempvar [(newvarname)],tempvar ...]
[/EXTERNAL]
```

**\*\*Default if subcommand or keyword is omitted.**

*Temporary variables created by COXREG are:*

*SURVIVAL*

*SE*

*HAZARD*

*RESID*

*LML*

*DFBETA*

*PRESID*

## *XBETA*

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Example**

```
TIME PROGRAM.  
COMPUTE Z=AGE + T_ .  
  
COXREG SURVIVAL WITH Z  
  /STATUS SURVSTA EVENT(1) .
```

## **Overview**

COXREG applies Cox proportional hazards regression to analysis of survival times—that is, the length of time before the occurrence of an event. COXREG supports continuous and categorical independent variables (covariates), which can be time dependent. Unlike SURVIVAL and KM, which compare only distinct subgroups of cases, COXREG provides an easy way of considering differences in subgroups as well as analyzing effects of a set of covariates.

### **Options**

**Processing of Independent Variables.** You can specify which of the independent variables are categorical with the CATEGORICAL subcommand and control treatment of these variables with the CONTRAST subcommand. You can select one of seven methods for entering independent variables into the model using the METHOD subcommand. You can also indicate interaction terms using the keyword BY between variable names on either the VARIABLES subcommand or the METHOD subcommand.

**Specifying Termination and Model-Building Criteria.** You can specify the criteria for termination of iteration and control variable entry and removal with the CRITERIA subcommand.

**Adding New Variables to Active Dataset.** You can use the SAVE subcommand to save the cumulative survival, standard error, cumulative hazard, log-minus-log-of-survival function, residuals, XBeta, and, wherever available, partial residuals and DfBeta.

**Output.** You can print optional output using the PRINT subcommand, suppress or request plots with the PLOT subcommand, and, with the OUTFILE subcommand, write data files containing coefficients from the final model or a survival table. When only time-constant covariates are used, you can use the PATTERN subcommand to specify a pattern of covariate values in addition to the covariate means to use for the plots and the survival table.

### **Basic Specification**

- The minimum specification on COXREG is a dependent variable with the STATUS subcommand.
- To analyze the influence of time-constant covariates on the survival times, the minimum specification requires either the WITH keyword followed by at least one covariate (independent variable) on the VARIABLES subcommand or a METHOD subcommand with at least one independent variable.

- To analyze the influence of time-dependent covariates on the survival times, the `TIME PROGRAM` command and transformation language are required to define the functions for the time-dependent covariate(s).

### **Subcommand Order**

- The `VARIABLES` subcommand must be specified first; the subcommand keyword is optional.
- Remaining subcommands can be named in any order.

### **Syntax Rules**

- Only one dependent variable can be specified for each `COXREG` command.
- Any number of covariates (independent variables) can be specified. The dependent variable cannot appear on the covariate list.
- The covariate list is required if any of the `METHOD` subcommands are used without a variable list or if the `METHOD` subcommand is not used.
- Only one status variable can be specified on the `STATUS` subcommand. If multiple `STATUS` subcommands are specified, only the last specification is in effect.
- You can use the `BY` keyword to specify interaction between covariates.

### **Operations**

- `TIME PROGRAM` computes the values for time-dependent covariates. [For more information, see the topic TIME PROGRAM on p. 1962.](#)
- `COXREG` replaces covariates specified on `CATEGORICAL` with sets of contrast variables. In stepwise analyses, the set of contrast variables associated with one categorical variable is entered or removed from the model as a block.
- Covariates are screened to detect and eliminate redundancies.
- `COXREG` deletes all cases that have negative values for the dependent variable.

### **Limitations**

- Only one dependent variable is allowed.
- Maximum 100 covariates in a single interaction term.
- Maximum 35 levels for a `BY` variable on `PATTERN`.

## **Examples**

```
COXREG
  tenure /STATUS=churn(1)
  /PATTERN BY custcat
  /CONTRAST (marital)=Indicator /CONTRAST (ed)=Indicator /CONTRAST
  (retire)=Indicator /CONTRAST (gender)=Indicator /CONTRAST
  (custcat)=Indicator
  /METHOD=FSTEP(LR) age marital address ed employ retire gender reside
  /METHOD=ENTER custcat
  /PLOT SURVIVAL HAZARD
  /CRITERIA=PIN(.05) POUT(.10) ITERATE(20) .
```

- The procedure fits a Cox regression model to the variable *tenure*.



- The `STATUS` subcommand specifies that a value of 1 on the variable *churn* indicates the event of interest has occurred.
- The `PATTERN` subcommand specifies that separate lines be produced for each value of *custcat* on the requested plots.
- The `CONTRAST` subcommand specifies that *marital*, *ed*, *retire*, *gender*, and *custcat* should be treated as categorical variables using indicator contrasts.
- The first `METHOD` subcommand specifies that *age*, *marital*, *address*, *ed*, *employ*, *retire*, *gender*, and *reside* should be tested for entry into the model using forward stepwise selection using the likelihood ratio statistic.
- The second `METHOD` subcommand specifies that *custcat* should be entered into the model after the forward stepwise selection is performed in the previous `METHOD` subcommand.
- The `PLOT` subcommand requests plots of the cumulative survival and cumulative hazard functions.
- All other options are set to their default values.

### **Using a Time-Dependent Covariate**

```
TIME PROGRAM.
COMPUTE T_COV_ = T_*age .
COXREG
  time /STATUS=arrest2(1)
  /METHOD=ENTER age T_COV_
  /CRITERIA=PIN(.05) POUT(.10) ITERATE(20) .
```

- `TIME PROGRAM` defines the time-dependent covariate *T\_COV\_* as the interaction between the current time and *age*.
- `COXREG` fits a Cox regression model to the variable *time*.
- The `STATUS` subcommand specifies that a value of 1 on the variable *arrest2* indicates the event of interest (a second arrest) has occurred.
- The `METHOD` subcommand specifies that *age* and *T\_COV\_* should be entered into the model.
- All other options are set to their default values.

For more information, see the topic [TIME PROGRAM](#) on p. 1962.

## **VARIABLES Subcommand**

`VARIABLES` identifies the dependent variable and the covariates to be included in the analysis.

- The minimum specification is the dependent variable.
- Cases whose dependent variable values are negative are excluded from the analysis.
- You must specify the keyword `WITH` and a list of all covariates if no `METHOD` subcommand is specified or if a `METHOD` subcommand is specified without naming the variables to be used.
- If the covariate list is not specified on `VARIABLES` but one or more `METHOD` subcommands are used, the covariate list is assumed to be the union of the sets of variables listed on all of the `METHOD` subcommands.

- You can specify an interaction of two or more covariates using the keyword `BY`. For example, `A B BY C D` specifies the three terms  $A$ ,  $B*C$ , and  $D$ .
- The keyword `TO` can be used to specify a list of covariates. The implied variable order is the same as in the active dataset.

## **STATUS Subcommand**

To determine whether the event has occurred for a particular observation, `COXREG` checks the value of a status variable. `STATUS` lists the status variable and the code for the occurrence of the event.

- Only one status variable can be specified. If multiple `STATUS` subcommands are specified, `COXREG` uses the last specification and displays a warning.
- The keyword `EVENT` is optional, but the value list in parentheses must be specified.
- The value list must be enclosed in parentheses. All cases with non-negative times that do not have a code within the range specified after `EVENT` are classified as **censored cases**—that is, cases for which the event has not yet occurred.
- The value list can be one value, a list of values separated by blanks or commas, a range of values using the keyword `THRU`, or a combination.
- If missing values occur within the specified ranges, they are ignored if `MISSING=EXCLUDE` (the default) is specified, but they are treated as valid values for the range if `MISSING=INCLUDE` is specified.
- The status variable can be either numeric or string. If a string variable is specified, the `EVENT` values must be enclosed in apostrophes and the keyword `THRU` cannot be used.

### **Example**

```
COXREG VARIABLES = SURVIVAL WITH GROUP
  /STATUS SURVSTA (3 THRU 5, 8 THRU 10).
```

- `STATUS` specifies that `SURVSTA` is the status variable.
- A value between either 3 and 5 or 8 and 10, inclusive, means that the terminal event occurred.
- Values outside the specified ranges indicate censored cases.

## **STRATA Subcommand**

`STRATA` identifies a stratification variable. A different baseline survival function is computed for each stratum.

- The only specification is the subcommand keyword with one, and only one, variable name.
- If you have more than one stratification variable, create a new variable that corresponds to the combination of categories of the individual variables before invoking the `COXREG` command.
- There is no limit to the number of levels for the strata variable.

### **Example**

```
COXREG VARIABLES = SURVIVAL WITH GROUP
  /STATUS SURVSTA (1)
```

---

`/STRATA=LOCATION.`

- `STRATA` specifies *LOCATION* as the strata variable.
- Different baseline survival functions are computed for each value of *LOCATION*.

## ***CATEGORICAL Subcommand***

`CATEGORICAL` identifies covariates that are nominal or ordinal. Variables that are declared to be categorical are automatically transformed to a set of contrast variables (see [CONTRAST Subcommand on p. 373](#)). If a variable coded as 0–1 is declared as categorical, by default, its coding scheme will be changed to deviation contrasts.

- Covariates not specified on `CATEGORICAL` are assumed to be at least interval, except for strings.
- Variables specified on `CATEGORICAL` but not on `VARIABLES` or any `METHOD` subcommand are ignored.
- Variables specified on `CATEGORICAL` are replaced by sets of contrast variables. If the categorical variable has  $n$  distinct values,  $n-1$  contrast variables will be generated. The set of contrast variables associated with one categorical variable are entered or removed from the model together.
- If any one of the variables in an interaction term is specified on `CATEGORICAL`, the interaction term is replaced by contrast variables.
- All string variables are categorical. Only the first eight bytes of each value of a string variable are used in distinguishing among values. Thus, if two values of a string variable are identical for the first eight characters, the values are treated as though they were the same.

## ***CONTRAST Subcommand***

`CONTRAST` specifies the type of contrast used for categorical covariates. The interpretation of the regression coefficients for categorical covariates depends on the contrasts used. The default is `DEVIATION`. For illustration of contrast types, see the appendix.

- The categorical covariate is specified in parentheses following `CONTRAST`.
- If the categorical variable has  $n$  values, there will be  $n-1$  rows in the contrast matrix. Each contrast matrix is treated as a set of independent variables in the analysis.
- Only one variable can be specified per `CONTRAST` subcommand, but multiple `CONTRAST` subcommands can be specified.
- You can specify one of the contrast keywords in parentheses following the variable specification to request a specific contrast type.

The following contrast types are available:

<b>DEVIATION(refcat)</b>	<i>Deviations from the overall effect.</i> This is the default. The effect for each category of the independent variable except one is compared to the overall effect. <i>Refcat</i> is the category for which parameter estimates are not displayed (they must be calculated from the others). By default, <i>refcat</i> is the last category. To omit a category other than the last, specify the sequence number of the omitted category (which is not necessarily the same as its value) in parentheses following the keyword <code>DEVIATION</code> .
<b>SIMPLE(refcat)</b>	<i>Each category of the independent variable except the last is compared to the last category.</i> To use a category other than the last as the omitted reference category, specify its sequence number (which is not necessarily the same as its value) in parentheses following the keyword <code>SIMPLE</code> .
<b>DIFFERENCE</b>	<i>Difference or reverse Helmert contrasts.</i> The effects for each category of the covariate except the first are compared to the mean effect of the previous categories.
<b>HELMERT</b>	<i>Helmert contrasts.</i> The effects for each category of the independent variable except the last are compared to the mean effects of subsequent categories.
<b>POLYNOMIAL(metric)</b>	<i>Polynomial contrasts.</i> The first degree of freedom contains the linear effect across the categories of the independent variable, the second contains the quadratic effect, and so on. By default, the categories are assumed to be equally spaced; unequal spacing can be specified by entering a metric consisting of one integer for each category of the independent variable in parentheses after the keyword <code>POLYNOMIAL</code> . For example, <code>CONTRAST (STIMULUS) = POLYNOMIAL(1,2,4)</code> indicates that the three levels of <i>STIMULUS</i> are actually in the proportion 1:2:4. The default metric is always (1,2,... <i>k</i> ), where <i>k</i> categories are involved. Only the relative differences between the terms of the metric matter: (1,2,4) is the same metric as (2,3,5) or (20,30,50) because, in each instance, the difference between the second and third numbers is twice the difference between the first and second.
<b>REPEATED</b>	<i>Comparison of adjacent categories.</i> Each category of the independent variable except the last is compared to the next category.
<b>SPECIAL(matrix)</b>	<i>A user-defined contrast.</i> After this keyword, a matrix is entered in parentheses with <i>k</i> -1 rows and <i>k</i> columns, where <i>k</i> is the number of categories of the independent variable. The rows of the contrast matrix contain the special contrasts indicating the desired comparisons between categories. If the special contrasts are linear combinations of each other, <code>COXREG</code> reports the linear dependency and stops processing. If <i>k</i> rows are entered, the first row is discarded and only the last <i>k</i> -1 rows are used as the contrast matrix in the analysis.
<b>INDICATOR(refcat)</b>	<i>Indicator variables.</i> Contrasts indicate the presence or absence of category membership. By default, <i>refcat</i> is the last category (represented in the contrast matrix as a row of zeros). To omit a category other than the last, specify the sequence number of the category (which is not necessarily the same as its value) in parentheses after the keyword <code>INDICATOR</code> .

### Example

```
COXREG VARIABLES = SURVIVAL WITH GROUP
/STATUS SURVSTA (1)
/STRATA=LOCATION
/CATEGORICAL = GROUP
/CONTRAST (GROUP) = SPECIAL (2 -1 -1
                                0 1 -1) .
```

- The specification of *GROUP* on *CATEGORICAL* replaces the variable with a set of contrast variables.
- *GROUP* identifies whether a case is in one of the three treatment groups.
- A *SPECIAL* type contrast is requested. A three-column, two-row contrast matrix is entered in parentheses.

## ***METHOD Subcommand***

*METHOD* specifies the order of processing and the manner in which the covariates enter the model. If no *METHOD* subcommand is specified, the default method is *ENTER*.

- The subcommand keyword *METHOD* can be omitted.
- You can list all covariates to be used for the method on a variable list. If no variable list is specified, the default is *ALL*; all covariates named after *WITH* on the *VARIABLES* subcommand are used for the method.
- The keyword *BY* can be used between two variable names to specify an interaction term.
- Variables specified on *CATEGORICAL* are replaced by sets of contrast variables. The contrast variables associated with a categorical variable are entered or removed from the model together.
- Three keywords are available to specify how the model is to be built:

**ENTER**      *Forced entry.* All variables are entered in a single step. This is the default if the *METHOD* subcommand is omitted.

**FSTEP**      *Forward stepwise.* The covariates specified on *FSTEP* are tested for entry into the model one by one based on the significance level of the score statistic. The variable with the smallest significance less than *PIN* is entered into the model. After each entry, variables that are already in the model are tested for possible removal based on the significance of the Wald statistic, likelihood ratio, or conditional criterion. The variable with the largest probability greater than the specified *POUT* value is removed and the model is reestimated. Variables in the model are then again evaluated for removal. Once no more variables satisfy the removal criteria, covariates not in the model are evaluated for entry. Model building stops when no more variables meet entry or removal criteria, or when the current model is the same as a previous one.

**BSTEP**      *Backward stepwise.* As a first step, the covariates specified on *BSTEP* are entered into the model together and are tested for removal one by one. Stepwise removal and entry then follow the same process as described for *FSTEP* until no more variables meet entry and removal criteria, or when the current model is the same as a previous one.

- Multiple *METHOD* subcommands are allowed and are processed in the order in which they are specified. Each method starts with the results from the previous method. If *BSTEP* is used, all eligible variables are entered at the first step. All variables are then eligible for entry and removal unless they have been excluded from the *METHOD* variable list.

- The statistic used in the test for removal can be specified by an additional keyword in parentheses following `FSTEP` or `BSTEP`. If `FSTEP` or `BSTEP` is specified by itself, the default is `COND`.

<b>COND</b>	<i>Conditional statistic.</i> This is the default if <code>FSTEP</code> or <code>BSTEP</code> is specified by itself
<b>WALD</b>	<i>Wald statistic.</i> The removal of a covariate from the model is based on the significance of the Wald statistic.
<b>LR</b>	<i>Likelihood ratio.</i> The removal of a covariate from the model is based on the significance of the change in the log-likelihood. If <code>LR</code> is specified, the model must be reestimated without each of the variables in the model. This can substantially increase computational time. However, the likelihood-ratio statistic is better than the Wald statistic for deciding which variables are to be removed.

### Example

```
COXREG VARIABLES = SURVIVAL WITH GROUP SMOKE DRINK
  /STATUS SURVSTA (1)
  /CATEGORICAL = GROUP SMOKE DRINK
  /METHOD ENTER GROUP
  /METHOD BSTEP (LR) SMOKE DRINK SMOKE BY DRINK.
```

- `GROUP`, `SMOKE`, and `DRINK` are specified as covariates and as categorical variables.
- The first `METHOD` subcommand enters `GROUP` into the model.
- Variables in the model at the termination of the first `METHOD` subcommand are included in the model at the beginning of the second `METHOD` subcommand.
- The second `METHOD` subcommand adds `SMOKE`, `DRINK`, and the interaction of `SMOKE` with `DRINK` to the previous model.
- Backward stepwise regression analysis is then done using the likelihood-ratio statistic as the removal criterion. The variable `GROUP` is not eligible for removal because it was not specified on the `BSTEP` subcommand.
- The procedure continues until the removal of a variable will result in a decrease in the log-likelihood with a probability smaller than `POUT`.

## MISSING Subcommand

`MISSING` controls missing value treatments. If `MISSING` is omitted, the default is `EXCLUDE`.

- Cases with negative values on the dependent variable are automatically treated as missing and are excluded.
- To be included in the model, a case must have nonmissing values for the dependent, status, strata, and all independent variables specified on the `COXREG` command.

<b>EXCLUDE</b>	<i>Exclude user-missing values.</i> User-missing values are treated as missing. This is the default if <code>MISSING</code> is omitted.
<b>INCLUDE</b>	<i>Include user-missing values.</i> User-missing values are included in the analysis.

## **PRINT Subcommand**

By default, COXREG prints a full regression report for each step. You can use the PRINT subcommand to request specific output. If PRINT is not specified, the default is DEFAULT.

<b>DEFAULT</b>	<i>Full regression output including overall model statistics and statistics for variables in the equation and variables not in the equation. This is the default when PRINT is omitted.</i>
<b>SUMMARY</b>	<i>Summary information. The output includes <math>-2</math> log-likelihood for the initial model, one line of summary for each step, and the final model printed with full detail.</i>
<b>CORR</b>	<i>Correlation/covariance matrix of parameter estimates for the variables in the model.</i>
<b>BASELINE</b>	<i>Baseline table. For each stratum, a table is displayed showing the baseline cumulative hazard, as well as survival, standard error, and cumulative hazard evaluated at the covariate means for each observed time point in that stratum.</i>
<b>CI (value)</b>	<i>Confidence intervals for <math>e^\beta</math>. Specify the confidence level in parentheses. The requested intervals are displayed whenever a variables-in-equation table is printed. The default is 95%.</i>
<b>ALL</b>	<i>All available output.</i>

Estimation histories showing the last 10 iterations are printed if the solution fails to converge.

### **Example**

```
COXREG VARIABLES = SURVIVAL WITH GROUP
/STATUS = SURVSTA (1)
/STRATA = LOCATION
/CATEGORICAL = GROUP
/METHOD = ENTER
/PRINT ALL.
```

PRINT requests summary information, a correlation matrix for parameter estimates, a baseline survival table for each stratum, and confidence intervals for  $e^\beta$  with each variables-in-equation table, in addition to the default output.

## **CRITERIA Subcommand**

CRITERIA controls the statistical criteria used in building the Cox Regression models. The way in which these criteria are used depends on the method specified on the METHOD subcommand. The default criteria are noted in the description of each keyword below. Iterations will stop if any of the criteria for BCON, LCON, or ITERATE are satisfied.

<b>BCON(value)</b>	<i>Change in parameter estimates for terminating iteration. Alias PCON. Iteration terminates when the parameters change by less than the specified value. BCON defaults to <math>1E-4</math>. To eliminate this criterion, specify a value of 0.</i>
<b>ITERATE(value)</b>	<i>Maximum number of iterations. If a solution fails to converge after the maximum number of iterations has been reached, COXREG displays an iteration history showing the last 10 iterations and terminates the procedure. The default for ITERATE is 20.</i>

<b>LCON(value)</b>	<i>Percentage change in the log-likelihood ratio for terminating iteration. If the log-likelihood decreases by less than the specified value, iteration terminates. LCON defaults to 1E-5. To eliminate this criterion, specify a value of 0.</i>
<b>PIN(value)</b>	<i>Probability of score statistic for variable entry. A variable whose significance level is greater than PIN cannot enter the model. The default for PIN is 0.05.</i>
<b>POUT(value)</b>	<i>Probability of Wald, LR, or conditional LR statistic to remove a variable. A variable whose significance is less than POUT cannot be removed. The default for POUT is 0.1.</i>

**Example**

```
COXREG VARIABLES = SURVIVAL WITH GROUP AGE BP TMRSZ
  /STATUS = SURVSTA (1)
  /STRATA = LOCATION
  /CATEGORICAL = GROUP
  /METHOD BSTEP
  /CRITERIA BCON(0) ITERATE(10) PIN(0.01) POUT(0.05) .
```

- A backward stepwise Cox Regression analysis is performed.
- CRITERIA alters four of the default statistical criteria that control the building of a model.
- Zero specified on BCON indicates that change in parameter estimates is not a criterion for termination. BCON can be set to 0 if only LCON and ITER are to be used.
- ITERATE specifies that the maximum number of iterations is 10. LCON is not changed and the default remains in effect. If either ITERATE or LCON is met, iterations will terminate.
- POUT requires that the probability of the statistic used to test whether a variable should remain in the model be smaller than 0.05. This is more stringent than the default value of 0.1.
- PIN requires that the probability of the score statistic used to test whether a variable should be included be smaller than 0.01. This makes it more difficult for variables to be included in the model than does the default PIN, which has a value of 0.05.

**PLOT Subcommand**

You can request specific plots to be produced with the PLOT subcommand. Each requested plot is produced once for each pattern specified on the PATTERN subcommand. If PLOT is not specified, the default is NONE (no plots are printed). Requested plots are displayed at the end of the final model.

- The set of plots requested is displayed for the functions at the mean of the covariates and at each combination of covariate values specified on PATTERN.
- If time-dependent covariates are included in the model, no plots are produced.
- Lines on a plot are connected as step functions.

<b>NONE</b>	<i>Do not display plots.</i>
<b>SURVIVAL</b>	<i>Plot the cumulative survival distribution.</i>
<b>HAZARD</b>	<i>Plot the cumulative hazard function.</i>
<b>LML</b>	<i>Plot the log-minus-log-of-survival function.</i>
<b>OMS</b>	<i>Plot the one-minus-survival function.</i>



## ***PATTERN Subcommand***

PATTERN specifies the pattern of covariate values to be used for the requested plots and coefficient tables.

- A value must be specified for each variable specified on PATTERN.
- Continuous variables that are included in the model but not named on PATTERN are evaluated at their means.
- Categorical variables that are included in the model but not named on PATTERN are evaluated at the means of the set of contrasts generated to replace them.
- You can request separate lines for each category of a variable that is in the model. Specify the name of the categorical variable after the keyword BY. The BY variable must be a categorical covariate. You cannot specify a value for the BY covariate.
- Multiple PATTERN subcommands can be specified. COXREG produces a set of requested plots for each specified pattern.
- PATTERN cannot be used when time-dependent covariates are included in the model.

## ***OUTFILE Subcommand***

OUTFILE writes data to an external IBM® SPSS® Statistics data file or a previously declared dataset (DATASET DECLARE command). COXREG writes two types of data files. You can specify the file type to be created with one of the two keywords, followed by a quoted file specification in parentheses. It also saves model information in XML format.

### **COEFF('savfile' | 'dataset')**

*Write a data file containing the coefficients from the final model.*

### **TABLE('savfile' | 'dataset')**

*Write the survival table to a data file. The file contains cumulative survival, standard error, and cumulative hazard statistics for each uncensored time within each stratum evaluated at the baseline and at the mean of the covariates. Additional covariate patterns can be requested on PATTERN.*

### **PARAMETER('file')**

*Write parameter estimates only to an XML file. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#)*

## ***SAVE Subcommand***

SAVE saves the temporary variables created by COXREG. The temporary variables include:

<b>SURVIVAL</b>	<i>Survival function evaluated at the current case.</i>
<b>SE</b>	<i>Standard error of the survival function.</i>
<b>HAZARD</b>	<i>Cumulative hazard function evaluated at the current case. Alias RESID.</i>
<b>LML</b>	<i>Log-minus-log-of-survival function.</i>
<b>DFBETA</b>	<i>Change in the coefficient if the current case is removed. There is one DFBETA for each covariate in the final model. If there are time-dependent covariates, only DFBETA can be requested. Requests for any other temporary variable are ignored.</i>

<b>PRESID</b>	<i>Partial residuals.</i> There is one residual variable for each covariate in the final model. If a covariate is not in the final model, the corresponding new variable has the system-missing value.
<b>XBETA</b>	<i>Linear combination of mean corrected covariates times regression coefficients from the final model.</i>

- To specify variable names for the new variables, assign the new names in parentheses following each temporary variable name.
- Assigned variable names must be unique in the active dataset. Scratch or system variable names cannot be used (that is, the variable names cannot begin with # or \$).
- If new variable names are not specified, COXREG generates default names. The default name is composed of the first three characters of the name of the temporary variable (two for *SE*), followed by an underscore and a number to make it unique.
- A temporary variable can be saved only once on the same *SAVE* subcommand.

**Example**

```
COXREG VARIABLES = SURVIVAL WITH GROUP
  /STATUS = SURVSTA (1)
  /STRATA = LOCATION
  /CATEGORICAL = GROUP
  /METHOD = ENTER
  /SAVE SURVIVAL HAZARD.
```

COXREG saves cumulative survival and hazard in two new variables, *SUR\_1* and *HAZ\_1*, provided that neither of the two names exists in the active dataset. If one does, the numeric suffixes will be incremented to make a distinction.

**EXTERNAL Subcommand**

EXTERNAL specifies that the data for each split-file group should be held in an external scratch file during processing. This helps conserve working space when running analyses with large datasets.

- The EXTERNAL subcommand takes no other keyword and is specified by itself.
- If time-dependent covariates exist, external data storage is unavailable, and EXTERNAL is ignored.

# CREATE

```
CREATE new series={CSUM (series)           }
                  {DIFF (series, order)    }
                  {FFT (series)            }
                  {IFFT (series)           }
                  {LAG (series, order [,order ])}
                  {LEAD (series, order [,order ])}
                  {MA (series, span [,minimum span]}
                  {PMA (series, span)      }
                  {RMED (series, span [,minimum span]}
                  {SDIFF (series, order [,periodicity]}
                  {T4253H (series)        }

[/new series=function (series {,span {,minimum span}})
                        {,order {,order }}
                        {,periodicity }
```

## *Function keywords:*

CSUM	Cumulative sum
DIFF	Difference
FFT	Fast Fourier transform
IFFT	Inverse fast Fourier transform
LAG	Lag
LEAD	Lead
MA	Centered moving averages
PMA	Prior moving averages
RMED	Running medians
SDIFF	Seasonal difference
T4253H	Smoothing

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
CREATE NEWVAR1 NEWVAR2 = CSUM(TICKETS RNDTRP).
```

## **Overview**

CREATE produces new series as a function of existing series. You can also use CREATE to replace the values of existing series.

CREATE displays a list of the new series, the case numbers of the first and last nonmissing cases, the number of valid cases, and the functions used to create the variables.

### **Basic Specification**

The basic specification is a new series name, an equals sign, a function, and the existing series, along with any additional specifications needed.

**Syntax Rules**

- The existing series together with any additional specifications (order, span, or periodicity) must be enclosed in parentheses.
- The equals sign is required.
- Series names and additional specifications must be separated by commas or spaces.
- You can specify only one function per equation.
- You can create more than one new series per equation by specifying more than one new series name on the left side of the equation and either multiple existing series names or multiple orders on the right.
- The number of new series named on the left side of the equation must equal the number of series created on the right. Note that the `FFT` function creates two new series for each existing series, and `IFFT` creates one series from two existing series.
- You can specify more than one equation on a `CREATE` command. Equations are separated by slashes.
- A newly created series can be specified in subsequent equations on the same `CREATE` command.

**Operations**

- Each new series created is added to the active dataset.
- If the new series named already exist, their values are replaced.
- If the new series named do not already exist, they are created.
- Series are created in the order in which they are specified on the `CREATE` command.
- If multiple series are created by a single equation, the first new series named is assigned the values of the first series created, the second series named is assigned the values of the second series created, and so on.
- `CREATE` automatically generates a variable label for each new series describing the function and series used to create it.
- The format of the new series is based on the function specified and the format of the existing series.
- `CREATE` honors the `TSET MISSING` setting that is currently in effect.
- `CREATE` does not honor the `USE` command.
- When an even-length span is specified for the functions `MA` and `RMED`, the centering algorithm uses an average of two spans of the specified length. The first span ranges from  $\text{span}/2$  cases before the current observation to the span length. The second span ranges from  $(\text{span}/2)-1$  cases before the current observation to the span length.

**Limitations**

- A maximum of 1 function per equation.
- There is no limit on the number of series created by an equation.
- There is no limit on the number of equations.

## Examples

```
CREATE NEWVAR1 = DIFF(OLDVAR,1) .
```

- In this example, the series *NEWVAR1* is created by taking the first-order difference of *OLDVAR*.

## CSUM Function

CSUM produces new series based on the cumulative sums of the existing series. Cumulative sums are the inverse of first-order differencing.

- The only specification on CSUM is the name or names of the existing series in parentheses.
- Cases with missing values in the existing series are not used to compute values for the new series. The values of these cases are system-missing in the new series.

### Example

```
CREATE NEWVAR1 NEWVAR2 = CSUM(TICKETS RNDTRP) .
```

- This example produces a new series called *NEWVAR1*, which is the cumulative sum of the series *TICKETS*, and a new series called *NEWVAR2*, which is the cumulative sum of the series *RNDTRP*.

## DIFF Function

DIFF produces new series based on nonseasonal differences of existing series.

- The specification on DIFF is the name or names of the existing series and the degree of differencing, in parentheses.
- The degree of differencing must be specified; there is no default.
- Since one observation is lost for each order of differencing, system-missing values will appear at the beginning of the new series.
- You can specify only one degree of differencing per DIFF function.
- If either of the pair of values involved in a difference computation is missing, the result is set to system-missing in the new series.

### Example

```
CREATE ADIF2 = DIFF(VARA,2) /
      YDIF1 ZDIF1 = DIFF(VARY VARZ,1) .
```

- The series *ADIF2* is created by differencing *VARA* twice.
- The series *YDIF1* is created by differencing *VARY* once.
- The series *ZDIF1* is created by differencing *VARZ* once.

## FFT Function

FFT produces new series based on fast Fourier transformations of existing series (Brigham, 1974).

- The only specification on FFT is the name or names of the existing series in parentheses.
- FFT creates two series, the cosine and sine parts (also called real and imaginary parts), for each existing series named. Thus, you must specify two new series names on the left side of the equation for each existing series specified on the right side.
- The first new series named becomes the real series, and the second new series named becomes the imaginary series.
- The existing series cannot have embedded missing values.
- The existing series must be of even length. If an odd-length series is specified, FFT pads it with a 0 to make it even. Alternatively, you can make the series even by adding or dropping an observation.
- The new series will be only half as long as the existing series. The remaining cases are assigned the system-missing value.

### Example

```
CREATE A B = FFT(C) .
```

- Two series, *A* (real) and *B* (imaginary), are created by applying a fast Fourier transformation to series *C*.

## IFFT Function

IFFT produces new series based on the inverse Fourier transformation of existing series.

- The only specification on IFFT is the name or names of the existing series in parentheses.
- IFFT needs two existing series to compute each new series. Thus, you must specify two existing series names on the right side of the equation for each new series specified on the left.
- The first existing series specified is the real series and the second series is the imaginary series.
- The existing series cannot have embedded missing values.
- The new series will be twice as long as the existing series. Thus, the last half of each existing series must be system-missing to allow enough room to create the new series.

### Example

```
CREATE C = IFFT(A B) .
```

- This command creates one new series, *C*, from the series *A* (real) and *B* (imaginary).

## LAG Function

LAG creates new series by copying the values of the existing series and moving them forward the specified number of observations. This number is called the **lag order**. The table below shows a first-order lag for a hypothetical dataset.

- The specification on `LAG` is the name or names of the existing series and one or two lag orders, in parentheses.
- At least one lag order must be specified; there is no default.
- Two lag orders indicate a range. For example, 2,6 indicates lag orders two through six. A new series is created for each lag order in the range.
- The number of new series specified must equal the number of existing series specified times the number of lag orders in the range.
- The first  $n$  cases at the beginning of the new series, where  $n$  is the lag order, are assigned the system-missing value.
- Missing values in the existing series are lagged and are assigned the system-missing value in the new series.
- A first-order lagged series can also be created using `COMPUTE`. `COMPUTE` does not cause a data pass (see `COMPUTE`).

Table 40-1  
First-order lag and lead of series  $X$

<b>X</b>	<b>Lag</b>	<b>Lead</b>
198	.	220
220	198	305
305	220	470
470	305	.

### Example

```
CREATE LAGVAR2 TO LAGVAR5 = LAG(VARA, 2, 5) .
```

- Four new variables are created based on lags on `VARA`. `LAGVAR2` is `VARA` lagged two steps, `LAGVAR3` is `VARA` lagged three steps, `LAGVAR4` is `VARA` lagged four steps, and `LAGVAR5` is `VARA` lagged five steps.

## LEAD Function

`LEAD` creates new series by copying the values of the existing series and moving them back the specified number of observations. This number is called the lead order.

- The specification on `LEAD` is the name or names of the existing series and one or two lead orders, in parentheses.
- At least one lead order must be specified; there is no default.
- Two lead orders indicate a range. For example, 1,5 indicates lead orders one through five. A new series is created for each lead order in the range.
- The number of new series must equal the number of existing series specified times the number of lead orders in the range.
- The last  $n$  cases at the end of the new series, where  $n$  equals the lead order, are assigned the system-missing value.
- Missing values in the existing series are moved back and are assigned the system-missing value in the new series.

**Example**

```
CREATE LEAD1 TO LEAD4 = LEAD(VARA, 1, 4) .
```

- Four new series are created based on leads of *VARA*. *LEAD1* is *VARA* led one step, *LEAD2* is *VARA* led two steps, *LEAD3* is *VARA* led three steps, and *LEAD4* is *VARA* led four steps.

**MA Function**

MA produces new series based on the centered moving averages of existing series.

- The specification on MA is the name or names of the existing series and the span to be used in averaging, in parentheses.
- A span must be specified; there is no default.
- If the specified span is odd, the MA is naturally associated with the middle term. If the specified span is even, the MA is centered by averaging each pair of uncentered means (Velleman and Hoaglin, 1981).
- After the initial span, a second span can be specified to indicate the minimum number of values to use in averaging when the number specified for the initial span is unavailable. This makes it possible to produce nonmissing values at or near the ends of the new series.
- The second span must be greater than or equal to 1 and less than or equal to the first span.
- The second span should be even (or 1) if the first span is even; it should be odd if the first span is odd. Otherwise, the next higher span value will be used.
- If no second span is specified, the minimum span is simply the value of the first span.
- If the number of values specified for the span or the minimum span is not available, the case in the new series is set to system-missing. Thus, unless a minimum span of 1 is specified, the endpoints of the new series will contain system-missing values.
- When MA encounters an embedded missing value in the existing series, it creates two subsets, one containing cases before the missing value and one containing cases after the missing value. Each subset is treated as a separate series for computational purposes.
- The endpoints of these subset series will have missing values according to the rules described above for the endpoints of the entire series. Thus, if the minimum span is 1, the endpoints of the subsets will be nonmissing; the only cases that will be missing in the new series are cases that were missing in the original series.

**Example**

```
CREATE TICKMA = MA(TICKETS, 4, 2) .
```

- This example creates the series *TICKMA* based on centered moving average values of the series *TICKETS*.
- A span of 4 is used for computing averages. At the endpoints, where four values are not available, the average is based on the specified minimum of two values.



## ***PMA Function***

PMA creates new series based on the prior moving averages of existing series. The prior moving average for each case in the original series is computed by averaging the values of a span of cases preceding it.

- The specification on PMA is the name or names of the existing series and the span to be used, in parentheses.
- Only one span can be specified and it is required. There is no default span.
- If the number of values specified for the span is not available, the case is set to system-missing. Thus, the number of cases with system-missing values at the beginning of the new series equals the number specified for the span.
- When PMA encounters an imbedded missing value in the existing series, it creates two subsets, one containing cases before the missing value and one containing cases after the missing value. Each subset is treated as a separate series for computational purposes. The first  $n$  cases in the second subset will be system-missing, where  $n$  is the span.

### ***Example***

```
CREATE PRIORA = PMA(VARA, 3) .
```

- This command creates the series *PRIORA* by computing prior moving averages for the series *VARA*. Since the span is 3, the first three cases in the series *PRIORA* are system-missing. The fourth case equals the average of cases 1, 2, and 3 of *VARA*, the fifth case equals the average of cases 2, 3, and 4 of *VARA*, and so on.

## ***RMED Function***

RMED produces new series based on the centered running medians of existing series.

- The specification on RMED is the name or names of the existing series and the span to be used in finding the median, in parentheses.
- A span must be specified; there is no default.
- If the specified span is odd, RMED is naturally the middle term. If the specified span is even, the RMED is centered by averaging each pair of uncentered medians (Velleman et al., 1981).
- After the initial span, a second span can be specified to indicate the minimum number of values to use in finding the median when the number specified for the initial span is unavailable. This makes it possible to produce nonmissing values at or near the ends of the new series.
- The second span must be greater than or equal to 1 and less than or equal to the first span.
- The second span should be even (or 1) if the first span is even; it should be odd if the first span is odd. Otherwise, the next higher span value will be used.
- If no second span is specified, the minimum span is simply the value of the first span.
- If the number of values specified for the span or the minimum span is not available, the case in the new series is set to system-missing. Thus, unless a minimum span of 1 is specified, the endpoints of the new series will contain system-missing values.

- When `RMED` encounters an imbedded missing value in the existing series, it creates two subsets, one containing cases before the missing value and one containing cases after the missing value. Each subset is treated as a separate series for computational purposes.
- The endpoints of these subset series will have missing values according to the rules described above for the endpoints of the entire series. Thus, if the minimum span is 1, the endpoints of the subsets will be nonmissing; the only cases that will be missing in the new series are cases that were missing in the original series.

### Example

```
CREATE TICKRMED = RMED(TICKETS,4,2).
```

- This example creates the series `TICKRMED` using centered running median values of the series `TICKETS`.
- A span of 4 is used for computing medians. At the endpoints, where four values are not available, the median is based on the specified minimum of two values.

## SDIFF Function

`SDIFF` produces new series based on seasonal differences of existing series.

- The specification on `SDIFF` is the name or names of the existing series, the degree of differencing, and, optionally, the periodicity, all in parentheses.
- The degree of differencing must be specified; there is no default.
- Since the number of seasons used in the calculations decreases by 1 for each order of differencing, system-missing values will appear at the beginning of the new series.
- You can specify only one degree of differencing per `SDIFF` function.
- If no periodicity is specified, the periodicity established on `TSET PERIOD` is in effect. If `TSET PERIOD` has not been specified, the periodicity established on the `DATE` command is used. If periodicity was not established anywhere, the `SDIFF` function cannot be executed.
- If either of the pair of values involved in a seasonal difference computation is missing, the result is set to system-missing in the new series.

### Example

```
CREATE SDVAR = SDIFF(VARA,1,12).
```

- The series `SDVAR` is created by applying one seasonal difference with a periodicity of 12 to the series `VARA`.

## T4253H Function

`T4253H` produces new series by applying a compound data smoother to the original series. The smoother starts with a running median of 4, which is centered by a running median of 2. It then resmooths these values by applying a running median of 5, a running median of 3, and hanning (running weighted averages). Residuals are computed by subtracting the smoothed series from the original series. This whole process is then repeated on the computed residuals. Finally, the

smoothed residuals are added to the smoothed values obtained the first time through the process (Velleman et al., 1981).

- The only specification on T4253H is the name or names of the existing series in parentheses.
- The existing series cannot contain imbedded missing values.
- Endpoints are smoothed through extrapolation and are not system-missing.

### **Example**

```
CREATE SMOOTHA = T4253H(VARA) .
```

- The series *SMOOTHA* is a smoothed version of the series *VARA*.

## **References**

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# CROSSTABS

CROSSTABS is available in the Statistics Base option.

## *General mode:*

CROSSTABS is available in the Statistics Base option.

```
CROSSTABS [TABLES=]varlist BY varlist [BY...] [/varlist...]
  [/MISSING={TABLE**}]
             {INCLUDE}
  [/WRITE[={NONE**}]]
             {CELLS }
  [/HIDESMALLCOUNTS [COUNT = {5          }]] ]
                               {integer}
  [/SHOWDIM = integer]
```

## *Integer mode :*

```
CROSSTABS VARIABLES=varlist(min,max) [varlist...]
  /TABLES=varlist BY varlist [BY...] [/varlist...]
  [/MISSING={TABLE**}]
             {INCLUDE}
             {REPORT }
  [/WRITE[={NONE**}]]
             {CELLS }
             {ALL   }
```

## *Both modes:*

```
[/FORMAT= {AVALUE**} {TABLES**}]
           {DVALUE  } {NOTABLES}
[/COUNT = [{ASIS}] [{ROUND  }]]
           {CASE   } {TRUNCATE}
           {CELL   }
[/CELLS=[{COUNT**}] [ROW    ] [EXPECTED] [SRESID ] [PROP  ]]
           {NONE   } [COLUMN] [RESID  ] [ASRESID] [BPROP]
                   [TOTAL ] [ALL    ]
[/STATISTICS=[CHISQ] [LAMBDA] [BTAU ] [GAMMA ] [ETA   ]]
              [PHI  ] [UC    ] [CTAU ] [D      ] [CORR  ]
              [CC   ] [RISK  ] [KAPPA] [MCNEMAR] [CMH(1*)]
              [ALL  ] [NONE  ]
[/METHOD={MC [CIN({99.0 })] [SAMPLES({10000})]}}]++
          {value} {value}
          {EXACT [TIMER({5  })] }
          {value}
[/BARCHART]
```

\*\*Default if the subcommand is omitted.

†† The `METHOD` subcommand is available only if the Exact Tests option is installed (available only on Windows operating systems).

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 19.0

- `HIDESMALLCOUNTS` subcommand introduced.
- `SHOWDIM` subcommand introduced.
- `PROP` and `BPROP` keywords introduced on the `CELLS` subcommand.

### **Example**

```
CROSSTABS TABLES=FEAR BY SEX  
/CELLS=ROW COLUMN EXPECTED RESIDUALS  
/STATISTICS=CHISQ.
```

## **Overview**

`CROSSTABS` produces contingency tables showing the joint distribution of two or more variables that have a limited number of distinct values. The frequency distribution of one variable is subdivided according to the values of one or more variables. The unique combination of values for two or more variables defines a cell.

`CROSSTABS` can operate in two different modes: *general* and *integer*. Integer mode builds some tables more efficiently but requires more specifications than general mode. Some subcommand specifications and statistics are available only in integer mode.

### **Options**

**Methods for building tables.** To build tables in general mode, use the `TABLES` subcommand. Integer mode requires the `TABLES` and `VARIABLES` subcommands and minimum and maximum values for the variables.

**Cell contents.** By default, `CROSSTABS` displays only the number of cases in each cell. You can request row, column, and total percentages, and also expected values and residuals, by using the `CELLS` subcommand.

**Statistics.** In addition to the tables, you can obtain measures of association and tests of hypotheses for each subtable using the `STATISTICS` subcommand.

**Formatting options.** With the `FORMAT` subcommand, you can control the display order for categories in rows and columns of subtables and suppress crosstabulation. With the `SHOWDIM` subcommand you can display a subset of the variables as table layers in the crosstabulation table.

**Writing and reproducing tables.** You can write cell frequencies to a file and reproduce the original tables with the `WRITE` subcommand.

**Basic specification**

In general mode, the basic specification is `TABLES` with a table list. The actual keyword `TABLES` can be omitted. In integer mode, the minimum specification is the `VARIABLES` subcommand, specifying the variables to be used and their value ranges, and the `TABLES` subcommand with a table list.

- The minimum table list specifies a list of row variables, the keyword `BY`, and a list of column variables.
- In integer mode, all variables must be numeric with integer values. In general mode, variables can be numeric (integer or non-integer) or string.
- The default table shows cell counts.

**Subcommand order**

- In general mode, the table list must be first if the keyword `TABLES` is omitted. If the keyword `TABLES` is explicitly used, subcommands can be specified in any order.
- In integer mode, `VARIABLES` must precede `TABLES`. The keyword `TABLES` must be explicitly specified.

**Operations**

- Integer mode builds tables more quickly but requires more workspace if a table has many empty cells.
- Statistics are calculated separately for each two-way table or two-way subtable. Missing values are reported for the table as a whole.
- In general mode, the keyword `TO` on the `TABLES` subcommand refers to the order of variables in the active dataset. `ALL` refers to all variables in the active dataset. In integer mode, `TO` and `ALL` refer to the position and subset of variables specified on the `VARIABLES` subcommand.

**Limitations**

The following limitations apply to `CROSSTABS` in *general mode*:

- A maximum of 200 variables named or implied on the `TABLES` subcommand
- A maximum of 1000 non-empty rows or columns for each table
- A maximum of 20 table lists per `CROSSTABS` command
- A maximum of 10 dimensions (9 `BY` keywords) per table
- A maximum of 400 value labels displayed on any single table

The following limitations apply to `CROSSTABS` in *integer mode*:

- A maximum of 100 variables named or implied on the `VARIABLES` subcommand
- A maximum of 100 variables named or implied on the `TABLES` subcommand
- A maximum of 1000 non-empty rows or columns for each table
- A maximum of 20 table lists per `CROSSTABS` command
- A maximum of 8 dimensions (7 `BY` keywords) per table

- A maximum of 20 rows or columns of missing values when REPORT is specified on MISSING
- The minimum value that can be specified is -99,999
- The maximum value that can be specified is 999,999

## Examples

### **Nominal by nominal relationships**

```
CROSSTABS
  /TABLES=store BY service
  /FORMAT= AVALUE TABLES
  /STATISTIC=CHISQ CC PHI LAMBDA UC
  /CELLS= COUNT .
```

- The procedure produces a crosstabulation table for *store* by *service*, with row and column variables displayed in ascending value.
- The CELLS subcommand specifies that only counts are displayed in the cells of the table.
- The STATISTICS subcommand requests a chi-square test and nominal-by-nominal measures of association .

### **Ordinal by ordinal relationships**

```
CROSSTABS
  /TABLES=regular BY overall
  /FORMAT= AVALUE TABLES
  /STATISTIC=D BTAU CTAU GAMMA
  /CELLS= COUNT .
```

- The procedure produces a crosstabulation table for *regular* by *overall*, with row and column variables displayed in ascending value.
- The CELLS subcommand specifies that only counts are displayed in the cells of the table.
- The STATISTICS subcommand requests measures of ordinal-by-ordinal association.

### **Relative risk**

```
CROSSTABS
  /TABLES=news BY response
  /FORMAT= AVALUE TABLES
  /STATISTIC= RISK
  /CELLS= COUNT ROW.
```

- The procedure produces a crosstabulation table for *news* by *response*, with row and column variables displayed in ascending value.
- The CELLS subcommand specifies that counts and row percentages are displayed in the cells of the table.
- The STATISTICS subcommand requests relative risk.

### **Measuring agreement**

```
CROSSTABS
  /TABLES=cons1 BY cons2
```

## CROSSTABS

```

/FORMAT= AVALUE TABLES
/STATISTIC=KAPPA
/CELLS= COUNT .

```

- The procedure produces a crosstabulation table for *cons* by *cons2*, with row and column variables displayed in ascending value.
- The `CELLS` subcommand specifies that only counts are displayed in the cells of the table.
- The `STATISTICS` subcommand requests the Kappa measure of agreement.

**Column proportions tests**

```

CROSSTABS
/TABLES=empcat BY jobsat
/FORMAT=AVALUE TABLES
/STATISTICS=CHISQ
/CELLS=COLUMN BPROP
/COUNT ROUND CELL.

```

- The procedure produces a crosstabulation table for *empcat* by *jobsat*, with row and column variables displayed in ascending value.
- The `STATISTICS` subcommand requests a chi-square test.
- The `CELLS` subcommand requests a column proportions test that utilizes the Bonferroni correction and specifies that only column percentages are displayed in the cells of the table.
- The `COUNT` subcommand specifies that case weights are used as is but the accumulated weights in the cells are rounded before computing any statistics.

**VARIABLES subcommand**

The `VARIABLES` subcommand is required for integer mode. `VARIABLES` specifies a list of variables to be used in the crosstabulations and the lowest and highest values for each variable. Values are specified in parentheses and must be integers. Non-integer values are truncated.

- Variables can be specified in any order. However, the order in which they are named on `VARIABLES` determines their implied order on `TABLES` (see the `TABLES` subcommand below).
- A range must be specified for each variable. If several variables can have the same range, it can be specified once after the last variable to which it applies.
- `CROSSTABS` uses the specified ranges to allocate tables. One cell is allocated for each possible combination of values of the row and column variables before the data are read. Thus, if the specified ranges are larger than the actual ranges, workspace will be wasted.
- Cases with values outside the specified range are considered missing and are not used in the computation of the table. This allows you to select a subset of values within `CROSSTABS`.
- If the table is sparse because the variables do not have values throughout the specified range, consider using general mode or recoding the variables.

**Example**

```
CROSSTABS VARIABLES=FEAR SEX RACE (1,2) MOBILE16 (1,3)
```



```
/TABLES=FEAR BY SEX MOBILE16 BY RACE.
```

- `VARIABLES` defines values 1 and 2 for *FEAR*, *SEX*, and *RACE* and values 1, 2, and 3 for *MOBILE16*.

## ***TABLES* subcommand**

`TABLES` specifies the table lists and is required in both integer mode and general mode. The following rules apply to both modes:

- You can specify multiple `TABLES` subcommands on a single `CROSSTABS` command. The slash between the subcommands is required; the keyword `TABLES` is required only in integer mode.
- Variables named before the first `BY` on a table list are row variables, and variables named after the first `BY` on a table list are column variables.
- When the table list specifies two dimensions (one `BY` keyword), the first variable before `BY` is crosstabulated with each variable after `BY`, then the second variable before `BY` with each variable after `BY`, and so on.
- Each subsequent use of the keyword `BY` on a table list adds a new dimension to the tables requested. Variables named after the second (or subsequent) `BY` are control variables.
- When the table list specifies more than two dimensions, a two-way subtable is produced for each combination of values of control variables. The value of the last specified control variable changes the most slowly in determining the order in which tables are displayed.
- You can name more than one variable in each dimension.

### ***General mode***

- The actual keyword `TABLES` can be omitted in general mode.
- In general mode, both numeric and string variables can be specified.
- The keywords `ALL` and `TO` can be specified in any dimension. In general mode, `TO` refers to the order of variables in the active dataset and `ALL` refers to all variables defined in the active dataset.

#### ***Example***

```
CROSSTABS TABLES=FEAR BY SEX BY RACE.
```

- This example crosstabulates *FEAR* by *SEX* controlling for *RACE*. In each subtable, *FEAR* is the row variable and *SEX* is the column variable.
- A subtable is produced for each value of the control variable *RACE*.

#### ***Example***

```
CROSSTABS TABLES=CONFINAN TO CONARMY BY SEX TO REGION.
```

- This command produces crosstabulations of all variables in the active dataset between and including *CONFINAN* and *CONARMY* by all variables between and including *SEX* and *REGION*.

### Integer mode

- In integer mode, variables specified on TABLES must first be named on VARIABLES.
- The keywords TO and ALL can be specified in any dimension. In integer mode, TO and ALL refer to the position and subset of variables specified on the VARIABLES subcommand, not to the variables in the active dataset.

#### Example

```
CROSSTABS VARIABLES=FEAR (1,2) MOBILE16 (1,3)
/TABLES=FEAR BY MOBILE16.
```

- VARIABLES names two variables, *FEAR* and *MOBILE16*. Values 1 and 2 for *FEAR* are used in the tables, and values 1, 2, and 3 are used for the variable *MOBILE16*.
- TABLES specifies a Crosstabulation table with two rows (values 1 and 2 for *FEAR*) and three columns (values 1, 2, and 3 for *MOBILE16*). *FEAR* and *MOBILE16* can be named on TABLES because they were named on the previous VARIABLES subcommand.

#### Example

```
CROSSTABS VARIABLES=FEAR SEX RACE DEGREE (1,2)
/TABLES=FEAR BY SEX BY RACE BY DEGREE.
```

- This command produces four subtables. The first subtable crosstabulates *FEAR* by *SEX*, controlling for the first value of *RACE* and the first value of *DEGREE*; the second subtable controls for the second value of *RACE* and the first value of *DEGREE*; the third subtable controls for the first value of *RACE* and the second value of *DEGREE*; and the fourth subtable controls for the second value of *RACE* and the second value of *DEGREE*.

### CELLS subcommand

By default, CROSSTABS displays only the number of cases in each cell of the Crosstabulation table. Use CELLS to display row, column, or total percentages, expected counts, or residuals. These are calculated separately for each Crosstabulation table or subtable.

- CELLS specified without keywords displays cell counts plus row, column, and total percentages for each cell.
- If CELLS is specified with keywords, CROSSTABS displays only the requested cell information.
- Scientific notation is used for cell contents when necessary.
- BPROP overrides PROP if both are specified.
- If BPROP or PROP is specified without specifying COUNT or COLUMN, then the observed cell counts are included in the Crosstabulation table, with APA-style subscripts indicating the results of the column proportions tests.

<b>COUNT</b>	<i>Observed cell counts.</i> This is the default if CELLS is omitted.
<b>ROW</b>	<i>Row percentages.</i> The number of cases in each cell in a row is expressed as a percentage of all cases in that row.
<b>COLUMN</b>	<i>Column percentages.</i> The number of cases in each cell in a column is expressed as a percentage of all cases in that column.

<b>TOTAL</b>	<i>Two-way table total percentages.</i> The number of cases in each cell of a subtable is expressed as a percentage of all cases in that subtable.
<b>EXPECTED</b>	<i>Expected counts.</i> Expected counts are the number of cases expected in each cell if the two variables in the subtable are statistically independent.
<b>RESID</b>	<i>Residuals.</i> Residuals are the difference between the observed and expected cell counts.
<b>SRESID</b>	<i>Standardized residuals</i> (Haberman, 1978).
<b>ASRESID</b>	<i>Adjusted standardized residuals</i> (Haberman, 1978).
<b>ALL</b>	<i>All cell information.</i> This includes cell counts; row, column, and total percentages; expected counts; residuals; standardized residuals; adjusted standardized residuals; and pairwise comparison of column proportions using the Bonferroni correction.
<b>NONE</b>	<i>No cell information.</i> Use NONE when you want to write tables to a procedure output file without displaying them. <a href="#">For more information, see the topic WRITE subcommand on p. 401.</a> This is the same as specifying NOTABLES on FORMAT.
<b>PROP</b>	<i>Pairwise comparison of column proportions.</i> Indicates which pairs of columns (for a given row) are significantly different. Significant differences (at the 0.05 level) are indicated with APA-style formatting using subscript letters. PROP is only available in general mode.
<b>BPROP</b>	<i>Pairwise comparison of column proportions using the Bonferroni correction.</i> Indicates which pairs of columns (for a given row) are significantly different, making use of the Bonferroni correction. Significant differences (at the 0.05 level) are indicated with APA-style formatting using subscript letters. BPROP is only available in general mode.

### **Example: pairwise comparison of column proportions**

```
CROSSTABS
  /TABLES= news BY inccat
  /FORMAT=AVALUE TABLES
  /CELLS=COLUMN BPROP
  /COUNT ROUND CELL.
```

Figure 41-1

#### **Newspaper subscription \* Income category in thousands Crosstabulation**

% within Income category in thousands

		Income category in thousands				Total
		Under \$25	\$25 - \$49	\$50 - \$74	\$75+	
Newspaper subscription	Yes	35.9% <sub>a</sub>	36.7% <sub>a</sub>	43.4% <sub>b</sub>	57.3% <sub>c</sub>	43.3%
	No	64.1% <sub>a</sub>	63.3% <sub>a</sub>	56.6% <sub>b</sub>	42.7% <sub>c</sub>	56.8%
Total		100.0%	100.0%	100.0%	100.0%	100.0%

The column proportions test assigns a subscript letter to the categories of the column variable. For each pair of columns, the column proportions are compared using a z test. If a pair of values is significantly different, the values have different subscript letters assigned to them.

The table in this example is a crosstabulation of survey respondents who have a newspaper subscription by the income category of the respondent, with column percentages shown as the summary statistic. The percentages in the *Under \$25* and *\$25 - \$49* categories both have the subscript *a* so the percentages in those columns are not significantly different. However, the subscripts in the *\$50 - \$74* and *\$75+* categories differ from each other as well as from the subscript for the *Under \$25* and *\$25 - \$49* categories. This means that the percentages in the *\$50 - \$74*

and \$75+ categories are significantly different from each other as well as from the percentages in the *Under \$25* and *\$25 - \$49* categories.

## **STATISTICS subcommand**

STATISTICS requests measures of association and related statistics. By default, CROSSTABS does not display any additional statistics.

- STATISTICS without keywords displays the chi-square test.
- If STATISTICS is specified with keywords, CROSSTABS calculates only the requested statistics.
- In integer mode, values that are not included in the specified range are *not* used in the calculation of the statistics, even if these values exist in the data.
- If user-missing values are included with MISSING, cases with user-missing values are included in the calculation of statistics as well as in the tables.

<b>CHISQ</b>	<i>Display the Chi-Square Test table.</i> Chi-square statistics include Pearson chi-square, likelihood-ratio chi-square, and Mantel-Haenszel chi-square (linear-by-linear association). Mantel-Haenszel is valid only if both variables are numeric. Fisher's exact test and Yates' corrected chi-square are computed for all $2 \times 2$ tables. This is the default if STATISTICS is specified with no keywords.
<b>PHI</b>	<i>Display phi and Cramér's V in the Symmetric Measures table.</i>
<b>CC</b>	<i>Display contingency coefficient in the Symmetric Measures table.</i>
<b>LAMBDA</b>	<i>Display lambda (symmetric and asymmetric) and Goodman and Kruskal's tau in the Directional Measures table.</i>
<b>UC</b>	<i>Display uncertainty coefficient (symmetric and asymmetric) in the Directional Measures table.</i>
<b>BTAU</b>	<i>Display Kendall's tau-b in the Symmetric Measures table.</i>
<b>CTAU</b>	<i>Display Kendall's tau-c in the Symmetric Measures table.</i>
<b>GAMMA</b>	<i>Display gamma in the Symmetric Measures table or Zero-Order and Partial Gammas table.</i> The Zero-Order and Partial Gammas table is produced only for tables with more than two variable dimensions.
<b>D</b>	<i>Display Somers' d (symmetric and asymmetric) in the Directional Measures table.</i>
<b>ETA</b>	<i>Display eta in the Directional Measures table.</i> Available for numeric data only.
<b>CORR</b>	<i>Display Pearson's r and Spearman's correlation coefficient in the Symmetric Measures table.</i> This is available for numeric data only.
<b>KAPPA</b>	<i>Display kappa coefficient (Kraemer, 1982) in the Symmetric Measures table.</i> Kappa is based on a square table in which row and column values represent the same scale. Any cell that has observed values for one variable but not the other is assigned a count of 0. Kappa is not computed if the data storage type (string or numeric) is not the same for the two variables. For string variable, both variables must have the same defined length.
<b>RISK</b>	<i>Display relative risk (Bishop, Feinberg, and Holland, 1975) in the Risk Estimate table.</i> Relative risk can be calculated only for $2 \times 2$ tables.
<b>MCNEMAR</b>	<i>Display a test of symmetry for square tables.</i> The McNemar test is displayed for $2 \times 2$ tables, and the McNemar-Bowker test, for larger tables.

<b>CMH(1*)</b>	<i>Conditional independence and homogeneity tests.</i> Cochran's and the Mantel-Haenszel statistics are computed for the test for conditional independence. The Breslow-Day and Tarone's statistics are computed for the test for homogeneity. For each test, the chi-squared statistic with its degrees of freedom and asymptotic <i>p</i> value are computed. <i>Mantel-Haenszel relative risk (common odds ratio) estimate.</i> The Mantel-Haenszel relative risk (common odds ratio) estimate, the natural log of the estimate, the standard error of the natural log of the estimate, the asymptotic <i>p</i> value, and the asymptotic confidence intervals for common odds ratio and for the natural log of the common odds ratio are computed. The user can specify the null hypothesis for the common odds ratio in parentheses after the keyword. The passive default is 1. (The parameter value must be positive.)
<b>ALL</b>	<i>All statistics available.</i>
<b>NONE</b>	<i>No summary statistics.</i> This is the default if <code>STATISTICS</code> is omitted.

## **METHOD subcommand**

`METHOD` displays additional results for each statistic requested. If no `METHOD` subcommand is specified, the standard asymptotic results are displayed. If fractional weights have been specified, results for all methods will be calculated on the weight rounded to the nearest integer. This subcommand is available only if you have the Exact Tests add-on option installed, which is only available on Windows operating systems.

<b>MC</b>	Displays an unbiased point estimate and confidence interval based on the Monte Carlo sampling method, for all statistics. Asymptotic results are also displayed. When exact results can be calculated, they will be provided instead of the Monte Carlo results.
<b>CIN(n)</b>	Controls the confidence level for the Monte Carlo estimate. <code>CIN</code> is available only when <code>/METHOD=MC</code> is specified. <code>CIN</code> has a default value of 99.0. You can specify a confidence interval between 0.01 and 99.9, inclusive.
<b>SAMPLES</b>	Specifies the number of tables sampled from the reference set when calculating the Monte Carlo estimate of the exact <i>p</i> value. Larger sample sizes lead to narrower confidence limits but also take longer to calculate. You can specify any integer between 1 and 1,000,000,000 as the sample size. <code>SAMPLES</code> has a default value of 10,000.
<b>EXACT</b>	Computes the exact significance level for all statistics in addition to the asymptotic results. <code>EXACT</code> and <code>MC</code> are mutually exclusive alternatives (you cannot specify both on the same command). Calculating the exact <i>p</i> value can be memory-intensive. If you have specified <code>/METHOD=EXACT</code> and find that you have insufficient memory to calculate results, you should first close any other applications that are currently running in order to make more memory available. You can also enlarge the size of your swap file (see your Windows documentation for more information). If you still cannot obtain exact results, specify <code>/METHOD=MC</code> to obtain the Monte Carlo estimate of the exact <i>p</i> value. An optional <code>TIMER</code> keyword is available if you choose <code>/METHOD=EXACT</code> .
<b>TIMER(n)</b>	Specifies the maximum number of minutes allowed to run the exact analysis for each statistic. If the time limit is reached, the test is terminated, no exact results are provided, and the program begins to calculate the next test in the analysis. <code>TIMER</code> is available only when <code>/METHOD=EXACT</code> is specified. You can specify any integer value for <code>TIMER</code> . Specifying a value of 0 for <code>TIMER</code> turns the timer off completely. <code>TIMER</code> has a default value of 5 minutes. If a test exceeds a time limit of 30 minutes, it is recommended that you use the Monte Carlo, rather than the exact, method.

**Example**

```
CROSSTABS TABLES=FEAR BY SEX
/CELLS=ROW COLUMN EXPECTED RESIDUALS
/STATISTICS=CHISQ
/METHOD=MC SAMPLES(10000) CIN(95).
```

- This example requests chi-square statistics.
- An unbiased point estimate and confidence interval based on the Monte Carlo sampling method are displayed with the asymptotic results.

**MISSING subcommand**

By default, CROSSTABS deletes cases with missing values on a table-by-table basis. Cases with missing values for any variable specified for a table are not used in the table or in the calculation of statistics. Use MISSING to specify alternative missing-value treatments.

- The only specification is a single keyword.
- The number of missing cases is always displayed in the Case Processing Summary table.
- If the missing values are not included in the range specified on VARIABLES, they are excluded from the table regardless of the keyword you specify on MISSING.

<b>TABLE</b>	<i>Delete cases with missing values on a table-by-table basis.</i> When multiple table lists are specified, missing values are handled separately for each list. This is the default.
<b>INCLUDE</b>	<i>Include user-missing values.</i>
<b>REPORT</b>	<i>Report missing values in the tables.</i> This option includes missing values in tables but not in the calculation of percentages or statistics. The missing status is indicated on the categorical label. REPORT is available only in integer mode.

**FORMAT subcommand**

By default, CROSSTABS displays tables and subtables. The values for the row and column variables are displayed in order from lowest to highest. Use FORMAT to modify the default table display.

<b>AVALUE</b>	<i>Display row and column variables from lowest to highest value.</i> This is the default.
<b>DVALUE</b>	<i>Display row variables from highest to lowest.</i> This setting has no effect on column variables.
<b>TABLES</b>	<i>Display tables.</i> This is the default.
<b>NOTABLES</b>	<i>Suppress Crosstabulation tables.</i> NOTABLES is useful when you want to write tables to a file without displaying them or when you want only the Statistics table. This is the same as specifying NONE on CELLS.

## ***COUNT subcommand***

The `COUNT` subcommand controls how case weights are handled.

<b>ASIS</b>	<i>The case weights are used as is. However, when Exact Statistics are requested, the accumulated weights in the cells are either truncated or rounded before computing the Exact test statistics.</i>
<b>CASE</b>	<i>The case weights are either rounded or truncated before use.</i>
<b>CELL</b>	<i>The case weights are used as is but the accumulated weights in the cells are either truncated or rounded before computing any statistics.</i>
<b>ROUND</b>	<i>Performs Rounding operation.</i>
<b>TRUNCATE</b>	<i>Performs Truncation operation.</i>

## ***BARCHART subcommand***

`BARCHART` produces a clustered bar chart where bars represent categories defined by the first variable in a crosstabulation while clusters represent categories defined by the second variable in a crosstabulation. Any controlling variables in a crosstabulation are collapsed over before the clustered bar chart is created.

- `BARCHART` takes no further specification.
- If integer mode is in effect and `MISSING=REPORT`, `BARCHART` displays valid and user-missing values. Otherwise only valid values are used.

## ***WRITE subcommand***

Use the `WRITE` subcommand to write cell frequencies to a file for subsequent use by the current program or another program. `CROSSTABS` can also use these cell frequencies as input to reproduce tables and compute statistics. When `WRITE` is specified, an Output File Summary table is displayed before all other tables. See the [OMS](#) command for alternative and more flexible methods for writing out results to external files in various formats.

- The only specification is a single keyword.
- The name of the file must be specified on the `PROCEDURE OUTPUT` command preceding `CROSSTABS`.
- If you include missing values with `INCLUDE` or `REPORT` on `MISSING`, no values are considered missing and all non-empty cells, including those with missing values, are written, even if `CELLS` is specified.
- If you exclude missing values on a table-by-table basis (the default), no records are written for combinations of values that include a missing value.
- If multiple tables are specified, the tables are written in the same order as they are displayed.

## CROSSTABS

- `WRITE` is not supported for long string variables (defined width greater than 8 bytes) and will result in an error if any of the variables are long strings.
- `WRITE` is not supported for tables with more than eight dimensions (seven `BY` variables) and will result in an error if the table has more than eight dimensions.

<b>NONE</b>	<i>Do not write cell counts to a file.</i> This is the default.
<b>CELLS</b>	<i>Write cell counts for non-empty and nonmissing cells to a file.</i> Combinations of values that include a missing value are not written to the file.
<b>ALL</b>	<i>Write cell counts for all cells to a file.</i> A record for each combination of values defined by <code>VARIABLES</code> and <code>TABLES</code> is written to the file. <code>ALL</code> is available only in integer mode.

The file contains one record for each cell. Each record contains the following:

Columns	Contents
1–4	<i>Split-file group number, numbered consecutively from 1.</i> Note that this is not the value of the variable or variables used to define the splits.
5–8	<i>Table number.</i> Tables are defined by the <code>TABLES</code> subcommand.
9–16	<i>Cell frequency.</i> The number of times this combination of variable values occurred in the data, or, if case weights are used, the sum of case weights for cases having this combination of values.
17–24	<i>The value of the row variable</i> (the one named before the first <code>BY</code> ).
25–32	<i>The value of the column variable</i> (the one named after the first <code>BY</code> ).
33–40	<i>The value of the first control variable</i> (the one named after the second <code>BY</code> ).
41–48	<i>The value of the second control variable</i> (the one named after the third <code>BY</code> ).
49–56	<i>The value of the third control variable</i> (the one named after the fourth <code>BY</code> ).
57–64	<i>The value of the fourth control variable</i> (the one named after the fifth <code>BY</code> ).
65–72	<i>The value of the fifth control variable</i> (the one named after the sixth <code>BY</code> ).
73–80	<i>The value of the sixth control variable</i> (the one named after the seventh <code>BY</code> ).

- The split-file group number, table number, and frequency are written as integers.
- In integer mode, the values of variables are also written as integers. In general mode, the values are written according to the print format specified for each variable. Alphanumeric values are written at the left end of any field in which they occur.
- Within each table, records are written from one column of the table at a time, and the value of the last control variable changes the most slowly.

**Example**

```
PROCEDURE OUTPUT  OUTFILE='/data/celldata.txt'.
CROSSTABS VARIABLES=FEAR SEX (1,2)
  /TABLES=FEAR BY SEX
  /WRITE=ALL.
```

- `CROSSTABS` writes a record for each cell in the table *FEAR* by *SEX* to the file *celldata.txt*.

**Example**

```
PROCEDURE OUTPUT  OUTFILE='/data/xtabdata.txt'.
CROSSTABS  TABLES=V1 TO V3 BY V4 BY V10 TO V15
```



```
/WRITE=CELLS.
```

- CROSSTABS writes a set of records for each table to file *xtabdata.txt*.
- Records for the table *V1* by *V4* by *V10* are written first, followed by records for *V1* by *V4* by *V11*, and so on. The records for *V3* by *V4* by *V15* are written last.

### Reading a CROSSTABS Procedure Output file

You can use the file created by `WRITE` in a subsequent session to reproduce a table and compute statistics for it. Each record in the file contains all of the information used to build the original table. The cell frequency information can be used as a weight variable on the `WEIGHT` command to replicate the original cases.

#### Example

```
DATA LIST FILE='/celldata.txt'
  /WGHT 9-16 FEAR 17-24 SEX 25-32.
VARIABLE LABELS FEAR 'AFRAID TO WALK AT NIGHT IN NEIGHBORHOODS'.
VALUE LABELS FEAR 1 'YES' 2 'NO' / SEX 1 'MALE' 2 'FEMALE'.
WEIGHT BY WGHT.
CROSSTABS TABLES=FEAR BY SEX
  /STATISTICS=ALL.
```

- `DATA LIST` reads the cell frequencies and row and column values from the *celldata.txt* file. The cell frequency is read as a weighting factor (variable *WGHT*). The values for the rows are read as *FEAR*, and the values for the columns are read as *SEX*, the two original variables.
- The `WEIGHT` command recreates the sample size by weighting each of the four cases (cells) by the cell frequency.

If you do not have the original data or the `CROSSTABS` procedure output file, you can reproduce a crosstabulation and compute statistics simply by entering the values from the table:

```
DATA LIST /FEAR 1 SEX 3 WGHT 5-7.
VARIABLE LABELS FEAR 'AFRAID TO WALK AT NIGHT IN NEIGHBORHOOD'.
VALUE LABELS FEAR 1 'YES' 2 'NO' / SEX 1 'MALE' 2 'FEMALE'.
WEIGHT BY WGHT.
BEGIN DATA
1 1 55
2 1 172
1 2 180
2 2 89
END DATA.
CROSSTABS TABLES=FEAR BY SEX
  /STATISTICS=ALL.
```

### HIDESMALLCOUNTS Subcommand

`HIDESMALLCOUNTS` allows you to hide counts displayed in tables for count values that are less than a specified integer. Hidden values are displayed as <N, where N is the specified integer.

**COUNT Keyword**

The COUNT keyword specifies the cutoff, below which count values are hidden. Values equal to the cutoff are not hidden. The value of COUNT must be an integer and greater than or equal to 2, although the value 0 is permitted and specifies that there is no cutoff. The default is 5.

**Notes**

- Percentages associated with hidden counts are also hidden.
- HIDE SMALL COUNTS is only available in general mode.

**SHOWDIM Subcommand**

SHOWDIM specifies the number of dimensions to be displayed as rows and columns in the crosstabulation table, with the remainder of the dimensions displayed as table layers.

- The specification must be a positive integer. For a specified value of  $n$ , the dimensions corresponding to the variables before the  $n$ th BY keyword, as well as cell statistics (such as Count and Expected Count), will be displayed as rows or columns in the crosstabulation table. Variables after the  $n$ th BY keyword will be displayed as layers in the crosstabulation table.
- For SHOWDIM = 1, values of the variables before the first BY keyword are shown as row headings and cell statistics are shown as column headings, in the crosstabulation table. All other dimensions appear in layers of the crosstabulation table.
- When SHOWDIM is omitted all dimensions will be displayed as rows or columns.
- An integer value greater than or equal to the command dimensionality (the number of BY keywords plus one) is equivalent to omitting the SHOWDIM subcommand.

*Note:* SHOWDIM is only available in general mode.

See [SET TABLERENDER](#) for information on how multiple layer dimensions are rendered in lightweight tables.

**References**

- Bishop, Y. M., S. E. Feinberg, and P. W. Holland. 1975. *Discrete multivariate analysis: Theory and practice*. Cambridge, Mass.: MIT Press.
- Haberman, S. J. 1978. *Analysis of qualitative data*. London: Academic Press.
- Kraemer, H. C. 1982. Kappa Coefficient. In: *Encyclopedia of Statistical Sciences*, S. Kotz, and N. L. Johnson, eds. New York: John Wiley and Sons.

# CSCOXREG

CSCOXREG is available in the Complex Samples option.

*Note:* Square brackets used in the CSCOXREG syntax chart are required parts of the syntax and are not used to indicate optional elements. Equals signs (=) used in the syntax chart are required elements. All subcommands are optional.

```
CSCOXREG starttime endtime BY factor list WITH covariate list

/VARIABLES STATUS = varname(valuelist)
           ID = varname BASELINESTRATA = varname

/PLAN FILE = 'file'

/JOINTPROB FILE = 'savfile' | 'dataset'

/MODEL effect list
/CUSTOM LABEL = 'label'
           LMATRIX = {effect list, effect list ...; ... }
                   {effect list, effect list ... }
                   {ALL list; ALL ... }
                   {ALL list }

           KMATRIX = {number; number; ...}
                   {number }

/CUSTOM ...

/CRITERIA CILEVEL = {95** } DF = number
                   {number}

           LCONVERGE = [{0** } {RELATIVE**}]
                   {number} {ABSOLUTE }

           MXITER = {100** } MXSTEP = {5** }
                   {integer} {integer}

           PCONVERGE = [{1E-6**} {RELATIVE**}]
                   {number} {ABSOLUTE }

           SINGULAR = {1E-12**}
                   {number }

           TIES = {EFRON**}
                   {BRESLOW}

/STATISTICS PARAMETER EXP SE TTEST CINTERVAL DEFF DEFFSQRT

/TEST TYPE = {F** } PADJUST = {LSD** }
            {ADJF } {BONFERRONI }
            {CHISQUARE } {SEQBONFERRONI}
            {ADJCHISQUARE} {SIDAK }
            {SEQSIDAK }

/TESTASSUMPTIONS PROPHAZARD = {KM** }
                               {IDENTITY }
                               {LOG }
                               {RANK }
PARAMETER COVB

/DOMAIN VARIABLE = varname (value)

/MISSING CLASSMISSING = {EXCLUDE**}
                       {INCLUDE }

/SURVIVALMETHOD BASELINE = {BRESLOW } CI = {ORIGINAL}
                          {EFRON } {LOG** }
                          {PRODUCTLIMIT} {LML }
```

```

/PRINT SAMPLEINFO** EVENTINFO** RISKINFO
      HISTORY({1**    }) GEF LMATRIX COVB CORB BASELINE NONE
          {integer}

/SAVE SCHOENFELD(rootname:{25**  }) MARTINGALE(varname) DEVIANCE(varname)
          {integer}
      COXSNEILL(varname) SCORE(rootname:{25**  }) DFBETA(rootname:{25**  })
          {integer}          {integer}
      AGGMARTINGALE(varname) AGGDEVIANCE(varname) AGGCOXSNEILL(varname)
      AGGSCORE(rootname:{25**  }) AGGDFBETA(rootname:{25**  })
          {integer}          {integer}
      SURVIVAL(varname) LCL_SURVIVAL(varname) UCL_SURVIVAL(varname)
      CUMHAZARD(varname) LCL_CUMHAZARD(varname) UCL_CUMHAZARD(varname)
      XBETA(varname)

/PLOT SURVIVAL CUMHAZARD LML OMS CI = {YES }
          {NO**}

/PATTERN {varname(value)...          }
          {endtime(valuelist) varname({valuelist}) ...}
          {value          }
      BY factor

/PATTERN ...

/OUTFILE {COVB = 'savfile' | 'dataset'    } {MODEL = 'file'    }
          {CORB = 'savfile' | 'dataset'    }
          {SURVIVAL = 'savfile' | 'dataset'}

```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

#### Release 16.0

- Command introduced.

#### Release 17.0

- Added support for SET THREADS.

### **Example**

```

CSCOXREG endtime_var BY a b c WITH x y z
/VARIABLES STATUS=status_var(1)
/PLAN FILE='/survey/myfile.csplan'.

```

## **Overview**

For samples drawn by complex sampling methods, CSCOXREG applies Cox proportional hazards regression to analysis of survival times—that is, the length of time before the occurrence of an event. CSCOXREG supports scale and categorical predictors, which can be time dependent. CSCOXREG provides an easy way of considering differences in subgroups as well as analyzing effects of a set of predictors. The procedure estimates variances by taking into account the sample design used to select the sample, including equal probability and probability proportional to size (PPS) methods and with replacement (WR) and without replacement (WOR) sampling procedures. Optionally, CSCOXREG performs analyses for a subpopulation.

### **Basic Specification**

- The basic specification is a variable list identifying the time variables (at least one but can be up to two), the factors (if any), and the covariates (if any); a `VARIABLES` subcommand specifying the event status variable; and a `PLAN` subcommand with the name of a complex sample analysis plan file, which may be generated by the `CSPLAN` procedure.
- The default model includes main effects for any factors and any covariates.
- The basic specification displays summary information about the sample and all analysis variables, model summary statistics, and Wald F tests for all model effects. Additional subcommands must be used for other output.
- Minimum syntax is a time variable, a status variable, and the `PLAN` subcommand. This specification fits a baseline-only model.

### **Syntax Rules**

- The *endtime* variable, `STATUS` in `VARIABLES` subcommand, and `PLAN` subcommand are required. All other variables and subcommands are optional.
- Multiple `CUSTOM` and `PATTERN` subcommands may be specified; each is treated independently. All other subcommands may be specified only once.
- Empty subcommands are not allowed; all subcommands must be specified with options.
- Each keyword may be specified only once within a subcommand.
- Subcommand names and keywords must be spelled in full.
- Equals signs (=) and slashes shown in the syntax chart are required.
- Bold square brackets shown in the syntax chart are required parts of the syntax and are not used to indicate optional elements.
- Subcommands may be specified in any order.
- The factors, `ID` variable, baseline strata variable, and the subpopulation can be numeric or string variables, but covariates must be numeric.
- Across the time variables, factor, and covariate variable lists, a variable may be specified only once.
- The status variable, `ID` variable, baseline strata variable, and subpopulation variables may not be specified on the variable list
- Minimum syntax is a time variable, a status variable, and the `PLAN` subcommand. This specification fits a baseline-only model.
- Only factors and covariates can be defined by the `TIME PROGRAM`, all other variables cannot be defined there. [For more information, see the topic TIME PROGRAM on p. 1962.](#)

### **Operations**

- `TIME PROGRAM` computes the values for time-dependent predictors (see `TIME PROGRAM` syntax help).
- `CSCOXREG` performs Cox proportional hazards regression analysis for sampling designs supported by the `CSPLAN` and `CSSELECT` procedures.

- The input dataset must contain the variables to be analyzed and variables related to the sampling design.
- The complex sample analysis plan file provides an analysis plan based on the sampling design.
- By default, CSCOXREG uses a model that includes main effects for any factors and any covariates.
- Other effects, including interaction and nested effects, may be specified using the MODEL subcommand.
- The default output for the specified model is summary information about the sample and all analysis variables, model summary statistics, and Wald F tests for all model effects.
- This procedure uses the multithreaded options specified by SET THREADS.

### Limitations

WEIGHT and SPLIT FILE settings are ignored with a warning by the CSCOXREG procedure.

## Examples

```
CSCOXREG t BY a b c WITH x
/VARIABLES STATUS = dead(1)
/PLAN FILE='c:\survey\myfile.csplan'.
```

- *t* is the time variable; *a*, *b*, and *c* are factors; *x* is a covariate.
- The status variable is *dead* with a value of 1 representing the terminal event.
- The complex sampling plan is given in the file *c:\survey\myfile.csplan*.
- CSCOXREG will fit the default model including the main effects for factors *a*, *b*, and *c* and the covariate *x*.

### Multiple Cases per Subject

```
* Complex Samples Cox Regression.
CSCOXREG start_time time_to_event BY mi is hs
/PLAN FILE='samplesDirectory\srs.csaplan'
/VARIABLES STATUS=event(4) ID=patid
/MODEL mi is hs
/PRINT SAMPLEINFO EVENTINFO
/STATISTICS PARAMETER EXP SE CINTERVAL
/PLOT LML CI=NO
/PATTERN is(1) hs(0) BY mi
/TEST TYPE=F PADJUST=LSD
/CRITERIA MXITER=100 MXSTEP=5 PCONVERGE=[1E-006 RELATIVE] LCONVERGE=[0] TIES=BRESLOW CILEVEL=95
/SURVIVALMETHOD BASELINE=EFRON CI=LOG
/MISSING CLASSMISSING=EXCLUDE.
```

- The CSCOXREG procedure creates a Cox regression model for survival times defined by *start\_time* and *time\_to\_event*, using *mi*, *is*, and *hs* as factors. The sampling design is defined in *srs.csaplan*.
- The VARIABLES subcommand specifies on the STATUS keyword that a value of 4 for *event* indicates that the terminal event (death) has occurred. The ID keyword specifies *patid* as the subject ID variable. All cases sharing the same value of *patid* belong to the same subject.

- The `STATISTICS` subcommand requests estimates, exponentiated estimates, standard errors, and confidence intervals for model parameters.
- The `PLOT` subcommand requests log-minus-log plots of the estimated survival for the reference pattern (which uses the highest value of each factor), plus each pattern defined in any `PATTERN` subcommands.
- The `PATTERN` subcommand requests a plot to be produced using 1 as the value for *is* and 0 as the value for *hs*. Separate lines in the plot will be produced for each value of *mi*.
- The `CRITERIA` subcommand requests that the Breslow method be used for breaking ties.
- All other options are set to their default values.

### ***Time-Dependent Covariates***

```
* Complex Samples Cox Regression.
CLEAR TIME PROGRAM.
TIME PROGRAM.
COMPUTE t_age=ln(T_)*age.
CSCOXREG time_to_event WITH age t_age
  /PLAN FILE='samplesDirectory\recidivism_cs.csplan'
  /JOINTPROB FILE='samplesDirectory\recidivism_cs_jointprob.sav'
  /VARIABLES STATUS=arrest2(1)
  /MODEL age t_age
  /PRINT SAMPLEINFO EVENTINFO
  /STATISTICS PARAMETER SE CINTERVAL DEFF
  /TEST TYPE=F PADJUST=LSD
  /CRITERIA MXITER=100 MXSTEP=5 PCONVERGE=[1E-006 RELATIVE]
  LCONVERGE=[0] TIES=EFRON CILEVEL=95
  /SURVIVALMETHOD BASELINE=EFRON CI=LOG
  /MISSING CLASSMISSING=EXCLUDE.
```

- The `TIME PROGRAM` command indicates that the following `COMPUTE` statement defines a time-dependent predictor for use with `CSCOXREG`. The time-dependent predictor is the interaction between the covariate *age* and the natural log of the internal time variable *T\_*.
- The `CSCOXREG` procedure fits a model for *time\_to\_event* given covariates *age* and *t\_age*. The sampling design is defined in *recidivism\_cs.csplan*, and joint probabilities are stored in *recidivism\_cs\_jointprob.sav*.
- The `VARIABLES` subcommand specifies that a value of 1 for *arrest2* indicates that the event of interest (rearrest) has occurred.
- The `STATISTICS` subcommand requests estimates, standard errors, confidence intervals, and design effects for model parameters.
- All other options are set to their default values.

## ***Variable List Subcommand***

The variable list specifies the time variable(s), the factors, and the covariates in the model.

- The time variables *starttime* (if specified) and *endtime* must be listed first. These variables represent the endpoints of a time interval (*starttime*, *endtime*) during which the case is at risk.

- When *starttime* is not specified, it is implied that *starttime* = 0 for all cases if an ID variable is not specified. If an ID variable is specified, it is assumed that *starttime* = 0 for the first observation for that subject and *starttime* for following observations equals *endtime* of the previous observation. See the example below.
- The time variables must be numeric and non-negative.
- If the time variables are of Date or Time type, the internal numeric representation will be used and a warning will be given. For example, November 8, 1957, is 1.2E+10 (the number of seconds from midnight, October 14, 1582). See [Date and time functions](#) for detailed internal numeric representation of different date and time formats.
- The names of the factors and covariates, if any, follow the time variables. Specify any factors following the keyword BY. Specify any covariates following the keyword WITH.
- Factors can be numeric or string variables, but covariates must be numeric.
- Each variable may be specified only once on the variable list.
- The status variable, ID variable, baseline strata variable, and subpopulation variables may not be specified on the variable list.

### Example

```
CSCOXREG tstart tend BY a b c WITH x
/VARIABLES STATUS = dead(1) ID = SSN
/PLAN FILE='c:\survey\myfile.csplan'
/MODEL a b c a*b a*c b*c x.
```

- Two time variables, *tstart* and *tend*, are specified.
- ID specifies *SSN* as the subject ID variable. All cases sharing the same value of *SSN* belong to the same subject.
- This example fits a model that includes the main effects for factors *a*, *b*, and *c*; all two-way interactions among the factors; and the covariate *x*.

```
CSCOXREG tend BY a b c WITH x
/VARIABLES STATUS = dead(1) ID = SSN
/PLAN FILE='c:\survey\myfile.csplan'
/MODEL a b c a*b a*c b*c x.
```

- This is the same as the above example except that only one time variable, *tend*, is specified.
- The values of *starttime* derived from *tend*. For example, for the following subject,

SSN	tend	
123456789	10	...
123456789	20	...
123456789	25	...

it is implied that *starttime* = 0, 10, and 20 for these three cases.



## VARIABLES Subcommand

VARIABLES specifies the status variable, ID variable, and baseline strata variable.

### STATUS = varname

*Event status variable.* To determine whether the event has occurred at *endtime* for a particular observation, CSCOXREG checks the value of a status variable. STATUS lists the status variable and the values that indicate the occurrence of the event.

The value list must be enclosed in parentheses. All cases with non-negative times that do not have a value within the range specified are classified as censored cases—that is, cases for which the event has not yet occurred at *endtime*.

The value list can be one value, a list of values separated by blanks or commas, a range of values using the keyword THRU, or a combination.

The status variable can be either numeric or string. If a string variable is specified, the event values must be enclosed in apostrophes and the keyword THRU cannot be used.

### ID = varname

*ID variable.* Cases with the same ID value are repeated observations from the same subject. If ID is not specified, each case represents one subject.

### BASELINESTRATA = varname

*Baseline stratification variable.* A separate baseline hazard and survival function is computed for each value of this variable, while a single set of model coefficients is estimated across strata.

### Example

```
CSCOXREG SURVIVAL by GROUP
/VARIABLES STATUS=SRVSTA(3 THRU 5, 8 THRU 10) BASELINESTRATA=LOCATION
/PLAN FILE='c:\survey\myfile.csplan'.
```

- STATUS specifies that *SRVSTA* is the status variable.
- A value between either 3 and 5 or 8 and 10, inclusive, means that the terminal event occurred.
- Values outside the specified ranges indicate censored cases.
- BASELINESTRATA specifies *LOCATION* as the strata variable.
- Different baseline survival functions are computed for each value of *LOCATION*.

## PLAN Subcommand

The PLAN subcommand specifies the name of an XML file containing analysis design specifications. This file is written by the CSPLAN procedure.

- The PLAN subcommand is required.

**FILE** *Specifies the name of an external file.*

## **JOINTPROB Subcommand**

The JOINTPROB subcommand is used to specify the file or dataset containing the first stage joint inclusion probabilities for UNEQUAL\_WOR estimation. The CSSELECT procedure writes this file in the same location and with the same name (but a different extension) as the plan file. When UNEQUAL\_WOR estimation is specified, the procedure will use the default location and name of the file unless the JOINTPROB subcommand is used to override them.

**FILE** *Specifies the name of the file or dataset containing the joint inclusion probabilities.*

## **MODEL Subcommand**

The MODEL subcommand is used to specify the effects to be included in the model.

- Specify a list of terms to be included in the model, separated by spaces or commas.
- If the MODEL subcommand is not specified, CSCOXREG uses a model that includes main effects for any factors, and any covariates, in the order specified on the variable list.
- To include a term for the main effect of a factor, enter the name of the factor.
- To include a term for an interaction among factors, use the keyword BY or the asterisk (\*) to join the factors involved in the interaction. For example, A\*B means a two-way interaction effect of A and B, where A and B are factors. A\*A is not allowed because factors inside an interaction effect must be distinct.
- To include a term for nesting one factor within another, use a pair of parentheses. For example, A(B) means that A is nested within B. A(A) is not allowed because factors inside a nested effect must be distinct.
- Multiple nesting is allowed. For example, A(B(C)) means that B is nested within C, and A is nested within B(C). When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, A(B)(C) is not valid.
- Nesting within an interaction effect is valid. For example, A(B\*C) means that A is nested within B\*C.
- Interactions among nested effects are allowed. The correct syntax is the interaction followed by the common nested effect inside the parentheses. For example, interaction between A and B within levels of C should be specified as A\*B(C) instead of A(C)\*B(C).
- To include a covariate term in the design, enter the name of the covariate.
- Covariates can be connected, but not nested, through the \* operator or using the keyword BY to form another covariate effect. Interactions among covariates such as X1\*X1 and X1\*X2 are valid, but X1(X2) is not.
- Factor and covariate effects can be connected in various ways except that no effects can be nested within a covariate effect. Suppose A and B are factors and X1 and X2 are covariates, examples of valid combinations of factor and covariate effects are A\*X1, A\*B\*X1, X1(A), X1(A\*B), X1\*A(B), X1\*X2(A\*B), and A\*B\*X1\*X2.

## ***CUSTOM Subcommand***

The `CUSTOM` subcommand defines custom hypothesis tests by specifying the **L** matrix (contrast coefficients matrix) and the **K** matrix (contrast results matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ . The vector **B** is the parameter vector in the linear model.

- Multiple `CUSTOM` subcommands are allowed. Each subcommand is treated independently.
- An optional label may be specified by using the `LABEL` keyword. The label is a string with a maximum length of 255 characters. Only one label can be specified.
- Either the `LMATRIX` or `KMATRIX` keyword, or both, must be specified.

**LMATRIX** *Contrast coefficients matrix.* This matrix specifies coefficients of contrasts, which can be used for studying the effects in the model. An **L** matrix can be specified by using the `LMATRIX` keyword.

**KMATRIX** *Contrast results matrix.* This matrix specifies the results of the linear hypothesis. A **K** matrix can be specified by using the `KMATRIX` keyword.

- The number of rows in the **L** and **K** matrices must be equal.
- A custom hypothesis test can be formed by specifying an **L** or **K** matrix, or both. If only one matrix is specified, the unspecified matrix uses the defaults described below.
- If `KMATRIX` is specified but `LMATRIX` is not specified, the **L** matrix is assumed to be the row vector corresponding to the intercept in the estimable function, provided that `INCLUDE = YES` or `ONLY` is specified on the `INTERCEPT` subcommand. In this case, the **K** matrix can be only a scalar matrix.
- The default **K** matrix is a zero matrix—that is,  $\mathbf{LB} = \mathbf{0}$  is assumed.
- There are three general formats that can be used on the `LMATRIX` keyword: (1) Specify a coefficient value for the intercept, followed optionally by an effect name and a list of real numbers. (2) Specify an effect name and a list of real numbers. (3) Specify keyword `ALL` and a list of real numbers. In all three formats, there can be multiple effect names (or instances of the keyword `ALL`) and number lists.
- Only valid effects in the default model or on the `MODEL` subcommand can be specified on the `LMATRIX` keyword.
- The length of the list of real numbers on the `LMATRIX` keyword must be equal to the number of parameters (including the redundant parameters) corresponding to the specified effect. For example, if the effect `A*B` takes up six columns in the design matrix, the list after `A*B` must contain exactly six numbers.
- When `ALL` is specified, the length of the list that follows `ALL` must be equal to the total number of parameters (including the redundant parameters) in the model.
- Effects that are in the model but not specified on the `LMATRIX` keyword are assumed to have entries of 0 in the corresponding columns of the **L** matrix.
- When an **L** matrix is being defined, a number can be specified as a fraction with a positive denominator. For example, `1/3` and `-1/3` are valid, but `1/-3` is invalid.
- A semicolon (;) indicates the end of a row in the **L** matrix.

- The format for the `KMATRIX` keyword is one or more real numbers. If more than one number is specified, then separate adjacent numbers using a semicolon (;). Each semicolon indicates the end of a row in the **K** matrix. Each number is the hypothesized value for a contrast, which is defined by a row in the **L** matrix.
- For the `KMATRIX` keyword to be valid, either the `LMATRIX` keyword, or `INCLUDE = YES` on the `INTERCEPT` subcommand, must be specified.

### Example

Suppose that factors A and B each have three levels.

```
CSCOXREG t BY a b
/VARIABLES STATUS=death(1)
/PLAN FILE='c:\survey\myfile.csplan'
/MODEL a b a*b
/CUSTOM LABEL = "Effect A"
  LMATRIX = a 1 0 -1
            a*b 1/3 1/3 1/3
                0 0 0
                -1/3 -1/3 -1/3;
            a 0 1 -1
            a*b 0 0 0
                1/3 1/3 1/3
                -1/3 -1/3 -1/3.
```

- The preceding syntax specifies a test of effect A.
- Because there are three levels in effect A, two independent contrasts can be formed at most; thus, there are two rows in the **L** matrix separated by a semicolon (;).
- There are three levels each in effects A and B; thus, the interaction effect A\*B takes nine columns in the design matrix.
- The first row in the **L** matrix tests the difference between levels 1 and 3 of effect A; the second row tests the difference between levels 2 and 3 of effect A.
- The `KMATRIX` keyword is not specified, so the null hypothesis value for both tests is 0.

### Example

Suppose that factor A has three levels.

```
CSCOXREG t BY a
/VARIABLES STATUS=death(1)
/PLAN FILE='c:\survey\myfile.csplan'
/MODEL a
/CUSTOM LABEL = 'Effect A'
  LMATRIX = a 1 0 -1;
            a 0 1 -1
  KMATRIX = 1; 1.
```

- The syntax specifies a model with a main effect for factor A and a custom hypothesis test of effect A.
- The equivalent `LMATRIX` keyword using the `ALL` option follows.

```
LMATRIX = ALL 1 0 -1;
```

ALL 0 1 -1

- The `KMATRIX` keyword is specified and the hypothesis that the difference between levels 1 and 3 and levels 2 and 3 of effect A are both equal to 1 is tested.

## **CRITERIA Subcommand**

The `CRITERIA` subcommand controls the iterative algorithm used for estimation, specifies the numerical tolerance for checking for singularity, and specifies the ties breaking method used in estimating regression parameters.

**CILEVEL = number**

*Confidence interval level for coefficient estimates, exponentiated coefficient estimates, survival function estimates, and cumulative hazard function estimates.* Specify a value greater than or equal to 0 and less than 100. The default value is 95.

**DF = number**

*Sampling design degrees of freedom to use in computing p values for all test statistics.* Specify a positive number. The default value is the difference between the number of primary sampling units and the number of strata in the first stage of sampling.

**LCONVERGE = [number RELATIVE|ABSOLUTE]**

*Log-likelihood function convergence criterion.* Convergence is assumed if the relative or absolute change in the log-likelihood function is less than the given value. This criterion is not used if the value is 0.

Specify square brackets containing a non-negative number followed optionally by keyword `RELATIVE` or `ABSOLUTE`, which indicates the type of change. The default value is 0; the default type is `RELATIVE`.

**MXITER = integer**

*Maximum number of iterations.* Specify a non-negative integer. The default value is 100.

**MXSTEP = integer**

*Maximum step-halving allowed.* Specify a positive integer. The default value is 5.

**PCONVERGE = [number RELATIVE|ABSOLUTE]**

*Parameter estimates convergence criterion.* Convergence is assumed if the relative or absolute change in the parameter estimates is less than the given value. This criterion is not used if the value is 0.

Specify square brackets containing a non-negative number followed optionally by keyword `RELATIVE` or `ABSOLUTE`, which indicates the type of change. The default value is  $10^{-6}$ ; the default type is `RELATIVE`.

**SINGULAR = number**

*Tolerance value used to test for singularity.* Specify a positive value. The default value is  $10^{-12}$ .

**TIES = EFRON|BRESLOW**

*Tie breaking method in estimating parameters.* The default Efron method is specified by the keyword `EFRON`; the default Breslow method is specified by the keyword `BRESLOW`.

## **STATISTICS Subcommand**

The `STATISTICS` subcommand requests various statistics associated with the parameter estimates.

There are no default keywords on the `STATISTICS` subcommand. If this subcommand is not specified, then none of the statistics listed below are displayed.

<b>PARAMETER</b>	<i>Parameter estimates.</i>
<b>EXP</b>	<i>The exponentiated parameter estimates.</i>
<b>SE</b>	<i>Standard error for each parameter estimate.</i>
<b>TTEST</b>	<i>t test for each parameter estimate.</i>
<b>CINTERVAL</b>	<i>Confidence interval for each parameter estimate and/or exponentiated parameter estimate.</i>
<b>DEFF</b>	<i>Design effect for each parameter estimate.</i>
<b>DEFFSQRT</b>	<i>Square root of design effect for each parameter estimate.</i>

## **TEST Subcommand**

The `TEST` subcommand specifies the type of test statistic and the method of adjusting the significance level to be used for hypothesis tests requested on the `MODEL`, `CUSTOM`, and `PRINT` subcommands.

### **TYPE Keyword**

The `TYPE` keyword indicates the type of test statistic.

<b>F</b>	<i>Wald F test.</i> This is the default test statistic if the <code>TYPE</code> keyword is not specified.
<b>ADJF</b>	<i>Adjusted Wald F test.</i>
<b>CHISQUARE</b>	<i>Wald chi-square test.</i>
<b>ADJCHISQUARE</b>	<i>Adjusted Wald chi-square test.</i>

### **PADJUST Keyword**

The `PADJUST` keyword indicates the method of adjusting the significance level.

<b>LSD</b>	<i>Least significant difference.</i> This method does not control the overall probability of rejecting the hypotheses that some linear contrasts are different from the null hypothesis value(s). This is the default.
<b>BONFERRONI</b>	<i>Bonferroni.</i> This method adjusts the observed significance level for the fact that multiple contrasts are being tested.
<b>SEQBONFERRONI</b>	<i>Sequential Bonferroni.</i> This is a sequentially step-down rejective Bonferroni procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.
<b>SIDAK</b>	<i>Sidak.</i> This method provides tighter bounds than the Bonferroni approach.
<b>SEQSIDAK</b>	<i>Sequential Sidak.</i> This is a sequentially step-down rejective Sidak procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.

## **TESTASSUMPTIONS Subcommand**

The TESTASSUMPTIONS subcommand produces tests of the proportional hazards and covariate form model assumptions. You can request various statistics associated with the alternative models.

### **PROPHAZARD Keyword**

The PROPHAZARD keyword produces a test for proportional hazards assumption. The time function used in testing for proportional hazards is specified in parentheses. Specify one of the following options.

<b>KM</b>	<i>Kaplan-Meier estimation of survival function.</i> This is the default.
<b>IDENTITY</b>	<i>Identity function of time.</i>
<b>LOG</b>	<i>Log function of time.</i>
<b>RANK</b>	<i>Rank of death time.</i>

### **PARAMETER Keyword**

The PARAMETER keyword displays the parameter estimates of the alternative model. The alternative model is estimated using the same convergence criteria as the original model. Both parameters and their standard errors are estimated.

### **COVB Keyword**

The COVB keyword displays the covariance matrix for the alternative model parameters

## **DOMAIN Subcommand**

The DOMAIN subcommand specifies the subpopulation for which the analysis is to be performed.

- The keyword VARIABLE, followed by an equals sign, a variable, and a value in parentheses are required. Put the value inside a pair of quotes if the value is formatted (such as date or currency) or if the factor is of string type.
- The subpopulation is defined by all cases having the given value on the specified variable.
- Analyses are performed only for the specified subpopulation.
- For example, DOMAIN VARIABLE = myvar (1) defines the subpopulation by all cases for which variable MYVAR has value 1.
- The specified variable may be numeric or string and must exist at the time the CSCOXREG procedure is invoked.
- Stratification or cluster variables may be specified, but no other plan file variables are allowed on the DOMAIN subcommand.
- Analysis variables may not be specified on the DOMAIN subcommand.

## **MISSING Subcommand**

The MISSING subcommand specifies how missing values are handled.

- In general, cases must have valid data for all design variables as well as for the dependent variable and any covariates. Cases with invalid data for any of these variables are excluded from the analysis.
- There is one important exception to the preceding rule. This exception applies when an inclusion probability or population size variable is defined in an analysis plan file. Within a stratum at a given stage, if the inclusion probability or population size values are unequal across cases or missing for a case, then the first valid value found within that stratum is used as the value for the stratum. If strata are not defined, then the first valid value found in the sample is used. If the inclusion probability or population size values are missing for all cases within a stratum (or within the sample if strata are not defined) at a given stage, then an error message is issued.
- The CLASSMISSING keyword specifies whether user-missing values are treated as valid. This specification is applied to categorical design variables (that is, strata, cluster, and subpopulation variables), the dependent variable, and any factors.

**EXCLUDE**        *Exclude user-missing values among the strata, cluster, subpopulation, dependent variable, and factor variables. This is the default.*

**INCLUDE**        *Include user-missing values among the strata, cluster, subpopulation, dependent variable, and factor variables. Treat user-missing values for these variables as valid data.*

## ***SURVIVALMETHOD Subcommand***

The SURVIVALMETHOD subcommand controls the methods for estimating baseline functions and the confidence interval of the survival function.

### ***BASELINE Keyword***

The BASELINE keyword controls the method for estimating baseline functions. Specify one of the following options.

**EFRON**            *Efron method. Default if EFRON is chosen in TIES.*

**BRESLOW**        *Breslow method. Default if BRESLOW is chosen in TIES.*

**PRO-  
DUCTLIMIT**      *Product limit method.*



**CI Keyword**

The CI keyword controls the method for estimating the confidence interval of the survival function. Specify one of the following options.

<b>ORIGINAL</b>	<i>Based on original scale.</i> Calculate the confidence interval for the survival function directly.
<b>LOG</b>	<i>Based on log scale.</i> Calculate the confidence interval for $\ln(\text{survival})$ first, then back transform to get the confidence interval for the survival function.
<b>LML</b>	<i>Based on log-log scale.</i> Calculate the confidence interval for $\ln(-\ln(\text{survival}))$ first, then back transform to get the confidence interval for the survival function.

**PRINT Subcommand**

The PRINT subcommand is used to display optional output.

- If the PRINT subcommand is not specified, then the default output includes sample information, variable and factor information, and model summary statistics.
- If the PRINT subcommand is specified, then CSCOXREG displays output only for those keywords that are specified.

<b>SAMPLEINFO</b>	<i>Sample information table.</i> Displays summary information about the sample, including the unweighted count, the event and censoring counts, and the population size. This is default output if the PRINT subcommand is not specified.
<b>EVENTINFO</b>	<i>Event and censoring information.</i> Displays event and censoring information for each baseline stratum. This is the default output if the PRINT subcommand is not specified.
<b>RISKINFO</b>	<i>Risk and event information.</i> Displays number of events and number at risk for each event time in each baseline stratum.
<b>HISTORY(n)</b>	<i>Iteration history.</i> Displays coefficient estimates and statistics at every $n$ th iteration beginning with the 0th iteration (the initial estimates). The default is to print every iteration ( $n = 1$ ). The last iteration is always printed if HISTORY is specified, regardless of the value of $n$ .
<b>GEF</b>	<i>General estimable function table.</i>
<b>LMATRIX</b>	<i>Set of contrast coefficients (L) matrices.</i> These are the Type III contrast matrices used in testing model effects.
<b>COVB</b>	<i>Covariance matrix for model parameters.</i>
<b>CORB</b>	<i>Correlation matrix for model parameters.</i>
<b>BASELINE</b>	<i>Baseline functions.</i> Displays the baseline survival function, baseline cumulative hazards function and their standard errors. If time-dependent covariates defined by TIME PROGRAM are included in the model, no baseline functions are produced.
<b>NONE</b>	<i>No PRINT output.</i> None of the PRINT subcommand default output is displayed. However, if NONE is specified with one or more other keywords, then the other keywords override NONE.

## **SAVE Subcommand**

The `SAVE` subcommand writes optional model variables to the working data file.

- Specify one or more temporary variables, each followed by an optional new name in parentheses.
- The optional names must be valid variable names.
- If new names are not specified, `CSCOXREG` uses the default names.
- The optional variable name must be unique. If the default name is used and it conflicts with existing variable names, then a suffix is added to the default name to make it unique.
- If a subpopulation is defined on the `DOMAIN` subcommand, then `SAVE` applies only to cases within the subpopulation.
- Aggregated residuals are residuals aggregated over records with the same ID value. If ID is not specified, aggregated residuals are not available and a warning is issued if they are requested. The aggregated residual for a subject is written in the last case (or first case if it is easier) of that subject.
- If time-dependent covariates defined by `TIME PROGRAM` are included in the model, the following options are not available: `MARTINGALE`, `DEVIANCE`, `COXSNELL`, `AGGMARTINGALE`, `AGGDEVIANCE`, `AGGCOXSNELL`, `SURVIVAL`, `LCL_SURVIVAL`, `UCL_SURVIVAL`, `CUMHAZARD`, `LCL_CUMHAZARD`, and `UCL_CUMHAZARD`. A warning is issued if they are requested.
- In situations when rootname is needed, the rootname can be followed by a colon and a positive integer giving the maximum number of variables with the same rootname to be saved. The first  $n$  variables are saved. The default  $n$  is 25. To specify  $n$  without a rootname, enter a colon before the number.

### **SCHOENFELD(rootname:n)**

*Schoenfeld residual.* A separate variable is saved for each nonredundant parameter and calculated only for noncensored observations. The default variable name is *Resid\_Schoenfeld*.

### **MARTINGALE(varname)**

*Martingale residual.* The default variable name is *Resid\_Martingale*.

### **DEVIANCE(varname)**

*Deviance residual.* The default variable name is *Resid\_Deviance*.

### **COXSNELL(varname)**

*Cox\_Snell residual.* The default variable name is *Resid\_CoxSnell*.

### **SCORE(rootname:n)**

*Score residual.* A separate variable is saved for each nonredundant parameter. The default variable name is *Resid\_Score*.

### **DFBETA(rootname:n)**

*DFBETA.* A separate variable is saved for each nonredundant parameter. The default variable name is *Resid\_DFBETA*.

### **AGGMARTINGALE(varname)**

*Aggregated Martingale residual.* The default variable name is *AggResid\_Martingale*.

### **AGGDEVIANCE(varname)**

*Aggregated Deviance residual.* The default variable name is *AggResid\_Deviance*.

### **AGGCOXSNELL(varname)**

*Aggregated CoxSnell residual.* The default variable name is *AggResid\_CoxSnell*.

**AGGSCORE(rootname:n)**

*Aggregated Score residual.* A separate variable is saved for each nonredundant parameter. The default variable name is *AggResid\_Score*.

**AGGDFBETA(rootname:n)**

*Aggregated DFBETA.* A separate variable is saved for each nonredundant parameter. The default variable name is *AggResid\_DFBETA*.

**XBETA(varname)**

*Linear combination of reference value corrected predictors times regression coefficients.* The default variable name is *XBETA*.

**SURVIVAL(varname)**

*Survival function.* For one-time input data, it is the survival function at the observed time and predictor pattern for each record. For two-time input data, it is the survival function at *endtime* assuming that the predictor is fixed. The default variable name is *Survival*.

**LCL\_SURVIVAL(varname)**

*Lower confidence level of survival function.* The default variable name is *LCL\_Survival*.

**UCL\_SURVIVAL(varname)**

*Upper confidence level of survival function.* The default variable name is *UCL\_Survival*.

**CUMHAZARD(varname)**

*Cumulative hazards function.* The default variable name is *CumHazard*.

**LCL\_CUMHAZARD(varname)**

*Lower confidence level of cumulative hazards function.* The default variable name is *LCL\_CumHazard*.

**UCL\_CUMHAZARD(varname)**

*Upper confidence level of cumulative hazards function.* The default variable name is *SECumHazard*.

## ***PLOT Subcommand***

You can request specific plots to be produced with the `PLOT` subcommand. Each requested plot is produced once for each pattern specified on the `PATTERN` subcommand.

- The set of plots requested is displayed for the functions at the mean of the covariates and at each combination of covariate values specified on `PATTERN`.
- Lines on a plot are connected as step functions.

**SURVIVAL**     *Plot the survival function.*

**HAZARD**     *Plot the cumulative hazard function.*

**LML**         *Plot the log-minus-log-of-survival function.*

**OMS**         *Plot the one-minus-survival function.*

**CI = NO | YES**     *Plot confidence intervals along with the specified functions.* NO is the default.

## ***PATTERN Subcommand***

PATTERN specifies the pattern of predictor values to be used for requested plots on the PLOT subcommand and the exported survival file on the OUTFILE subcommand. PATTERN cannot be used when time-dependent predictors calculated by TIME PROGRAM are included in the model.

- A value must be specified for each variable specified on PATTERN.
- Covariates that are included in the model but not named on PATTERN are evaluated at their means.
- Factors that are included in the model but not named on PATTERN are evaluated at the reference category.
- You can request separate lines for each category of a factor that is in the model. Specify the name of the categorical variable after the keyword BY. The BY variable must be a categorical variable. You cannot specify a value for the BY variable.
- Multiple PATTERN subcommands can be specified. CSCOXREG produces a set of requested plots for each specified pattern.
- Piecewise constant predictor paths are also allowed. The path is specified by `endtime(valuelist) varname(valuelist) varname(value)...`. If `varname(valuelist)` is used, the length of `valuelist` must be the same as that for `endtime`. The `varname(value)` means that the value of the variable is constant over time.

### ***Example***

```
CSCOXREG t by A with x1 x2
/VARIABLES STATUS=dead(1)
/PLAN FILE='c:\survey\myfile.csplan'
/PATTERN x1(0.1) x2(3) A(3)
/PLOT SURVIVAL.
```

- Predictor pattern  $x1 = 0.1$ ,  $x2 = 3$ ,  $A = 3$  is specified by PATTERN.
- The survival function is plotted for the specified pattern.

### ***Example: Piecewise constant predictor path***

```
CSCOXREG t1 t2 by A with x1 x2
/VARIABLES STATUS=dead(1)
/PLAN FILE='c:\survey\myfile.csplan'.
/PATTERN t2(10 20 30 50) x1(1 1.2 1.7 1.9) x2(3) BY A
/OUTFILE SURVIVAL='surv.sav'.
```

- Two time variables are specified on the CSCOXREG variable list.
- PATTERN defines the following predictor paths for  $x1$  and  $x2$ .

starttime	endtime	x1	x2
0	10	1.0	3
10	20	1.2	3
20	30	1.7	3
30	50	1.9	3

- `PATTERN`, through `BY A`, also specifies that each category of factor A is considered separately. Combining different categories of A with the paths for `x1` and `x2`, the total number of paths considered here actually equals the number of categories of A.
- The survival table for the specified predictor paths are calculated and written to the file `surv.sav`.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand saves an external IBM® SPSS® Statistics data file or previously declared dataset containing the parameter covariance or correlation matrix with parameter estimates, standard errors, significance values, and sampling design degrees of freedom. It also saves the parameter estimates and the parameter covariance matrix in XML format.

- At least one keyword and a filename are required.
- The `COVB` and `CORB` keywords are mutually exclusive, as are the `MODEL` and `PARAMETER` keywords.
- The filename must be specified in full. `CSCOXREG` does not supply an extension.

**COVB = 'savfile']'dataset'**

*Writes the parameter covariance matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**CORB = 'savfile']'dataset'**

*Writes the parameter correlation matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**MODEL = 'file'**

*Writes all information needed to predict the survival function, including the parameter estimates and baseline survival function, to a PMML file.*

**SURVIVAL= 'savfile']'dataset'**

*Writes survival table to an external SPSS Statistics data file or previously declared dataset. The file contains the survival function, standard error of the survival function, upper and lower bounds of the confidence interval of the survival function, and the cumulative hazards function for each failure or event time evaluated at the baseline and at the covariate patterns specified on `PATTERN`. If time-dependent covariates are included in the model, no file is written.*

# CSDESCRIPTIVES

CSDESCRIPTIVES is available in the Complex Samples option.

```
CSDESCRIPTIVES
/PLAN FILE = file
[/JOINTPROB FILE = file]
[/SUMMARY VARIABLES = varlist]
[/MEAN [TTEST = {value    }]
      {valuelist}]
[/SUM [TTEST = {value    }]
      {valuelist}]
[/RATIO NUMERATOR = varlist DENOMINATOR = varlist
      [TTEST = {value    }]]
      {valuelist}]
[/RATIO...]

[/STATISTICS [COUNT] [POPSIZE] [SE] [CV] [DEFF] [DEFFSQRT]
             [CIN [({95** })]]]
             {value}]
[/SUBPOP TABLE = varname [BY varname [BY ...]] [DISPLAY = {LAYERED }]]
             {SEPARATE}]
[/MISSING [SCOPE = {ANALYSIS}] [CLASSMISSING = {EXCLUDE}]]
          {LISTWISE}           {INCLUDE}]
```

**\*\*Default if the subcommand is omitted.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
CSDESCRIPTIVES
  /PLAN FILE = '/survey/myfile.xml'
  /SUMMARY VARIABLES = y1 y2
  /MEAN.
```

## Overview

CSDESCRIPTIVES estimates means, sums, and ratios, and computes their standard errors, design effects, confidence intervals, and hypothesis tests, for samples that are drawn by complex sampling methods. The procedure estimates variances by taking into account the sample design that is used to select the sample, including equal probability and probability proportional to size (PPS) methods, and with replacement (WR) and without replacement (WOR) sampling procedures. Optionally, CSDESCRIPTIVES performs analyses for subpopulations.

**Basic Specification**

- The basic specification is a `PLAN` subcommand and the name of a complex sample analysis plan file (which may be generated by the `CSPLAN` procedure) and a `MEAN`, `SUM`, or `RATIO` subcommand. If a `MEAN` or `SUM` subcommand is specified, a `SUMMARY` subcommand must also be present.
- The basic specification displays the overall population size estimate. Additional subcommands must be used for other results.

**Operations**

- `CSDESCRIPTIVES` computes estimates for sampling designs that are supported by the `CSPLAN` and `CSSELECT` procedures.
- The input dataset must contain the variables to be analyzed and variables that are related to the sampling design.
- The complex sample analysis plan file provides an analysis plan based on the sampling design.
- The default output for each requested mean, sum, or ratio is the estimate and its standard error.
- `WEIGHT` and `SPLIT FILE` settings are ignored by the `CSDESCRIPTIVES` procedure.

**Syntax Rules**

- The `PLAN` subcommand is required. In addition, the `SUMMARY` subcommand and the `MEAN` or `SUM` subcommand must be specified, or the `RATIO` subcommand must be specified. All other subcommands are optional.
- Multiple instances of the `RATIO` subcommand are allowed—each instance is treated independently. All other subcommands may be specified only once.
- Subcommands can be specified in any order.
- All subcommand names and keywords must be spelled in full.
- Equals signs (=) that are shown in the syntax chart are required.
- The `MEAN` and `SUM` subcommands can be specified without further keywords, but no other subcommands may be empty.

**Examples****Example: Univariate Descriptive Statistics**

```
* Complex Samples Descriptives.
CSDESCRIPTIVES
  /PLAN FILE = 'samplesDirectory\nhis2000_subset.csaplan'
  /JOINTPROB FILE = 'samplesDirectory\nhis2000_subset.sav'
  /SUMMARY VARIABLES =VIGFREQW MODFREQW STRFREQW
  /SUBPOP TABLE = age_cat DISPLAY=LAYERED
  /MEAN
  /STATISTICS SE CIN (95)
  /MISSING SCOPE = ANALYSIS CLASSMISSING = EXCLUDE.
```

- The procedure computes estimates based on the complex sample analysis plan that is given in *nhis2000\_subset.csaplan*.

- The procedure estimates the mean, its standard error, and 95% confidence interval for variables *vigfreqw*, *modfreqw*, and *strfreqw*.
- In addition, these statistics are computed for the variables by values of *age\_cat*. The results for subpopulations are displayed in a single table.
- Other subcommands and keywords are set to their default values.

### **Example: Ratio Statistics**

```
* Complex Samples Ratios.
CSDESCRIPTIVES
  /PLAN FILE = 'samplesDirectory\property_assess.csplan'
  /RATIO NUMERATOR = currval DENOMINATOR = lastval TTEST = 1.3
  /STATISTICS SE COUNT POPSIZE CIN (95)
  /SUBPOP TABLE = county DISPLAY=LAYERED
  /MISSING SCOPE = ANALYSIS CLASSMISSING = EXCLUDE.
```

- The procedure computes estimates based on the complex sample analysis plan that is given in *property\_assess.csplan*.
- The ratio estimate for *currval/lastval*, its standard error, 95% confidence interval, observed count of cases used in the computations, and estimated population size are displayed.
- A *t* test of the ratio is performed against a hypothesized value of 1.3.
- In addition, these statistics are computed for the variables by values of *county*. The results for subpopulations are displayed in a single table.
- Other subcommands and keywords are set to their default values.

## **PLAN Subcommand**

PLAN specifies the name of an XML file containing analysis design specifications. This file is written by the CSPLAN procedure.

- The PLAN subcommand is required.

**FILE**                                      *Specifies the name of an external file.*

## **JOINTPROB Subcommand**

JOINTPROB is used to specify the file or dataset containing the first-stage joint inclusion probabilities for UNEQUAL\_WOR estimation. The CSSELECT procedure writes this file in the same location and with the same name (but different extension) as the plan file. When UNEQUAL\_WOR estimation is specified, the CSDESCRIPTIVES procedure will use the default location and name of the file unless the JOINTPROB subcommand is used to override them.

**FILE**                                      *Specifies the name of the file or dataset containing the joint inclusion probabilities.*

## **SUMMARY Subcommand**

SUMMARY specifies the analysis variables that are used by the MEAN and SUM subcommands.



- A variable list is required only if means or sums are to be estimated. If only ratios are to be estimated (that is, if the `RATIO` subcommand is specified but the `MEAN` and `SUM` subcommands are not specified), the `SUMMARY` subcommand is ignored.
- All specified variables must be numeric.
- All specified variables must be unique.
- Plan file and subpopulation variables may not be specified on the `SUMMARY` subcommand.

**VARIABLES**                    *Specifies the variables used by the `MEAN` and `SUM` subcommands.*

## ***MEAN Subcommand***

`MEAN` is used to request that means be estimated for variables that are specified on the `SUMMARY` subcommand.

The `TTEST` keyword requests  $t$  tests of the population means(s) and gives the null hypothesis value(s). If subpopulations are defined on the `SUBPOP` subcommand, null hypothesis values are used in the test(s) for each subpopulation, as well as for the entire population.

**value**                            The null hypothesis is that the population mean equals the specified value for all  $t$  tests.

**valuelist**                      This list gives the null hypothesis value of the population mean for each variable on the `SUMMARY` subcommand. The number and order of values must correspond to the variables on the `SUMMARY` subcommand.

- Commas or spaces must be used to separate the values.

## ***SUM Subcommand***

`SUM` is used to request that sums be estimated for variables specified on the `SUMMARY` subcommand.

The `TTEST` keyword requests  $t$  tests of the population sum(s) and gives the null hypothesis value(s). If subpopulations are defined on the `SUBPOP` subcommand, then null hypothesis values are used in the test(s) for each subpopulation as well as for the entire population.

**value**                            The null hypothesis is that the population sum equals the specified value for all  $t$  tests.

**valuelist**                      This list gives the null hypothesis value of the population sum for each variable on the `SUMMARY` subcommand. The number and order of values must correspond to the variables on the `SUMMARY` subcommand.

- Commas or spaces must be used to separate the values.

## ***RATIO Subcommand***

`RATIO` specifies ratios of variables to be estimated.

- Ratios are defined by crossing variables on the `NUMERATOR` keyword with variables on the `DENOMINATOR` keyword, with `DENOMINATOR` variables looping fastest, irrespective of the order of the keywords. For example, `/RATIO NUMERATOR = N1 N2 DENOMINATOR = D1 D2` yields the following ordered list of ratios:  $N1/D1$ ,  $N1/D2$ ,  $N2/D1$ ,  $N2/D2$ .

- Multiple `RATIO` subcommands are allowed. Each subcommand is treated independently.
- Variables that are specified on the `RATIO` subcommand do not need to be specified on the `SUMMARY` subcommand.
- All specified variables must be numeric.
- Within each variable list, all specified variables must be unique.
- Plan file and subpopulation variables may not be specified on the `RATIO` subcommand.

The `TTEST` keyword requests  $t$  tests of the population ratio(s) and gives the null hypothesis value(s). If subpopulations are defined on the `SUBPOP` subcommand, then null hypothesis values are used in the test(s) for each subpopulation as well as for the entire population.

<b>value</b>	The null hypothesis is that the population ratio equals the specified value for all $t$ tests.
<b>valuelist</b>	This list gives the null hypothesis value of the population ratio for each ratio specified on the <code>RATIO</code> subcommand. The number and order of values must correspond to the ratios defined on the <code>RATIO</code> subcommand.

- Commas or spaces must be used to separate the values.

## ***STATISTICS Subcommand***

`STATISTICS` requests various statistics that are associated with the mean, sum, or ratio estimates. If the `STATISTICS` subcommand is not specified, the standard error is computed for any displayed estimates. If the `STATISTICS` subcommand is specified, only statistics that are requested are computed.

<b>COUNT</b>	<i>The number of valid observations in the dataset for each mean, sum, or ratio estimate.</i>
<b>POPSIZE</b>	<i>The population size for each mean, sum, or ratio estimate.</i>
<b>SE</b>	<i>The standard error for each mean, sum, or ratio estimate. This output is default output if the <code>STATISTICS</code> subcommand is not specified.</i>
<b>CV</b>	<i>Coefficient of variation.</i>
<b>DEFF</b>	<i>Design effect.</i>
<b>DEFFSQRT</b>	<i>Square root of the design effect.</i>
<b>CIN [(value)]</b>	<i>Confidence interval. If the <code>CIN</code> keyword is specified alone, the default 95% confidence interval is computed. Optionally, <code>CIN</code> may be followed by a value in parentheses, where <math>0 \leq \text{value} &lt; 100</math>.</i>

## ***SUBPOP Subcommand***

`SUBPOP` specifies subpopulations for which analyses are to be performed.

- The set of subpopulations is defined by specifying a single categorical variable or specifying two or more categorical variables, separated by the `BY` keyword, whose values are crossed.
- For example, `/SUBPOP TABLE = A` defines subpopulations based on the levels of variable  $A$ .
- For example, `/SUBPOP TABLE = A BY B` defines subpopulations based on crossing the levels of variables  $A$  and  $B$ .

- A maximum of 17 variables may be specified.
- Numeric or string variables may be specified.
- All specified variables must be unique.
- Stratification or cluster variables may be specified, but no other plan file variables are allowed on the SUBPOP subcommand.
- Analysis variables may not be specified on the SUBPOP subcommand.
- The BY keyword is used to separate variables.

The DISPLAY keyword specifies the layout of results for subpopulations.

**LAYERED**            *Results for all subpopulations are displayed in the same table. This is the default.*  
**SEPARATE**         *Results for different subpopulations are displayed in different tables.*

## **MISSING Subcommand**

MISSING specifies how missing values are handled.

- All design variables must have valid data. Cases with invalid data for any design variable are deleted from the analysis.

The SCOPE keyword specifies which cases are used in the analyses. This specification is applied to analysis variables but not design variables.

**ANALYSIS**            *Each statistic is based on all valid data for the analysis variable(s) used in computing the statistic. Ratios are computed by using all cases with valid data for both of the specified variables. Statistics for different variables may be based on different sample sizes. This setting is the default.*

**LISTWISE**            *Only cases with valid data for all analysis variables are used in computing any statistics. Statistics for different variables are always based on the same sample size.*

The CLASSMISSING keyword specifies whether user-missing values are treated as valid. This specification is applied only to categorical design variables (strata, cluster, and subpopulation variables).

**EXCLUDE**            *Exclude user-missing values among the strata, cluster, and subpopulation variables. This setting is the default.*

**INCLUDE**            *Include user-missing values among the strata, cluster, and subpopulation variables. Treat user-missing values for these variables as valid data.*

# CSGLM

CSGLM is available in the Complex Samples option.

*Note:* Square brackets that are used in the CSGLM syntax chart are required parts of the syntax and are not used to indicate optional elements. Equals signs (=) that are used in the syntax chart are required elements. All subcommands, save the PLAN subcommand, are optional.

CSGLM dependent var BY factor list WITH covariate list

```
/PLAN FILE = file
/JOINTPROB FILE = file
/MODEL effect list
/INTERCEPT INCLUDE = {YES**} SHOW = {YES**}
                        {NO   }          {NO   }
                        {ONLY }
/CUSTOM LABEL = "label"
      LMATRIX = {number effect list effect list ...; ...}
                {number effect list effect list ...   }
                {effect list effect list ...; ...     }
                {effect list effect list ...         }
                {ALL list; ALL ...                   }
                {ALL list                             }
      KMATRIX = {number; number; ...}
                {number                             }
/CUSTOM ...
/EMMEANS TABLES = {factor           }
                  {factor*factor...}
      OTHER = [varname (value) varname (value) ...]
      COMPARE = factor
      CONTRAST = {SIMPLE** (value)      }
                 {DEVIATION (value)    }
                 {DIFFERENCE           }
                 {HELMERT               }
                 {REPEATED              }
                 {POLYNOMIAL (number list)}
/EMMEANS ...
/CRITERIA CILEVEL = {95** } DF = n SINGULAR = {1E-12**}
                  {value}                   {value  }
/STATISTICS PARAMETER SE TTEST CINTERVAL DEFF DEFFSQRT
/TEST TYPE = {F**           } PADJUST = {LSD**           }
             {ADJF          }           {BONFERRONI      }
             {CHISQUARE     }           {SEQBONFERRONI}
             {ADJCHISQUARE}           {SIDAK          }
             {ADJCHISQUARE}           {SEQSIDAK       }
/DOMAIN VARIABLE = varname (value)
/MISSING CLASSMISSING = {EXCLUDE**}
                       {INCLUDE  }
/PRINT SAMPLEINFO** VARIABLEINFO** SUMMARY**
```

```

      GEF LMATRIX COVB CORB NONE
/SAVE PRED(varname) RESID(varname)
/OUTFILE {COVB='savfile' | 'dataset'} {MODEL = 'file'      }
         {CORB='savfile' | 'dataset'} {PARAMETER = 'file' }

```

**\*\*Default if the keyword or subcommand is omitted.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- Command introduced.

### **Example**

```

CSGLM y BY a b c WITH x
      /PLAN FILE='/survey/myfile.csplan'.

```

## **Overview**

CSGLM performs linear regression analysis, as well as analysis of variance and covariance, for samples that are drawn by complex sampling methods. The procedure estimates variances by taking into account the sample design that is used to select the sample, including equal probability and probability proportional to size (PPS) methods, and with replacement (WR) and without replacement (WOR) sampling procedures. Optionally, CSGLM performs analyses for a subpopulation.

### **Basic Specification**

- The basic specification is a variable list (identifying the dependent variable, the factors, if any, and the covariates, if any) and a PLAN subcommand with the name of a complex sample analysis plan file, which may be generated by the CSPLAN procedure.
- The default model includes the intercept term, main effects for any factors, and any covariates.
- The basic specification displays summary information about the sample design,  $R$ -square and root mean square error for the model, regression coefficient estimates and  $t$  tests, and Wald  $F$  tests for all model effects. Additional subcommands must be used for other results.

### **Operations**

- CSGLM computes linear model estimates for sampling designs that are supported by the CSPLAN and CSSELECT procedures.
- The input dataset must contain the variables to be analyzed and variables that are related to the sampling design.
- The complex sample analysis plan file provides an analysis plan based on the sampling design.

- By default, CSGLM uses a model that includes the intercept term, main effects for any factors, and any covariates.
- Other effects, including interaction and nested effects, may be specified by using the MODEL subcommand.
- The default output for the specified model is summary information about the sample design, *R*-square and root mean square error, regression coefficient estimates and *t* tests, and Wald *F* tests for all effects.
- WEIGHT and SPLIT FILE settings are ignored by the CSGLM procedure.

### **Syntax Rules**

- The dependent variable and PLAN subcommand are required. All other variables and subcommands are optional.
- Multiple CUSTOM and EMMEANS subcommands may be specified; each subcommand is treated independently. All other subcommands may be specified only once.
- The EMMEANS subcommand may be specified without options. All other subcommands must be specified with options.
- Each keyword may be specified only once within a subcommand.
- Subcommand names and keywords must be spelled in full.
- Equals signs (=) that are shown in the syntax chart are required.
- Subcommands may be specified in any order.
- The dependent variable and covariates must be numeric, but factors and the subpopulation variable can be numeric or string variables.
- Across the dependent, factor, and covariate variable lists, a variable may be specified only once.
- Plan file and subpopulation variables may not be specified on the variable list.
- Minimum syntax is a dependent variable and the PLAN subcommand. This specification fits an intercept-only model.

### **Limitations**

- WEIGHT and SPLIT FILE settings are ignored with a warning by the CSGLM procedure.

## **Examples**

```
* Complex Samples General Linear Model.
CSGLM amtspent BY shopfor usecoup
  /PLAN FILE = 'samplesDirectory\grocery.csplan'
  /JOINTPROB FILE = 'samplesDirectory\grocery.sav'
  /MODEL shopfor usecoup shopfor*usecoup
  /INTERCEPT INCLUDE=YES SHOW=YES
  /STATISTICS PARAMETER SE CINTERVAL DEFF
  /PRINT SUMMARY VARIABLEINFO SAMPLEINFO
  /TEST TYPE=F PADJUST=LSD
  /EMMEANS TABLES=shopfor COMPARE CONTRAST=SIMPLE(3)
  /EMMEANS TABLES=usecoup COMPARE CONTRAST=SIMPLE(1)
  /EMMEANS TABLES=shopfor*usecoup
  /MISSING CLASSMISSING=EXCLUDE
```

```
/CRITERIA CILEVEL=95.
```

- The procedure fits a general linear model for the dependent variable *amtspent* using *shopfor* and *usecoup* as factors.
- The complex sampling plan is located in *grocery.csplan*; the file containing the joint inclusion probabilities is *grocery.sav*.
- The model specification calls for a full factorial model with intercept.
- Parameter estimates, their standard errors, 95% confidence intervals, and design effects will be displayed.
- Estimated marginal means are computed for each of the model effects. The third level of *shopfor* is specified as the reference category for contrast comparisons; the first level of *usecoup* is specified as the reference category.
- All other options are set to their default values.

## ***CSGLM Variable List***

The variable list specifies the dependent variable, the factors, and the covariates in the model.

- The dependent variable must be the first specification on CSGLM.
- The names of the factors and covariates, if any, follow the dependent variable. Specify any factors following the keyword `BY`. Specify any covariates following the keyword `WITH`.
- The dependent variable and covariates must be numeric, but factors can be numeric or string variables.
- Each variable may be specified only once on the variable list.
- Plan file and subpopulation variables may not be specified on the variable list.

## ***PLAN Subcommand***

The `PLAN` subcommand specifies the name of an XML file containing analysis design specifications. This file is written by the `CSPLAN` procedure.

- The `PLAN` subcommand is required.

**FILE** *Specifies the name of an external file.*

## ***JOINTPROB Subcommand***

The `JOINTPROB` subcommand is used to specify the file or dataset containing the first stage joint inclusion probabilities for `UNEQUAL_WOR` estimation. The `CSSELECT` procedure writes this file in the same location and with the same name (but different extension) as the plan file. When `UNEQUAL_WOR` estimation is specified, the `CSGLM` procedure will use the default location and name of the file unless the `JOINTPROB` subcommand is used to override them.

**FILE** *Specifies the name of the file or dataset containing the joint inclusion probabilities.*

## **MODEL Subcommand**

The `MODEL` subcommand is used to specify the effects to be included in the model. Use the `INTERCEPT` subcommand to control whether the intercept is included.

- The `MODEL` subcommand defines the cells in a design. In particular, cells are defined by all of the possible combinations of levels of the factors in the design. The number of cells equals the product of the number of levels of all the factors. A design is balanced if each cell contains the same number of cases. `CSGLM` can analyze balanced and unbalanced designs.
- The format is a list of effects to be included in the model, separated by spaces or commas.
- If the `MODEL` subcommand is not specified, `CSGLM` uses a model that includes the intercept term (unless it is excluded on the `INTERCEPT` subcommand), main effects for any factors, and any covariates.
- To include a term for the main effect of a factor, enter the name of the factor.
- To include a term for an interaction between factors, use the keyword `BY` or the asterisk (`*`) to join the factors that are involved in the interaction. For example, `A*B` means a two-way interaction effect of `A` and `B`, where `A` and `B` are factors. `A*A` is not allowed because factors inside an interaction effect must be distinct.
- To include a term for nesting one effect within another effect, use a pair of parentheses. For example, `A(B)` means that `A` is nested within `B`. When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.
- Multiple nesting is allowed. For example, `A(B(C))` means that `B` is nested within `C`, and `A` is nested within `B(C)`.
- Interactions between nested effects are not valid. For example, neither `A(C)*B(C)` nor `A(C)*B(D)` is valid.
- To include a covariate term in the design, enter the name of the covariate.
- Covariates can be connected, but not nested, through the `*` operator to form another covariate effect. Interactions among covariates such as `X1*X1` and `X1*X2` are valid, but `X1(X2)` is not.
- Factor and covariate effects can be connected only by the `*` operator. Suppose `A` and `B` are factors, and `X1` and `X2` are covariates. Examples of valid factor-by-covariate interaction effects are `A*X1`, `A*B*X1`, `X1*A(B)`, `A*X1*X1`, and `B*X1*X2`.

## **INTERCEPT Subcommand**

The `INTERCEPT` subcommand controls whether an intercept term is included in the model. This subcommand can also be used to display or suppress the intercept term in output tables.



## **INCLUDE Keyword**

The `INCLUDE` keyword specifies whether the intercept is included in the model, or the keyword requests the intercept-only model.

- YES**     *The intercept is included in the model.* This setting is the default.
- NO**       *The intercept is not included in the model.* If no factors or covariates are defined, specifying `INCLUDE = NO` is invalid syntax.
- ONLY**    *The intercept-only model is fit.* If the `MODEL` subcommand is specified, specifying `INCLUDE = ONLY` is invalid syntax.

## **SHOW Keyword**

The `SHOW` keyword specifies whether the intercept is displayed or suppressed in output tables.

- YES**     *The intercept is displayed in output tables.* This setting is the default.
- NO**       *The intercept is not displayed in output tables.* If `INCLUDE = NO` or `ONLY` is specified, `SHOW = NO` is ignored.

## **Example**

```
CSGLM y BY a b c
      /PLAN FILE='/survey/myfile.csplan'
      /INTERCEPT INCLUDE = ONLY.
```

- The preceding syntax defines the model space using factors A, B, and C but fits the intercept-only model.

## **CUSTOM Subcommand**

The `CUSTOM` subcommand defines custom hypothesis tests by specifying the **L** matrix (contrast coefficients matrix) and the **K** matrix (contrast results matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ . The vector **B** is the parameter vector in the linear model.

- Multiple `CUSTOM` subcommands are allowed. Each subcommand is treated independently.
- An optional label may be specified by using the `LABEL` keyword. The label is a string with a maximum length of 255 characters. Only one label can be specified.
- Either the `LMATRIX` or `KMATRIX` keyword, or both, must be specified.

**LMATRIX**                    *Contrast coefficients matrix.* This matrix specifies coefficients of contrasts, which can be used for studying the effects in the model. An **L** matrix can be specified by using the `LMATRIX` keyword.

**KMATRIX**                    *Contrast results matrix.* This matrix specifies the results of the linear hypothesis. A **K** matrix can be specified by using the `KMATRIX` keyword.

- The number of rows in the **L** and **K** matrices must be equal.
- A custom hypothesis test can be formed by specifying an **L** or **K** matrix, or both. If only one matrix is specified, the unspecified matrix uses the defaults described below.

- If `KMATRIX` is specified but `LMATRIX` is not specified, the **L** matrix is assumed to be the row vector corresponding to the intercept in the estimable function, provided that `INCLUDE = YES` or `ONLY` is specified on the `INTERCEPT` subcommand. In this case, the **K** matrix can be only a scalar matrix.
- The default **K** matrix is a zero matrix; that is,  $\mathbf{LB} = \mathbf{0}$  is assumed.
- There are three general formats that can be used on the `LMATRIX` keyword: (1) Specify a coefficient value for the intercept, followed optionally by an effect name and a list of real numbers. (2) Specify an effect name and a list of real numbers. (3) Specify keyword `ALL` and a list of real numbers. In all three formats, there can be multiple effect names (or instances of the keyword `ALL`) and number lists.
- Only valid effects in the default model or on the `MODEL` subcommand can be specified on the `LMATRIX` keyword.
- The length of the list of real numbers on the `LMATRIX` keyword must be equal to the number of parameters (including the redundant parameters) corresponding to the specified effect. For example, if the effect `A*B` takes up six columns in the design matrix, the list after `A*B` must contain exactly six numbers.
- When `ALL` is specified, the length of the list that follows `ALL` must be equal to the total number of parameters (including the redundant parameters) in the model.
- Effects that are in the model but not specified on the `LMATRIX` keyword are assumed to have entries of 0 in the corresponding columns of the **L** matrix.
- When an **L** matrix is being defined, a number can be specified as a fraction with a positive denominator. For example,  $1/3$  and  $-1/3$  are valid, but  $1/-3$  is invalid.
- A semicolon (;) indicates the end of a row in the **L** matrix.
- The format for the `KMATRIX` keyword is one or more real numbers. If more than one number is specified, then separate adjacent numbers using a semicolon (;). Each semicolon indicates the end of a row in the **K** matrix. Each number is the hypothesized value for a contrast, which is defined by a row in the **L** matrix.
- For the `KMATRIX` keyword to be valid, either the `LMATRIX` keyword, or `INCLUDE = YES` on the `INTERCEPT` subcommand, must be specified.

### Example

Suppose that factors A and B each have three levels.

```
CSGLM y BY a b
  /PLAN FILE='/survey/myfile.csplan'
  /MODEL a b a*b
  /CUSTOM LABEL = "Effect A"
    LMATRIX = a 1 0 -1
              a*b 1/3 1/3 1/3
                  0 0 0
                  -1/3 -1/3 -1/3;
              a 0 1 -1
              a*b 0 0 0
                  1/3 1/3 1/3
                  -1/3 -1/3 -1/3.
```

- The preceding syntax specifies a test of effect A.

- Because there are three levels in effect A, two independent contrasts can be formed at most; thus, there are two rows in the **L** matrix, separated by a semicolon (;).
- There are three levels each in effects A and B; thus, the interaction effect A\*B takes nine columns in the design matrix.
- The first row in the **L** matrix tests the difference between levels 1 and 3 of effect A; the second row tests the difference between levels 2 and 3 of effect A.
- The **KMATRIX** keyword is not specified, so the null hypothesis value for both tests is 0.

### Example

Suppose that factors A and B each have three levels.

```
CSGLM y BY a b
  /PLAN FILE='/survey/myfile.csplan'
  /CUSTOM LABEL = "Effect A"
    LMATRIX = a 1 0 -1; a 1 -1 0
  /CUSTOM LABEL = "Effect B"
    LMATRIX = b 1 0 -1; b 1 -1 0
    KMATRIX = 0; 0.
```

- The preceding syntax specifies tests of effects A and B.
- The **MODEL** subcommand is not specified, so the default model—which includes the intercept and main effects for A and B—is used.
- There are two **CUSTOM** subcommands; each subcommand specifies two rows in the **L** matrix.
- The first **CUSTOM** subcommand does not specify the **KMATRIX** keyword. By default, this subcommand tests whether the effect of factor A is 0.
- The second **CUSTOM** subcommand specifies the **KMATRIX** keyword. This subcommand tests whether the effect of factor B is 0.

## EMMEANS Subcommand

The **EMMEANS** subcommand displays estimated marginal means of the dependent variable in the cells for the specified factors. Note that these means are predicted, not observed, means.

- Multiple **EMMEANS** subcommands are allowed. Each subcommand is treated independently.
- The **EMMEANS** subcommand may be specified with no additional keywords. The output for an empty **EMMEANS** subcommand is the overall estimated marginal mean of the dependent variable, collapsing over any factors, and with any covariates held at their overall means.

### **TABLES = option**

Valid options are factors appearing on the factor list and crossed factors that are constructed of factors on the factor list. Crossed factors can be specified by using an asterisk (\*) or the keyword **BY**. All factors in a crossed factor specification must be unique. If a factor or a crossing of factors is specified on the **TABLES** keyword, **CSGLM** collapses over any other factors before computing the estimated marginal means for the dependent variable. If the **TABLES** keyword is not specified, the overall estimated marginal mean of the dependent variable, collapsing over any factors, is computed.

### **OTHER = [option]**

Specifies the covariate values to use when computing the estimated marginal means. If the **OTHER** keyword is used, it must be followed by

an equals sign and one or more elements enclosed in square brackets. Valid elements are covariates appearing on the CSGLM covariate list, each of which must be followed by a numeric value or the keyword MEAN in parentheses.

If a numeric value is used, the estimated marginal mean will be computed by holding the specified covariate at the supplied value. If the keyword MEAN is used, the estimated marginal mean will be computed by holding the covariate at its overall mean. If a covariate is not specified on the OTHER option, its overall mean will be used in estimated marginal mean calculations.

Any covariate may occur only once on the OTHER keyword.

**CONTRAST = type**

Specifies the type of contrast that is desired among the levels of the factor that is given on the COMPARE keyword. This keyword creates an L matrix such that the columns corresponding to the factor match the contrast that is given. The other columns are adjusted so that the L matrix is estimable. Available contrast types and their options are described in a separate table below. The CONTRAST keyword is ignored if the COMPARE keyword is not specified.

**COMPARE = factor**

Compares levels of a factor specified on the TABLES keyword and displays results for each individual comparison as well as for the overall set of comparisons. If only one factor is specified on TABLES, COMPARE can be specified by itself; otherwise, the factor specification is required. In the latter case, levels of the specified factor are compared for each level of the other factors that are specified on TABLES. The type of comparison that is performed is determined by the CONTRAST keyword. The TABLES keyword must be specified for the COMPARE keyword to be valid.

## ***CONTRAST Keyword***

The contrast types that may be specified on the CONTRAST keyword are described below.

The CSGLM procedure sorts levels of the factor in ascending order and defines the highest level as the last level. (If the factor is a string variable, the value of the highest level is locale-dependent.)

**SIMPLE (value)**

*Each level of the factor (except the highest level) is compared to the highest level. SIMPLE is the default contrast type if the COMPARE keyword is specified.*

The SIMPLE keyword may be followed optionally by parentheses containing a value. Put the value inside a pair of quotation marks if the value is formatted (such as date or currency) or if the factor is of string type. If a value is specified, the factor level with that value is used as the omitted reference category. If the specified value does not exist in the data, a warning is issued and the highest level is used.

An example is as follows: CSGLM y BY a ... /EMMEANS TABLES=a COMPARE=a CONTRAST=SIMPLE(1) . The specified contrast compares all levels of factor A (except level 1) to level 1. Simple contrasts are not orthogonal.

**DEVIATION (value)**

*Each level of the factor (except the highest level) is compared to the grand mean.*

The `DEVIATION` keyword may be followed optionally by parentheses containing a value. Put the value inside a pair of quotation marks if the value is formatted (such as date or currency) or if the factor is of string type. If a value is specified, the factor level with that value is used as the omitted reference category. If the specified value does not exist in the data, a warning is issued and the highest level is used.

An example is as follows: `CSGLM y BY a ... /EMMEANS TABLES=a COMPARE=a CONTRAST=DEVIATION(1)`. The specified contrast omits level 1 of A. Deviation contrasts are not orthogonal.

**DIFFERENCE**

*Each level of the factor (except the lowest level) is compared to the mean of previous levels.* In a balanced design, difference contrasts are orthogonal.

**HELMERT**

*Each level of the factor (except the highest level) is compared to the mean of subsequent levels.* In a balanced design, Helmert contrasts are orthogonal.

**REPEATED**

*Each level of the factor (except the highest level) is compared to the previous level.* Repeated contrasts are not orthogonal.

**POLYNOMIAL (number list)**

*Polynomial contrasts.* The first degree of freedom contains the linear effect across the levels of the factor, the second contains the quadratic effect, and so on. By default, the levels are assumed to be equally spaced; the default metric is (1 2 . . . k), where k levels are involved.

The `POLYNOMIAL` keyword may be followed optionally by parentheses containing a number list. Numbers in the list must be separated by spaces or commas. Unequal spacing may be specified by entering a metric consisting of one integer for each level of the factor. Only the relative differences between the terms of the metric matter. Thus, for example, (1 2 4) is the same metric as (2 3 5) or (20 30 50) because, in each instance, the difference between the second and third numbers is twice the difference between the first and second numbers. All numbers in the metric must be unique; thus, (1 1 2) is not valid.

An example is as follows: `CSGLM y BY a ... /EMMEANS TABLES=a COMPARE=a CONTRAST=POLYNOMIAL(1 2 4)`. Suppose that factor A has three levels. The specified contrast indicates that the three levels of A are actually in the proportion 1:2:4. In a balanced design, polynomial contrasts are orthogonal.

Orthogonal contrasts are particularly useful. In a balanced design, contrasts are orthogonal if the sum of the coefficients in each contrast row is 0 and if, for any pair of contrast rows, the products of corresponding coefficients sum to 0.

## **CRITERIA Subcommand**

The `CRITERIA` subcommand controls statistical criteria and specifies numerical tolerance for checking singularity.

**CILEVEL = value**

*Confidence interval level for coefficient estimates and estimated marginal means.* Specify a value that is greater than or equal to 0 and less than 100. The default value is 95.

**DF = value**

*Sampling design degrees of freedom to use in computing p values for all test statistics.* Specify a positive number. The default value is the difference between the number of primary sampling units and the number of strata in the first stage of sampling.

**SINGULAR = value**

*Tolerance value used to test for singularity.* Specify a positive value. The default value is  $10^{-12}$ .

## **STATISTICS Subcommand**

The `STATISTICS` subcommand requests various statistics associated with the coefficient estimates.

- There are no default keywords on the `STATISTICS` subcommand. If this subcommand is not specified, no statistics that are listed below are displayed.

<b>PARAMETER</b>	<i>Coefficient estimates.</i>
<b>SE</b>	<i>Standard error for each coefficient estimate.</i>
<b>TTEST</b>	<i>t test for each coefficient estimate.</i>
<b>CINTERVAL</b>	<i>Confidence interval for each coefficient estimate.</i>
<b>DEFF</b>	<i>Design effect for each coefficient estimate.</i>
<b>DEFFSQRT</b>	<i>Square root of the design effect for each coefficient estimate.</i>

## **TEST Subcommand**

The `TEST` subcommand specifies the type of test statistic and the method of adjusting the significance level to be used for hypothesis tests that are requested on the `MODEL`, `CUSTOM`, and `EMMEANS` subcommands.

### **TYPE Keyword**

The `TYPE` keyword indicates the type of test statistic.

<b>F</b>	<i>Wald F test. This is the default test statistic if the <code>TYPE</code> keyword is not specified.</i>
<b>ADJF</b>	<i>Adjusted Wald F test.</i>
<b>CHISQUARE</b>	<i>Wald chi-square test.</i>
<b>ADJCHISQUARE</b>	<i>Adjusted Wald chi-square test.</i>

### **PADJUST keyword**

The `PADJUST` keyword indicates the method of adjusting the significance level.

<b>LSD</b>	<i>Least significant difference. This method does not control the overall probability of rejecting the hypotheses that some linear contrasts are different from the null hypothesis value(s). This setting is the default.</i>
<b>BONFERRONI</b>	<i>Bonferroni. This method adjusts the observed significance level for the fact that multiple contrasts are being tested.</i>
<b>SEQBONFERRONI</b>	<i>Sequential Bonferroni. This procedure is a sequentially step-down rejective Bonferroni procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.</i>

---

<b>SIDAK</b>	<i>Sidak.</i> This method provides tighter bounds than the Bonferroni approach.
<b>SEQSIDAK</b>	<i>Sequential Sidak.</i> This procedure is a sequentially rejective step-down rejective Sidak procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.

## ***DOMAIN Subcommand***

The `DOMAIN` subcommand specifies the subpopulation for which the analysis is to be performed.

- Keyword `VARIABLE`, followed by an equals sign, a variable, and a value in parentheses, are required. Put the value inside a pair of quotation marks if the value is formatted (such as date or currency) or if the variable is of string type.
- The subpopulation is defined by all cases having the given value on the specified variable.
- Analyses are performed only for the specified subpopulation.
- For example, `DOMAIN VARIABLE = myvar (1)` defines the subpopulation by all cases for which variable `MYVAR` has value 1.
- The specified variable may be numeric or string and must exist at the time that the `CSGLM` procedure is invoked.
- Stratification or cluster variables may be specified, but no other plan file variables are allowed on the `DOMAIN` subcommand.
- Analysis variables may not be specified on the `DOMAIN` subcommand.

## ***MISSING Subcommand***

The `MISSING` subcommand specifies how missing values are handled.

- All design variables, as well as the dependent variable and any covariates, must have valid data. Cases with invalid data for any of these variables are deleted from the analysis.
- The `CLASSMISSING` keyword specifies whether user-missing values are treated as valid. This specification is applied to categorical design variables (i.e., strata, cluster, and subpopulation variables) and any factors.

<b>EXCLUDE</b>	<i>Exclude user-missing values among the strata, cluster, subpopulation, and factor variables.</i> This setting is the default.
----------------	---

<b>INCLUDE</b>	<i>Include user-missing values among the strata, cluster, subpopulation, and factor variables.</i> Treat user-missing values for these variables as valid data.
----------------	---

## ***PRINT Subcommand***

The `PRINT` subcommand is used to display optional output.

- If the `PRINT` subcommand is not specified, the default output includes sample information, variable and factor information, and model summary statistics.
- If the `PRINT` subcommand is specified, `CSGLM` displays output only for those keywords that are specified.

<b>SAMPLEINFO</b>	<i>Sample information table.</i> Displays summary information about the sample, including the unweighted count and the population size. This output is default output if the <code>PRINT</code> subcommand is not specified.
<b>VARIABLEINFO</b>	<i>Variable information.</i> Displays summary information about the dependent variable, covariates, and factors. This output is default output if the <code>PRINT</code> subcommand is not specified.
<b>SUMMARY</b>	<i>Model summary statistics.</i> Displays $R^2$ and root mean squared error statistics. This output is default output if the <code>PRINT</code> subcommand is not specified.
<b>GEF</b>	<i>General estimable function table.</i>
<b>LMATRIX</b>	<i>Set of contrast coefficients (L) matrices.</i>
<b>COVB</b>	<i>Covariance matrix for regression coefficients.</i>
<b>CORB</b>	<i>Correlation matrix for regression coefficients.</i>
<b>NONE</b>	<i>No PRINT subcommand output.</i> None of the <code>PRINT</code> subcommand output is displayed. However, if <code>NONE</code> is specified with one or more other keywords, the other keywords override <code>NONE</code> .

## ***SAVE Subcommand***

The `SAVE` subcommand adds predicted or residual values to the active dataset.

- Specify one or more temporary variables, each variable followed by an optional new name in parentheses.
- The optional names must be unique, valid variable names.
- If new names are not specified, `CSGLM` uses the default names. If the default names conflict with existing variable names, a suffix is added to the default name to make it unique.

<b>PRED</b>	<i>Saves predicted values.</i> The default variable name is Predicted.
<b>RESID</b>	<i>Saves residuals.</i> The default variable name is Residual.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand saves an external IBM® SPSS® Statistics data file or previously declared dataset containing the parameter covariance or correlation matrix with parameter estimates, standard errors, significance values, and sampling design degrees of freedom. It also saves the parameter estimates and the parameter covariance matrix in XML format.

- At least one keyword and a filename are required. Specify the keyword followed by a quoted file specification.
- The `COVB` and `CORB` keywords are mutually exclusive, as are the `MODEL` and `PARAMETER` keywords.



- The filename must be specified in full. CSGLM does not supply an extension.
- For COVB and CORB, you can specify a previously declared dataset name (DATASET DECLARE command) instead of a file specification.

**COVB = 'savfile'|'dataset'**      *Writes the parameter covariance matrix and other statistics to an external SPSS Statistics datafile or previously declared dataset.*

**CORB = 'savfile'|'dataset'**      *Writes the parameter correlation matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**MODEL = 'file'**                      *Writes the parameter estimates and the parameter covariance matrix to an XML file.*

**PARAMETER = 'file'**                *Writes the parameter estimates to an XML file.*

# CSLOGISTIC

CSLOGISTIC is available in the Complex Samples option.

*Note:* Square brackets that are used in the CSLOGISTIC syntax chart are required parts of the syntax and are not used to indicate optional elements. Equals signs (=) that are used in the syntax chart are required elements. All subcommands, save the PLAN subcommand, are optional.

```
CSLOGISTIC dependent var ({LOW }) BY factor list WITH covariate list
                        {HIGH**}
                        {value }

/PLAN FILE = file

/JOINTPROB FILE = file

/MODEL effect list

/INTERCEPT INCLUDE = {YES**} SHOW = {YES**}
                      {NO }          {NO }
                      {ONLY }

/CUSTOM LABEL = "label"

  LMATRIX = {number effect list effect list ...; ...}
            {number effect list effect list ... }
            {effect list effect list ...; ... }
            {effect list effect list ... }
            {ALL list; ALL ... }
            {ALL list }

  KMATRIX = {number; number; ...}
            {number }

/CUSTOM ...

/ODDSRATIOS {FACTOR = [varname ({LOW }) varname ...] }
            {HIGH**}
            {value }

            {COVARIATE = [varname ({1** }) varname ...]}
            {number list}

            CONTROL = [varname (value) varname (value) ...]

/ODDSRATIOS ...

/CRITERIA CHKSEP = {20**} CILEVEL = {95** } DF = n
                  {n }          {value}

LCONVERGE = [{0** } {RELATIVE**}] MXITER = {100**}
            {value} {ABSOLUTE }          {n }

MXSTEP = {5**} PCONVERGE = [{1E-6**} {RELATIVE**}]
                  {n }          {value } {ABSOLUTE }

SINGULAR = {1E-12**}
           {value }

/STATISTICS PARAMETER EXP SE TTEST CINTERVAL DEFF DEFFSQRT

/TEST TYPE = {F** } PADJUST = {LSD** }
             {ADJF }          {BONFERRONI }
             {CHISQUARE }     {SEQBONFERRONI}
             {ADJCHISQUARE}   {SIDAK }
             {ADJCHISQUARE}   {SEQSIDAK }
             {ADJCHISQUARE}   {SEQSIDAK }
```

```

/DOMAIN VARIABLE = varname (value)

/MISSING CLASSMISSING = {EXCLUDE**}
                        {INCLUDE  }

/PRINT SAMPLEINFO** VARIABLEINFO** SUMMARY**

        HISTORY({1**}) GEF LMATRIX COVB CORB CLASSTABLE NONE
                {n  }

/SAVE PREDPROB(rootname:{25**}) PREDVAL(varname)
                {n  }

/OUTFILE {COVB = 'savfile'|'dataset'} {MODEL = 'file'      }
         {CORB = 'savfile'|'dataset'} {PARAMETER = 'file'}

```

\*\*Default if the keyword or subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- Command introduced.

Release 17.0

- Added support for SET THREADS.

### **Example**

```

CSLOGISTIC y BY a b c WITH x
  /PLAN FILE='/survey/myfile.csplan'.

```

## **Overview**

CSLOGISTIC performs logistic regression analysis on a binary or multinomial dependent variable using the generalized link function for samples that are drawn by complex sampling methods. The procedure estimates variances by taking into account the sample design that is used to select the sample, including equal probability and probability proportional to size (PPS) methods, and with replacement (WR) and without replacement (WOR) sampling procedures. Optionally, CSLOGISTIC performs analyses for a subpopulation.

### **Basic Specification**

- The basic specification is a variable list (identifying the dependent variable, the factors, if any, and the covariates, if any) and a PLAN subcommand with the name of a complex sample analysis plan file, which may be generated by the CSPLAN procedure.
- The default model includes the intercept term, main effects for any factors, and any covariates.
- The basic specification displays summary information about the sample and all analysis variables, model summary statistics, and Wald  $F$  tests for all model effects. Additional subcommands must be used for other output.

**Operations**

- CSLOGISTIC performs logistic regression analysis for sampling designs that are supported by the CSPLAN and CSSELECT procedures.
- The input dataset must contain the variables to be analyzed and variables that are related to the sampling design.
- The complex sample analysis plan file provides an analysis plan based on the sampling design.
- By default, CSLOGISTIC uses a model that includes the intercept term, main effects for any factors, and any covariates.
- Other effects, including interaction and nested effects, may be specified by using the MODEL subcommand.
- The default output for the specified model is summary information about the sample and all analysis variables, model summary statistics, and Wald  $F$  tests for all model effects.
- WEIGHT and SPLIT FILE settings are ignored by the CSLOGISTIC procedure.
- This procedure uses the multithreaded options specified by SET THREADS.

**Syntax Rules**

- The dependent variable and PLAN subcommand are required. All other variables and subcommands are optional.
- Multiple CUSTOM and ODDS RATIOS subcommands may be specified; each subcommand is treated independently. All other subcommands may be specified only once.
- Empty subcommands are not allowed; all subcommands must be specified with options.
- Each keyword may be specified only once within a subcommand.
- Subcommand names and keywords must be spelled in full.
- Equals signs (=) that are shown in the syntax chart are required.
- Subcommands may be specified in any order.
- The dependent variable, factors, and the subpopulation variable can be numeric or string variables, but covariates must be numeric.
- Across the dependent, factor, and covariate variable lists, a variable may be specified only once.
- Plan file and subpopulation variables may not be specified on the variable list.
- Minimum syntax is a dependent variable and the PLAN subcommand. This specification fits an intercept-only model.

**Limitations**

- WEIGHT and SPLIT FILE settings are ignored with a warning by the CSLOGISTIC procedure.

**Examples**

```
* Complex Samples Logistic Regression.
CSLOGISTIC default(LOW) BY ed WITH age employ address income debtinc creddebt othdebt
/PLAN FILE = 'samplesDirectory\bankloan.csaplan'
/MODEL ed age employ address income debtinc creddebt othdebt
```

```

/INTERCEPT INCLUDE=YES SHOW=YES
/STATISTICS PARAMETER EXP SE CINTERVAL DEFF
/TEST TYPE=F PADJUST=LSD
/ODDSRATIOS FACTOR=[ed(HIGH)]
/ODDSRATIOS COVARIATE=[employ(1)]
/ODDSRATIOS COVARIATE=[debtinc(1)]
/MISSING CLASSMISSING=EXCLUDE
/CRITERIA MXITER=100 MXSTEP=5
PCONVERGE=[1e-006 RELATIVE] CHKSEP=20
CILEVEL=95
/PRINT SUMMARY CLASSTABLE VARIABLEINFO SAMPLEINFO .

```

- The procedure fits a logistic regression model for the dependent variable *default* (with the lowest value as the reference category) using *ed* as a factor and *age*, *employ*, *address*, *income*, *debtinc*, *creddebt*, and *othdebt* as covariates.
- The complex sampling analysis plan is contained in the file *bankloan.csaplan*.
- The model specification calls for a main effects model with intercept.
- Parameter estimates, their standard errors, 95% confidence intervals, and exponentiated parameter estimates and their 95% confidence intervals are requested.
- A classification table is requested in addition to the default model output.
- Odds ratios are produced for the factor *ed* and the covariates *employ* and *debtinc*, using the default reference category and change in value, respectively.
- All other options are set to their default values.

## CSLOGISTIC Variable List

The variable list specifies the dependent variable and reference category, the factors, and the covariates in the model.

- The dependent variable must be the first specification on CLOGISTIC.
- The dependent variable can be numeric or string.
- The CSLOGISTIC procedure sorts levels of the dependent variable in ascending order and defines the highest level as the last level. (If the dependent variable is a string variable, the value of the highest level is locale-dependent.) By default, the highest response category is used as the base (or reference) category.
- A custom reference category may be specified in parentheses immediately following the dependent variable.

<b>LOW</b>	<i>The lowest category is the reference category.</i>
<b>HIGH</b>	<i>The highest category is the reference category.</i> This setting is the default.
<b>value</b>	<i>User-specified reference category.</i> The category that corresponds to the specified value is the reference category. Put the value inside a pair of quotation marks if the value is formatted (such as date or time) or if the dependent variable is of string type. Note, however, that this does not work for custom currency formats.

- If a value is specified as the reference category of the dependent variable, but the value does not exist in the data, a warning is issued and the default HIGH is used.
- The names of the factors and covariates, if any, follow the dependent variable. Specify any factors following the keyword BY. Specify any covariates following the keyword WITH.

- Factors can be numeric or string variables, but covariates must be numeric.
- Each variable may be specified only once on the variable list.
- Plan file and subpopulation variables may not be specified on the variable list.

## ***PLAN Subcommand***

The `PLAN` subcommand specifies the name of an XML file containing analysis design specifications. This file is written by the `CSPLAN` procedure.

- The `PLAN` subcommand is required.

**FILE** *Specifies the name of an external file.*

## ***JOINTPROB Subcommand***

The `JOINTPROB` subcommand is used to specify the file or dataset containing the first stage joint inclusion probabilities for `UNEQUAL_WOR` estimation. The `CSSELECT` procedure writes this file in the same location and with the same name (but different extension) as the plan file. When `UNEQUAL_WOR` estimation is specified, the `CSLOGISTIC` procedure will use the default location and name of the file unless the `JOINTPROB` subcommand is used to override them.

**FILE** *Specifies the name of the file or dataset containing the joint inclusion probabilities.*

## ***MODEL Subcommand***

The `MODEL` subcommand is used to specify the effects to be included in the model. Use the `INTERCEPT` subcommand to control whether the intercept is included.

- The `MODEL` subcommand defines the cells in a design. In particular, cells are defined by all of the possible combinations of levels of the factors in the design. The number of cells equals the product of the number of levels of all the factors. A design is balanced if each cell contains the same number of cases. `CSLOGISTIC` can analyze balanced and unbalanced designs.
- The format is a list of effects to be included in the model, separated by spaces or commas.
- If the `MODEL` subcommand is not specified, `CSLOGISTIC` uses a model that includes the intercept term (unless it is excluded on the `INTERCEPT` subcommand), main effects for any factors, and any covariates.
- To include a term for the main effect of a factor, enter the name of the factor.
- To include a term for an interaction between factors, use the keyword `BY` or the asterisk (\*) to join the factors that are involved in the interaction. For example, `A*B` means a two-way interaction effect of `A` and `B`, where `A` and `B` are factors. `A*A` is not allowed because factors that are inside an interaction effect must be distinct.
- To include a term for nesting one effect within another effect, use a pair of parentheses. For example, `A(B)` means that `A` is nested within `B`. When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.

- Multiple nesting is allowed. For example, A(B(C)) means that B is nested within C, and A is nested within B(C).
- Interactions between nested effects are not valid. For example, neither A(C)\*B(C) nor A(C)\*B(D) is valid.
- To include a covariate term in the design, enter the name of the covariate.
- Covariates can be connected, but not nested, through the \* operator to form another covariate effect. Interactions among covariates such as X1\*X1 and X1\*X2 are valid, but X1(X2) is not.
- Factor and covariate effects can be connected only by the \* operator. Suppose A and B are factors, and X1 and X2 are covariates. Examples of valid factor-by-covariate interaction effects are A\*X1, A\*B\*X1, X1\*A(B), A\*X1\*X1, and B\*X1\*X2.

## ***INTERCEPT Subcommand***

The INTERCEPT subcommand controls whether an intercept term is included in the model. This subcommand can also be used to display or suppress the intercept term in output tables.

## ***INCLUDE Keyword***

The INCLUDE keyword specifies whether the intercept is included in the model, or the keyword requests the intercept-only model.

- YES**     *The intercept is included in the model.* This setting is the default.
- NO**       *The intercept is not included in the model.* If no factors or covariates are defined, specifying INCLUDE = NO is invalid syntax.
- ONLY**    *The intercept-only model is fit.* If the MODEL subcommand is specified, specifying INCLUDE = ONLY is invalid syntax.

## ***SHOW Keyword***

The SHOW keyword specifies whether the intercept is displayed or suppressed in output tables.

- YES**     *The intercept is displayed in output tables.* This setting is the default.
- NO**       *The intercept is not displayed in output tables.* If INCLUDE = NO or ONLY is specified, SHOW = NO is ignored.

## ***Example***

```
CSLOGISTIC y BY a b c
/PLAN FILE='/survey/myfile.csplan'
/INTERCEPT INCLUDE = ONLY.
```

- The preceding syntax defines the model space using factors A, B, and C but fits the intercept-only model.

## ***CUSTOM Subcommand***

The `CUSTOM` subcommand defines custom hypothesis tests by specifying the **L** matrix (contrast coefficients matrix) and the **K** matrix (contrast results matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ . The vector **B** is the parameter vector in the linear model.

For a binary dependent variable, `CSLOGISTIC` models a single logit. In this case, there is one set of parameters associated with the logit.

For a multinomial dependent variable with *K* levels, `CSLOGISTIC` models *K*–1 logits. In this case, there are *K*–1 sets of parameters, each associated with a different logit. The `CUSTOM` subcommand allows you to specify an **L** matrix in which the same or different contrast coefficients are used across logits.

- Multiple `CUSTOM` subcommands are allowed. Each subcommand is treated independently.
- An optional label may be specified by using the `LABEL` keyword. The label is a string with a maximum length of 255 characters. Only one label can be specified.
- Either the `LMATRIX` or `KMATRIX` keyword, or both, must be specified.

**LMATRIX**            *Contrast coefficients matrix.* This matrix specifies coefficients of contrasts, which can be used for studying the effects in the model. An **L** matrix can be specified by using the `LMATRIX` keyword.

**KMATRIX**            *Contrast results matrix.* This matrix specifies the results of the linear hypothesis. A **K** matrix can be specified by using the `KMATRIX` keyword.

- The number of rows in the **L** and **K** matrices must be equal.
- A custom hypothesis test can be formed by specifying an **L** or **K** matrix, or both. If only one matrix is specified, the unspecified matrix uses the defaults described below.
- If `KMATRIX` is specified but `LMATRIX` is not specified, the **L** matrix is assumed to be the row vector corresponding to the intercept in the estimable function, provided that `INCLUDE = YES` or `ONLY` is specified on the `INTERCEPT` subcommand.
- The default **K** matrix is a zero matrix; that is,  $\mathbf{LB} = \mathbf{0}$  is assumed.
- There are three general formats that can be used on the `LMATRIX` keyword: (1) Specify a coefficient value for the intercept, followed optionally by an effect name and a list of real numbers. (2) Specify an effect name and a list of real numbers. (3) Specify keyword `ALL` and a list of real numbers. In all three formats, there can be multiple effect names (or instances of the keyword `ALL`) and number lists.
- Only valid effects in the default model or on the `MODEL` subcommand can be specified on the `LMATRIX` keyword.
- The length of the list of real numbers on the `LMATRIX` keyword must be equal to the number of parameters (including the redundant parameters) corresponding to the specified effect. For example, if the effect `A*B` takes up six columns in the design matrix, the list after `A*B` must contain exactly six numbers.
- When `ALL` is specified, the length of the list that follows `ALL` must be equal to the total number of parameters (including the redundant parameters) in the model. For a binary dependent variable, the contrast coefficients for the one set of parameters must be listed following the `ALL` keyword. For a multinomial dependent variable with *K* levels, the contrast coefficients



for the  $K-1$  sets of parameters must be listed in order following the `ALL` keyword. That is, first list all parameters (including the redundant parameters) for the first logit, then list all parameters for the second logit, and so forth.

- In general, for a multinomial dependent variable with  $K$  levels, each contrast and its associated hypothesized value are generated separately for each of the  $K-1$  logits; that is, any given contrast is generated  $K-1$  times. However, if the `LMATRIX ALL` keyword is used to define a contrast, then that contrast and its associated hypothesized value are generated once, simultaneously covering all logits.
- Effects that are in the model but not specified on the `LMATRIX` keyword are assumed to have entries of 0 in the corresponding columns of the **L** matrix.
- When an **L** matrix is defined, a number can be specified as a fraction with a positive denominator. For example,  $1/3$  and  $-1/3$  are valid, but  $1/-3$  is invalid.
- A semicolon (;) indicates the end of a row in the **L** matrix.
- The format for the `KMATRIX` keyword is one or more real numbers. If more than one number is specified, then separate adjacent numbers using a semicolon (;). Each semicolon indicates the end of a row in the **K** matrix. Each number is the hypothesized value for a contrast, which is defined by a row in the **L** matrix.
- For the `KMATRIX` keyword to be valid, either the `LMATRIX` keyword, or `INCLUDE = YES` on the `INTERCEPT` subcommand, must be specified.

### Example

Suppose that dependent variable  $Y$  is binary, and factors  $A$  and  $B$  each have three levels.

```
CSLOGISTIC y BY a b
  /PLAN FILE='/survey/myfile.csplan'
  /MODEL a b a*b
  /CUSTOM LABEL = 'Effect A'
  LMATRIX = a 1 0 -1
            a*b 1/3 1/3 1/3
              0 0 0
              -1/3 -1/3 -1/3;
            a 0 1 -1
            a*b 0 0 0
              1/3 1/3 1/3
              -1/3 -1/3 -1/3.
```

- The preceding syntax specifies a test of effect  $A$ .
- Because there are three levels in effect  $A$ , two independent contrasts can be formed at most; thus, there are two rows in the **L** matrix, separated by a semicolon (;).
- There are three levels each in effects  $A$  and  $B$ ; thus, the interaction effect  $A*B$  takes nine columns in the design matrix.
- The first row in the **L** matrix tests the difference between levels 1 and 3 of effect  $A$ ; the second row tests the difference between levels 2 and 3 of effect  $A$ .
- The `KMATRIX` keyword is not specified, so the null hypothesis value for both tests is 0.

**Example**

Suppose that dependent variable  $Z$  and factor  $A$  each have three levels.

```
CSLOGISTIC z BY a
/PLAN FILE='/survey/myfile.csplan'
/MODEL a
/CUSTOM LABEL = 'Effect A'
  LMATRIX = a 1 0 -1;
            a 0 1 -1
```

- The dependent variable  $Z$  has three categories, so there will be two logits.
- The syntax specifies a model with an intercept and a main effect for factor  $A$  and a custom hypothesis test of effect  $A$ .
- Because the `ALL` option is not used on the `LMATRIX` keyword, the same set of contrast coefficients for the parameters will be used across both logits. That is, the resulting  $L$  matrix is block diagonal with the same 2-by-4 matrix of coefficients in each block. The equivalent `LMATRIX` keyword using the `ALL` option is as follows:

```
LMATRIX = ALL 0 1 0 -1 0 0 0 0;
           ALL 0 0 1 -1 0 0 0 0;
           ALL 0 0 0 0 0 1 0 -1;
           ALL 0 0 0 0 0 0 1 -1
```

**Example**

Suppose that dependent variable  $Z$  has three categories, and factors  $A$  and  $B$  each have three levels.

```
CSLOGISTIC z BY a b
/PLAN FILE='/survey/myfile.csplan'
/CUSTOM LABEL = 'Effect A for All Logits'
  LMATRIX = a 1 0 -1; a 0 1 -1
/CUSTOM LABEL = 'Effect A for 1st Logit, Effect B for 2nd Logit'
  LMATRIX = ALL 0 1 0 -1 0 0 0 0 0 0 0 0 1 0 -1;
           ALL 0 0 1 -1 0 0 0 0 0 0 0 0 0 1 -1
  KMATRIX = 0; 0.
```

- The dependent variable  $Z$  has three categories, so there will be two logits.
- The `MODEL` subcommand is not specified; thus the default model—which includes the intercept and main effects for  $A$  and  $B$ —is used.
- The first `CUSTOM` subcommand tests whether the effect of factor  $A$  is 0 across both logits.
- The second `CUSTOM` subcommand specifies different contrast coefficients for each logit. In particular, the  $L$  matrix tests the effect of factor  $A$  for the first logit and factor  $B$  for the second logit. The `KMATRIX` keyword explicitly states that each linear combination that is formed from the contrast coefficients and the parameter estimates is tested against the value 0.

**ODDSRATIOS Subcommand**

The `ODDSRATIOS` subcommand estimates odds ratios for the specified factor(s) or covariate(s). Note that these odds ratios are model-based and are not directly computed by using the observed data.

A separate set of odds ratios is computed for each category of the dependent variable (except the reference category).

If the `FACTOR` keyword is specified, the odds ratios compare the odds at each category  $j$  with the odds at category  $J$ , where  $J$  is the reference category defined in parentheses following the variable name of the factor. All other factors and covariates are fixed as defined on the `CONTROL` keyword.

If the `COVARIATE` keyword is specified, the odds ratios compare the odds at value  $x$  with the odds at value  $x + \Delta x$ , where  $\Delta x$  is the change in  $x$  defined in parentheses following the variable name of the covariate. To define the value  $x$ , specify the covariate and the value on the `CONTROL` keyword. All other factors and covariates are fixed as defined on the `CONTROL` keyword.

If a specified factor or covariate interacts with other predictors in the model, the odds ratios depend not only on the change in the specified variable but also on the values of the variables with which it interacts. If a specified covariate interacts with itself in the model (for example,  $X*X$ ), the odds ratios depend on both the change in the covariate and the value of the covariate. The values of interacting factors and covariates can be customized by using the `CONTROL` keyword.

The `CSLOGISTIC` procedure sorts levels of each factor in ascending order and defines the highest level as the last level. (If the factor is a string variable, the value of the highest level is locale-dependent.)

- Multiple `ODDSRATIOS` subcommands are allowed. Each subcommand is treated independently.
- Either the `FACTOR` keyword and one or more factors, or the `COVARIATE` keyword and one or more covariates, but not both, are required. All other keywords are optional.
- The `FACTOR`, `COVARIATE`, and `CONTROL` keywords must be followed by an equals sign and one or more elements enclosed in square brackets.
- If a variable is specified on the `FACTOR` or `COVARIATE` keyword and is also specified on the `CONTROL` keyword, the `CONTROL` specification for that variable is ignored when the variable's odds ratios are computed. Thus, `FACTOR = [A B] CONTROL = [A(1) B(2)]` estimates odds ratios for factor A holding factor B at level 2 and for factor B holding factor A at level 1.

**FACTOR = [option]**

Valid options are one or more factors appearing on the factor list. Optionally, each factor may be followed by parentheses containing the level to use as the reference category when computing odds ratios. Keyword `LOW` or `HIGH`, or a value, may be specified. Put the value inside a pair of quotes if the value is formatted (such as date or currency) or if the factor is of string type. By default, the highest category is used as the reference category.

If a value is specified but the value does not exist in the data, a warning is issued and the default `HIGH` is used.

Any factor may occur only once on the `FACTOR` keyword.

**COVARIATE = [option]**

Valid options are one or more covariates appearing on the covariate list. Optionally, each covariate may be followed by parentheses containing one or more nonzero numbers giving unit(s) of change to use for covariates when computing odds ratios. Odds ratios are estimated for each distinct value. The default value is 1.

Any covariate may occur only once on the `COVARIATE` keyword.

**CONTROL= [option]**

Specifies the factor and/or covariate values to use when computing odds ratios. Factors must appear on the factor list, and covariates must appear on the covariate list, of the `CSLOGISTIC` command.

Factors must be followed by the keyword `LOW` or `HIGH`, or a value, in parentheses. Put the value inside a pair of quotation marks if the value is formatted (such as date or currency) or if the factor is of string type. If keyword `LOW` or `HIGH` is used, each odds ratio is computed by holding the factor at its lowest or highest level, respectively. If a value is used, each odds ratio is computed by holding the specified factor at the supplied value. If a factor is not specified on the `CONTROL` option, its highest category is used in odds ratio calculations.

If a factor value is specified but the value does not exist in the data, a warning is issued and the default `HIGH` is used.

Covariates must be followed by the keyword `MEAN` or a number in parentheses. If the keyword `MEAN` is used, each odds ratio is computed by holding the covariate at its overall mean. If a number is used, each odds ratio is computed by holding the specified covariate at the supplied value. If a covariate is not specified on the `CONTROL` option, its overall mean is used in odds ratio calculations.

Any factor or covariate may occur only once on the `CONTROL` keyword.

## Example

Suppose that dependent variable *Y* is binary; factor *A* has two levels; and factor *B* has three levels coded 1, 2, and 3.

```
CSLOGISTIC y BY a b WITH x
  /PLAN FILE='/survey/myfile.csplan'
  /MODEL a b a*b x
  /ODDSRATIOS FACTOR=[a] CONTROL=[b(1)]
  /ODDSRATIOS FACTOR=[a] CONTROL=[b(2)]
  /ODDSRATIOS FACTOR=[a] CONTROL=[b(3)].
```

- The default reference category (the highest category) is used for the dependent variable.
- The model includes the intercept, main effects for factors *A* and *B*, the *A*\**B* interaction effect, and the covariate *X*.
- Odds ratios are requested for factor *A*. Assuming the *A*\**B* interaction effect is significant, the odds ratio for factor *A* will differ across levels of factor *B*. The specified syntax requests three odds ratios for factor *A*; each odds ratio is computed at a different level of factor *B*.

## Example

```
CSLOGISTIC y BY a b c WITH x
  /PLAN FILE='/survey/myfile.csplan'
  /MODEL a b c x
  /ODDSRATIOS COVARIATE=[x(1 3 5)].
```

- The preceding syntax will compute three odds ratios for covariate *X*.
- The parenthesized list following variable *X* provides the unit of change values to use when computing odds ratios. Odds ratios will be computed for *X* increasing by 1, 3, and 5 units.

## CRITERIA Subcommand

The CRITERIA subcommand offers controls on the iterative algorithm that is used for estimation, and the subcommand specifies numerical tolerance for checking singularity.

<b>CHKSEP = value</b>	<i>Starting iteration for checking complete separation.</i> Specify a non-negative integer. This criterion is not used if the value is 0. The default value is 20.
<b>CILEVEL = value</b>	<i>Confidence interval level for coefficient estimates, exponentiated coefficient estimates, and odds ratio estimates.</i> Specify a value that is greater than or equal to 0 and less than 100. The default value is 95.
<b>DF = value</b>	<i>Sampling design degrees of freedom to use in computing p values for all test statistics.</i> Specify a positive number. The default value is the difference between the number of primary sampling units and the number of strata in the first stage of sampling.
<b>LCONVERGE = [option]</b>	<i>Log-likelihood function convergence criterion.</i> Convergence is assumed if the absolute or relative change in the log-likelihood function is less than the given value. This criterion is not used if the value is 0. Specify square brackets containing a non-negative number followed optionally by keyword ABSOLUTE or RELATIVE, which indicates the type of change. The default value is 0, and the default type is RELATIVE.
<b>MXITER = value</b>	<i>Maximum number of iterations.</i> Specify a non-negative integer. The default value is 100.
<b>MXSTEP = value</b>	<i>Maximum step-halving allowed.</i> Specify a positive integer. The default value is 5.
<b>PCONVERGE = [option]</b>	<i>Parameter estimates convergence criterion.</i> Convergence is assumed if the absolute or relative change in the parameter estimates is less than the given value. This criterion is not used if the value is 0. Specify square brackets containing a non-negative number followed optionally by keyword ABSOLUTE or RELATIVE, which indicates the type of change. The default value is $10^{-6}$ , and the default type is RELATIVE.
<b>SINGULAR = value</b>	<i>Tolerance value used to test for singularity.</i> Specify a positive value. The default value is $10^{-12}$ .

## STATISTICS Subcommand

The STATISTICS subcommand requests various statistics that are associated with the coefficient estimates.

- There are no default keywords on the STATISTICS subcommand. If this subcommand is not specified, no statistics that are listed below are displayed.

<b>PARAMETER</b>	<i>Coefficient estimates.</i>
<b>EXP</b>	<i>The exponentiated coefficient estimates.</i>
<b>SE</b>	<i>Standard error for each coefficient estimate.</i>
<b>TTEST</b>	<i>t test for each coefficient estimate.</i>
<b>CINTERVAL</b>	<i>Confidence interval for each coefficient estimate and/or exponentiated coefficient estimate.</i>
<b>DEFFSQRT</b>	<i>Square root of the design effect for each coefficient estimate.</i>
<b>DEFF</b>	<i>Design effect for each coefficient estimate.</i>

## TEST Subcommand

The TEST subcommand specifies the type of test statistic and the method of adjusting the significance level to be used for hypothesis tests that are requested on the MODEL and CUSTOM subcommands.

### TYPE Keyword

The TYPE keyword indicates the type of test statistic.

<b>F</b>	<i>Wald F test.</i> This is the default test statistic if the TYPE keyword is not specified.
<b>ADJF</b>	<i>Adjusted Wald F test.</i>
<b>CHISQUARE</b>	<i>Wald chi-square test.</i>
<b>ADJCHISQUARE</b>	<i>Adjusted Wald chi-square test.</i>

### PADJUST Keyword

The PADJUST keyword indicates the method of adjusting the significance level.

<b>LSD</b>	<i>Least significant difference.</i> This method does not control the overall probability of rejecting the hypotheses that some linear contrasts are different from the null hypothesis value(s). This setting is the default.
<b>BONFERRONI</b>	<i>Bonferroni.</i> This method adjusts the observed significance level for the fact that multiple contrasts are being tested.
<b>SEQBONFERRONI</b>	<i>Sequential Bonferroni.</i> This procedure is a sequentially step-down rejective Bonferroni procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.
<b>SIDAK</b>	<i>Sidak.</i> This method provides tighter bounds than the Bonferroni approach.
<b>SEQSIDAK</b>	<i>Sequential Sidak.</i> This procedure is a sequentially rejective step-down rejective Sidak procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.

## DOMAIN Subcommand

The DOMAIN subcommand specifies the subpopulation for which the analysis is to be performed.

- Keyword VARIABLE, followed by an equals sign, a variable, and a value in parentheses, are required. Put the value inside a pair of quotation marks if the value is formatted (such as date or currency) or if the factor is of string type.
- The subpopulation is defined by all cases having the given value on the specified variable.
- Analyses are performed only for the specified subpopulation.
- For example, DOMAIN VARIABLE = myvar (1) defines the subpopulation by all cases for which variable MYVAR has value 1.
- The specified variable may be numeric or string and must exist at the time that the CSLOGISTIC procedure is invoked.

- Stratification or cluster variables may be specified, but no other plan file variables are allowed on the DOMAIN subcommand.
- Analysis variables may not be specified on the DOMAIN subcommand.

## **MISSING Subcommand**

The MISSING subcommand specifies how missing values are handled.

- All design variables, as well as the dependent variable and any covariates, must have valid data. Cases with invalid data for any of these variables are deleted from the analysis.
- The CLASSMISSING keyword specifies whether user-missing values are treated as valid. This specification is applied to categorical design variables (i.e., strata, cluster, and subpopulation variables), the dependent variable, and any factors.

<b>EXCLUDE</b>	<i>Exclude user-missing values among the strata, cluster, subpopulation, and factor variables. This setting is the default.</i>
<b>INCLUDE</b>	<i>Include user-missing values among the strata, cluster, subpopulation, and factor variables. Treat user-missing values for these variables as valid data.</i>

## **PRINT Subcommand**

The PRINT subcommand is used to display optional output.

- If the PRINT subcommand is not specified, the default output includes sample information, variable and factor information, and model summary statistics.
- If the PRINT subcommand is specified, CSLOGISTIC displays output only for those keywords that are specified.

<b>SAMPLEINFO</b>	<i>Sample information table.</i> Displays summary information about the sample, including the unweighted count and the population size. This output is default output if the PRINT subcommand is not specified.
<b>VARIABLEINFO</b>	<i>Variable information.</i> Displays summary information about the dependent variable, covariates, and factors. This output is default output if the PRINT subcommand is not specified.
<b>SUMMARY</b>	<i>Model summary statistics.</i> Displays pseudo- $R^2$ statistics. This output is default output if the PRINT subcommand is not specified.
<b>HISTORY(n)</b>	<i>Iteration history.</i> Displays coefficient estimates and statistics at every $n$ th iteration beginning with the zeroth iteration (the initial estimates). The default is to print every iteration ( $n = 1$ ). The last iteration is always printed if HISTORY is specified, regardless of the value of $n$ .
<b>GEF</b>	<i>General estimable function table.</i>
<b>LMATRIX</b>	<i>Set of contrast coefficients (L) matrices.</i>
<b>COVB</b>	<i>Covariance matrix for regression coefficients.</i>
<b>CORB</b>	<i>Correlation matrix for regression coefficients.</i>
<b>CLASSTABLE</b>	<i>Classification table.</i> Displays frequencies of observed versus predicted response categories.
<b>NONE</b>	<i>No PRINT subcommand output.</i> None of the PRINT subcommand output is displayed. However, if NONE is specified with one or more other keywords, the other keywords override NONE.

## SAVE Subcommand

The `SAVE` subcommand writes optional model variables to the active dataset.

- Specify one or more temporary variables, each variable followed by an optional new name in parentheses.
- The optional names must be unique, valid variable names.
- If new names are not specified, `CSLOGISTIC` generates a name using the temporary variable name with a suffix.

**PREDPROB**     *Predicted probability.* The user-specified or default name is treated as the rootname, and a suffix is added to get new unique variables names. The rootname can be followed by a colon and a positive integer giving the number of predicted probabilities to save. The predicted probabilities of the first  $n$  response categories are saved. One predicted probability variable can be saved for each category of the dependent variable. The default rootname is `PredictedProbability`. The default  $n$  of predicted probabilities to save is 25. To specify  $n$  without a rootname, enter a colon before the number.

**PREDVAL**     *Predicted value.* The class or value that is predicted by the model. The optional variable name must be unique. If the default name is used and it conflicts with existing variable names, a suffix is added to the default name to make it unique. The default variable name is `PredictedValue`.

## OUTFILE Subcommand

The `OUTFILE` subcommand saves an external IBM® SPSS® Statistics data file or previously declared dataset containing the parameter covariance or correlation matrix with parameter estimates, standard errors, significance values, and sampling design degrees of freedom. It also saves the parameter estimates and the parameter covariance matrix in XML format.

- At least one keyword and a file specification are required. The file specification should be enclosed in quotes.
- The `COVB` and `CORB` keywords are mutually exclusive, as are the `MODEL` and `PARAMETER` keywords.
- The filename must be specified in full. `CSLOGISTIC` does not supply an extension.
- For `COVB` and `CORB`, you can specify a previously declared dataset name (`DATASET DECLARE` command) instead of a file specification.

**COVB = 'savfile'|'dataset'**     *Writes the parameter covariance matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**CORB = 'savfile'|'dataset'**     *Writes the parameter correlation matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**MODEL = 'file'**     *Writes the parameter estimates and the parameter covariance matrix to an XML file.*

**PARAMETER = 'file'**     *Writes the parameter estimates to an XML file.*



# CSORDINAL

CSORDINAL is available in the Complex Samples option.

*Note:* Square brackets that are used in the CSORDINAL syntax chart are required parts of the syntax and are not used to indicate optional elements. Equals signs (=) that are used in the syntax chart are required elements. Except for the PLAN subcommand, all subcommands are optional.

```
CSORDINAL dependent varname ({ASCENDING**}) BY factor list
                        {DESCENDING }
      WITH covariate list

/PLAN FILE = 'file'

/JOINTPROB FILE = 'savfile' | 'dataset'

/MODEL effect-list

/LINK FUNCTION = {CAUCHIT}
                {CLOGLOG}
                {LOGIT**}
                {NLOGLOG}
                {PROBIT }

/CUSTOM LABEL = "label"

      LMATRIX = {list, effect list, effect list ...; ...}
                {list, effect list, effect list ... }
                {effect list, effect list ...; ... }
                {effect list, effect list ... }
                {ALL list; ALL ... }
                {ALL list }

      KMATRIX = {number; number; ...}
                {number }

/CUSTOM ...

/ODDSRATIOS {FACTOR = [varname ({LOW }) varname ...] }
            {HIGH**}
            {value }

            {COVARIATE = [varname ({1** }) varname ...]}
            {number list}

            CONTROL = [varname (value) varname (value) ...]

/ODDSRATIOS ...

/CRITERIA CHKSEP = {20**} CILEVEL = {95** } [DF = value]
                  {n } {value}

      LCONVERGE = [{0** } {RELATIVE**}] METHOD = {FISHER(n)}
                  {value} {ABSOLUTE } {NEWTON** }

      MXITER = {100**} MXSTEP = {5**}
               {n } {n }

      PCONVERGE = [{1E-6**} {RELATIVE**}] SINGULAR = {1E-12**}
                  {value } {ABSOLUTE } {value }

/STATISTICS PARAMETER EXP SE TTEST CINTERVAL DEFF DEFFSQRT

/NONPARALLEL TEST PARAMETER COVB

/TEST TYPE = {F** } PADJUST = {LSD** }
             {ADJF } {BONFERRONI }
```

*CSORDINAL*

```

                {CHISQUARE   }           {SEQBONFERRONI}
                {ADJCHISQUARE}         {SIDAK         }
                                           {SEQSIDAK      }

/DOMAIN VARIABLE = varname (value)

/MISSING CLASSMISSING = {EXCLUDE**}
                       {INCLUDE  }

/PRINT SAMPLEINFO** VARIABLEINFO** SUMMARY**
      HISTORY({1**}) GEF LMATRIX COVB CORB
              {n  }
      CLASSTABLE NONE

/SAVE CUMPROB(rootname:{25**}) PREDPROB(rootname:{25**})
      {n  } {n  }
      PREDVAL(varname) PREDVALPROB(varname) OBSVALPROB(varname)

/OUTFILE {COVB = 'savfile' | 'dataset'} {MODEL = 'file'   }
         {CORB = 'savfile' | 'dataset'} {PARAMETER = 'file'}

```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

**Release History**

Release 15.0

- Command introduced.

Release 17.0

- Added support for SET THREADS.

**Example**

```

CSORDINAL y BY a b c WITH x
  /PLAN FILE='/survey/myfile.csplan'.

```

**Overview**

CSORDINAL performs regression analysis on a binary or ordinal polytomous dependent variable using the selected cumulative link function for samples drawn by complex sampling methods. The procedure estimates variances by taking into account the sample design used to select the sample, including equal probability and probability proportional to size (PPS) methods and with replacement (WR) and without replacement (WOR) sampling procedures. Optionally, CSORDINAL performs analyses for a subpopulation.

**Basic Specification**

- The basic specification is a variable list identifying the dependent variable, the factors (if any), and the covariates (if any) and a PLAN subcommand with the name of a complex sample analysis plan file, which may be generated by the CSPLAN procedure.

- The default model includes threshold parameters, main effects for any factors, and any covariates.
- The basic specification displays summary information about the sample and all analysis variables, model summary statistics, and Wald  $F$  tests for all model effects. Additional subcommands must be used for other output.

### **Syntax Rules**

- The dependent variable and `PLAN` subcommand are required. All other variables and subcommands are optional.
- Multiple `CUSTOM` and `ODDSRATIOS` subcommands may be specified; each is treated independently. All other subcommands may be specified only once.
- Empty subcommands are not allowed; all subcommands must be specified with options.
- Each keyword may be specified only once within a subcommand.
- Subcommand names and keywords must be spelled in full.
- Equals signs (=) shown in the syntax chart are required.
- Square brackets shown in the syntax chart are required parts of the syntax and are not used to indicate optional elements. (See the `ODDSRATIOS` and `CRITERIA` subcommands.)
- Subcommands may be specified in any order.
- The dependent variable, factors, and the subpopulation variable can be numeric or string variables, but covariates must be numeric.
- Across the dependent, factor, and covariate variable lists, a variable may be specified only once.
- Plan file and subpopulation variables may not be specified on the variable list.
- Minimum syntax is a dependent variable and the `PLAN` subcommand. This specification fits a thresholds-only model.

### **Operations**

- `CSORDINAL` performs ordinal regression analysis for sampling designs supported by the `CSPLAN` and `CSSELECT` procedures.
- The input data set must contain the variables to be analyzed and variables related to the sampling design.
- The complex sample analysis plan file provides an analysis plan based on the sampling design.
- By default, `CSORDINAL` uses a model that includes thresholds, main effects for any factors, and any covariates.
- Other effects, including interaction and nested effects, may be specified using the `MODEL` subcommand.
- The default output for the specified model is summary information about the sample and all analysis variables, model summary statistics, and Wald  $F$  tests for all model effects.
- `WEIGHT` and `SPLIT FILE` settings are ignored by the `CSORDINAL` procedure.
- This procedure uses the multithreaded options specified by `SET THREADS`.

**Limitations**

- WEIGHT and SPLIT FILE settings are ignored with a warning by the CSORDINAL procedure.

**Examples**

```
* Complex Samples Ordinal Regression.
CSORDINAL opinion_gastax (ASCENDING) BY agecat gender votelast drivefreq
/PLAN FILE = 'samplesDirectory\poll.csplan'
/JOINTPROB FILE = 'samplesDirectory\poll_jointprob.sav'
/LINK FUNCTION=LOGIT
/MODEL agecat gender votelast drivefreq
/STATISTICS PARAMETER EXP SE CINTERVAL DEFF
/NONPARALLEL TEST PARAMETER
/TEST TYPE=ADJF PADJUST=SEQSIDAK
/ODDSRATIOS FACTOR=[agecat(HIGH)]
/ODDSRATIOS FACTOR=[drivefreq(3)]
/MISSING CLASSMISSING=EXCLUDE
/CRITERIA MXITER=100 MXSTEP=5
PCONVERGE=[1e-006 RELATIVE] LCONVERGE=[0]
METHOD=NEWTON CHKSEP=20
CILEVEL=95
/PRINT SUMMARY CLASSTABLE VARIABLEINFO SAMPLEINFO.
```

- The procedure builds a model for *opinion\_gastax* using *agecat*, *gender*, *votelast*, and *drivefreq* as factors.
- The complex sampling plan is located in *poll.csplan*, and the joint inclusion probabilities are in *poll\_jointprob.sav*.
- The model specifically calls for a main-effects model.
- Parameter estimates, their standard errors, 95% confidence intervals, and design effects are requested, along with exponentiated parameter estimates and their 95% confidence intervals.
- The test of parallel lines is requested, and the parameter estimates for the generalized cumulative model will be displayed.
- For all appropriate tests, the adjusted Wald *F* statistic will be computed, and *p* values for multiple comparisons will be adjusted according to the sequential Sidak method.
- Cumulative odds ratios are requested for *agecat*, with the highest level as the reference category, and *drivefreq*, with the third level as the reference category.
- A classification table is requested in addition to the default model output.
- All other options are set to their default values.

**Variable List**

The variable list specifies the dependent variable with the categories order, the factors, and the covariates in the model.

- The dependent variable must be the first specification on CSORDINAL.
- The dependent variable can be numeric or string.
- The CSORDINAL procedure sorts levels of the dependent variable in ascending or descending order. (If the dependent variable is a string variable, then the order is locale-dependent.)

- Sorting order for the values of the dependent variable may be specified in parentheses immediately following the dependent variable.

**ASCENDING**     *Sort dependent variable values in ascending order.* This is the default setting.

**DESCENDING**     *Sort dependent variable values in descending order.*

- The names of the factors and covariates, if any, follow the dependent variable. Specify any factors following the keyword `BY`. Specify any covariates following the keyword `WITH`.
- Factors can be numeric or string variables, but covariates must be numeric.
- Each variable may be specified only once on the variable list.
- Plan file and subpopulation variables may not be specified on the variable list.

## ***PLAN Subcommand***

The `PLAN` subcommand specifies the name of an XML file containing analysis design specifications. This file is written by the `CSPLAN` procedure.

- The `PLAN` subcommand is required.

**'FILE'**                     *Specifies the name of an external file.*

## ***JOINTPROB Subcommand***

The `JOINTPROB` subcommand is used to specify the file containing the first stage joint inclusion probabilities for `UNEQUAL_WOR` estimation. The `CSSELECT` procedure writes this file in the same location and with the same name (but different extension) as the plan file. When `UNEQUAL_WOR` estimation is specified, the `CSORDINAL` procedure will use the default location and name of the file unless the `JOINTPROB` subcommand is used to override them.

**'FILE' | 'dataset'**

*The name of the joint inclusion probabilities file. It can be an external file or an open dataset.*

## ***MODEL Subcommand***

The `MODEL` subcommand is used to specify the effects to be included in the model. Threshold parameters are included automatically. Their number is one less than the number of categories of the dependent variable found in the data.

- Specify a list of terms to be included in the model, separated by spaces or commas.
- If the `MODEL` subcommand is not specified, `CSORDINAL` uses a model that includes threshold parameters, main effects for any factors, and any covariates in the order specified on the variable list.
- To include a term for the main effect of a factor, enter the name of the factor.

- To include a term for an interaction among factors, use the keyword `BY` or the asterisk (\*) to join the factors involved in the interaction. For example, `A*B` means a two-way interaction effect of A and B, where A and B are factors. `A*A` is not allowed because factors inside an interaction effect must be distinct.
- To include a term for nesting one factor within another, use a pair of parentheses. For example, `A(B)` means that A is nested within B. `A(A)` is not allowed because factors inside a nested effect must be distinct.
- Multiple nesting is allowed. For example, `A(B(C))` means that B is nested within C, and A is nested within B(C). When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.
- Nesting within an interaction effect is valid. For example, `A(B*C)` means that A is nested within B\*C.
- Interactions among nested effects are allowed. The correct syntax is the interaction followed by the common nested effect inside the parentheses. For example, interaction between A and B within levels of C should be specified as `A*B(C)` instead of `A(C)*B(C)`.
- To include a covariate term in the design, enter the name of the covariate.
- Covariates can be connected, but not nested, through the \* operator or using the keyword `BY` to form another covariate effect. Interactions among covariates such as `X1*X1` and `X1*X2` are valid, but `X1(X2)` is not.
- Factor and covariate effects can be connected in various ways except that no effects can be nested within a covariate effect. Suppose A and B are factors and X1 and X2 are covariates. Examples of valid combinations of factor and covariate effects are `A*X1`, `A*B*X1`, `X1(A)`, `X1(A*B)`, `X1*A(B)`, `X1*X2(A*B)`, and `A*B*X1*X2`.

## ***LINK Subcommand***

The `LINK` subcommand offers the choice of a cumulative link function to specify the model.

- The keyword `FUNCTION`, followed by an equals sign, and a link function keyword are required.
- If the subcommand is not specified, `LOGIT` is the default cumulative link function.
- Only a single cumulative link function can be specified.

<b>CAUCHIT</b>	<i>Cauchit function.</i> $f(x)=\tan(\pi(x-0.5))$
<b>CLOGLOG</b>	<i>Complementary log-log function.</i> $f(x)=\log(-\log(1-x))$
<b>LOGIT</b>	<i>Logit function.</i> $f(x)=\log(x / (1-x))$ . This is the default link function.
<b>NLOGLOG</b>	<i>Negative log-log function.</i> $f(x)=-\log(-\log(x))$
<b>PROBIT</b>	<i>Probit function.</i> $f(x)=\Phi^{-1}(x)$ , where $\Phi^{-1}$ is the inverse standard normal cumulative distribution function.

## ***CUSTOM Subcommand***

The `CUSTOM` subcommand defines custom hypothesis tests by specifying the **L** matrix (contrast coefficients matrix) and the **K** matrix (contrast results matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ . The vector **B** is the parameter vector in the cumulative link model.

For a binary dependent variable, CSORDINAL models a single threshold parameter and a set of regression parameters. For a polytomous ordinal dependent variable with  $K$  levels, CSORDINAL models a threshold parameter for each category except the last and a single set of regression parameters for all response categories. The CUSTOM subcommand allows you to specify an **L** matrix with contrast coefficients for all thresholds and regression parameters.

- Multiple CUSTOM subcommands are allowed. Each is treated independently.
- An optional label may be specified using the LABEL keyword. The label is a string with a maximum length of 255 characters. Only one label can be specified.
- The **L** matrix is the contrast coefficients matrix. This matrix specifies coefficients of contrasts, which can be used for studying the effects in the model. An **L** matrix must always be specified using the LMATRIX keyword.
- The **K** matrix is the contrast results matrix. This matrix specifies the results of the linear hypothesis. A **K** matrix can be specified using the KMATRIX keyword.
- The number of rows in the **L** and **K** matrices must be equal.
- The default **K** matrix is a zero matrix; that is,  $\mathbf{LB} = \mathbf{0}$  is assumed.
- There are three general formats that can be used on the LMATRIX keyword: (1) Specify coefficient values for thresholds, followed optionally by an effect name and a list of real numbers. (2) Specify an effect name and a list of real numbers. (3) Specify the keyword ALL and a list of real numbers. In all three formats, there can be multiple effect names (or instances of the keyword ALL) and number lists.
- When specifying threshold coefficients in the first or the third general format, a complete list of  $K-1$  coefficient values must be given in the increasing threshold order.
- Only valid effects in the default model or on the MODEL subcommand can be specified on the LMATRIX keyword.
- The length of the list of real numbers on the LMATRIX keyword must be equal to the number of parameters (including the redundant ones) corresponding to the specified effect. For example, if the effect A\*B takes up six columns in the design matrix, then the list after A\*B must contain exactly six numbers.
- When ALL is specified, the length of the list that follows ALL must be equal to the total number of parameters (including the redundant ones) in the model. For a binary dependent variable, the contrast coefficients for the single threshold and all regression parameters must be listed following the ALL keyword. For a polytomous dependent variable with  $K$  levels, the contrast coefficients for the  $K-1$  thresholds and all regression parameters must be listed in order following the ALL keyword.
- Effects that are in the model but not specified on the LMATRIX keyword are assumed to have entries of 0 in the corresponding columns of the **L** matrix.
- When defining an **L** matrix, a number can be specified as a fraction with a positive denominator—for example, 1/3 and -1/3 are valid, but 1/-3 is invalid.
- A semicolon (;) indicates the end of a row in the **L** matrix.

- The format for the `KMATRIX` keyword is one or more real numbers. If more than one number is specified, then separate adjacent numbers using a semicolon (;). Each semicolon indicates the end of a row in the **K** matrix. Each number is the hypothesized value for a contrast, which is defined by a row in the **L** matrix.
- If rows of the **L** matrix are not independent, a submatrix of **L** with independent rows is used for testing. Tested rows are indicated when the **K** matrix is not a zero matrix.

### Example

Suppose that factors A and B each have three levels.

```
CSORDINAL y BY a b
/PLAN FILE='/survey/myfile.csplan'
/MODEL a b a*b
/CUSTOM LABEL = 'Effect A'
LMATRIX = a 1 0 -1
          a*b 1/3 1/3 1/3
              0 0 0
              -1/3 -1/3 -1/3;
a 0 1 -1
a*b 0 0 0
    1/3 1/3 1/3
    -1/3 -1/3 -1/3.
```

- The preceding syntax specifies a test of effect A.
- Because there are three levels in effect A, at most two independent contrasts can be formed; thus, there are two rows in the **L** matrix, separated by a semicolon (;).
- There are three levels each in effects A and B; thus, the interaction effect A\*B takes nine columns in the design matrix.
- The first row in the **L** matrix tests the difference between levels 1 and 3 of effect A; the second row tests the difference between levels 2 and 3 of effect A.
- The `KMATRIX` keyword is not specified, so the null hypothesis value for both tests is 0.

### Example

Suppose that dependent variable Z and factor A each have three levels.

```
CSORDINAL z BY a
/PLAN FILE='/survey/myfile.csplan'
/MODEL a
/CUSTOM LABEL = 'Effect A'
LMATRIX = a 1 0 -1;
          a 0 1 -1
KMATRIX = 1; 1.
```

- The dependent variable Z has three categories, so there will be two thresholds.
- The syntax specifies a model with thresholds and a main effect for factor A, and a custom hypothesis test of effect A.
- Because the `ALL` option is not used on the `LMATRIX` keyword, threshold coefficients are set to zero. The equivalent `LMATRIX` keyword using the `ALL` option follows.



```
LMATRIX = ALL 0 0 1 0 -1;
          ALL 0 0 0 1 -1
```

- The `KMATRIX` keyword is specified and the hypothesis that the difference between levels 1 and 3 and levels 2 and 3 of effect A are both equal to 1 is tested.

## ***ODDSRATIOS Subcommand***

The `ODDSRATIOS` subcommand estimates cumulative odds ratios for the specified factor(s) or covariate(s). The subcommand is available only for `LOGIT` link. For other link functions, the subcommand is ignored and a warning is issued. Note that these cumulative odds ratios are model-based and are not directly computed using the observed data.

A single cumulative odds ratio is computed for all categories of the dependent variable except the last; the proportional odds model postulates that they are all equal.

If the `FACTOR` keyword is specified, the cumulative odds ratios compare the cumulative odds at each factor category  $j$  with the cumulative odds at category  $J$ , where  $J$  is the reference category defined in parentheses following the variable name of the factor. All other factors and covariates are fixed as defined on the `CONTROL` keyword.

If the `COVARIATE` keyword is specified, the cumulative odds ratios compare the cumulative odds at value  $x$  with the cumulative odds at value  $x + \Delta x$ , where  $\Delta x$  is the change in  $x$  defined in parentheses following the variable name of the covariate. To define the value  $x$ , specify the covariate and the value on the `CONTROL` keyword. The value of all other factors and covariates are fixed as defined on the `CONTROL` keyword also.

If a specified factor or covariate interacts with other predictors in the model, then the cumulative odds ratios depend not only on the change in the specified variable but also on the values of the variables with which it interacts. If a specified covariate interacts with itself in the model (for example, `X*X`), then the cumulative odds ratios depend on both the change in the covariate and the value of the covariate. The values of interacting factors and covariates can be customized using the `CONTROL` keyword.

The `CSORDINAL` procedure sorts levels of each factor in ascending order and defines the highest level as the last level. (If the factor is a string variable, then the value of the highest level is locale-dependent.)

- Multiple `ODDSRATIOS` subcommands are allowed. Each is treated independently.
- Either the `FACTOR` keyword and one or more factors, or the `COVARIATE` keyword and one or more covariates, but not both, are required. All other keywords are optional.

- The `FACTOR`, `COVARIATE`, and `CONTROL` keywords must be followed by an equals sign and one or more elements enclosed in square brackets.
- If a variable is specified on the `FACTOR` keyword and is also specified on the `CONTROL` keyword, then the `CONTROL` specification for that variable is ignored when the variable's odds ratios are computed. Thus, `FACTOR = [A B] CONTROL = [A(1) B(2)]` estimates odds ratios for factor A holding factor B at level 2, and for factor B holding factor A at level 1.

**FACTOR = [option]**

Valid options are one or more factors appearing on the factor list. Optionally, each factor may be followed by parentheses containing the level to use as the reference category when computing cumulative odds ratios. The keyword `LOW` or `HIGH`, or a value, may be specified. Put the value inside a pair of quotes if the value is formatted (such as date or currency) or if the factor is of string type. By default, the highest category is used as the reference category.

If a value is specified but the value does not exist in the data, then a warning is issued and the default `HIGH` is used.

Any factor may occur only once on the `FACTOR` keyword.

**COVARIATE = [option]**

Valid options are one or more covariates appearing on the covariate list. Optionally, each covariate may be followed by parentheses containing one or more nonzero numbers giving unit(s) of change to use for covariates when computing cumulative odds ratios. Cumulative odds ratios are estimated for each distinct value. The default value is 1.

Any covariate may occur only once on the `COVARIATE` keyword.

**CONTROL = [option]**

Specifies the factor and/or covariate values to use when computing cumulative odds ratios. Factors must appear on the factor list, and covariates on the covariate list, of the `CSORDINAL` command.

Factors must be followed by the keyword `LOW` or `HIGH`, or a value, in parentheses. Put the value inside a pair of quotes if the value is formatted (such as date or currency) or if the factor is of string type. If keyword `LOW` or `HIGH` is used, then each cumulative odds ratio is computed by holding the factor at its lowest or highest level, respectively. If a value is used, then each cumulative odds ratio is computed by holding the specified factor at the supplied value. If a factor is not specified on the `CONTROL` option, then its highest category is used in cumulative odds ratio calculations.

If a factor value is specified but the value does not exist in the data, then a warning is issued and the default `HIGH` is used.

Covariates must be followed by keyword `MEAN` or a number in parentheses. If keyword `MEAN` is used, then each cumulative odds ratio is computed by holding the covariate at its overall mean. If a number is used, then each cumulative odds ratio is computed by holding the specified covariate at the supplied value. If a covariate is not specified on the `CONTROL` option, then its overall mean is used in cumulative odds ratio calculations.

Any factor or covariate may occur only once on the `CONTROL` keyword.

**Example**

Suppose that dependent variable Y has three levels; factor A has two levels; and factor B has three levels coded 1, 2, and 3.

```
CSORDINAL y BY a b WITH x
  /PLAN FILE='/survey/myfile.csplan'
```

```

/MODEL a b a*b x
/ODDSRATIOS FACTOR=[a] CONTROL=[b(1)]
/ODDSRATIOS FACTOR=[a] CONTROL=[b(2)]
/ODDSRATIOS FACTOR=[a] CONTROL=[b(3)].

```

- The default LOGIT cumulative link function is used and the cumulative odds ratios are computed. They are equal across all response levels by the model definition.
- The model includes two thresholds, main effects for factors A and B, the A\*B interaction effect, and the covariate X.
- Cumulative odds ratios are requested for factor A. Assuming the A\*B interaction effect is significant, the cumulative odds ratio for factor A will differ across levels of factor B. The specified syntax requests three cumulative odds ratios for factor A; each is computed at a different level of factor B.

### Example

```

CSORDINAL z BY a b c WITH x y
/PLAN FILE='/survey/myfile.csplan'
/MODEL a b c x*y
/ODDSRATIOS COVARIATE=[x(1 3 5)] CONTROL=[y(1)].

```

- The preceding syntax will compute three cumulative odds ratios for covariate X.
- The parenthesized list following variable X provides the unit of change values to use when computing cumulative odds ratios. Cumulative odds ratios will be computed for X increasing by 1, 3, and 5 units and holding covariate Y equal to 1.

## CRITERIA Subcommand

The CRITERIA subcommand offers controls on the iterative algorithm used for estimation, and specifies numerical tolerance for checking singularity.

### CHKSEP = integer

*Starting iteration for checking complete and quasi-complete separation.* Specify a non-negative integer. This criterion is not used if the value is 0. The default value is 20.

### CILEVEL = value

*Confidence interval level for coefficient estimates, exponentiated coefficient estimates, and cumulative odds ratio estimates.* Specify a value greater than or equal to 0, and less than 100. The default value is 95.

### DF = value

*Sampling design degrees of freedom to use in computing p values for all test statistics.* Specify a positive number. The default value is the difference between the number of primary sampling units and the number of strata in the first stage of sampling.

### LCONVERGE = [number (RELATIVE | ABSOLUTE)]

*Log-likelihood function convergence criterion.* Convergence is assumed if the relative or absolute change in the log-likelihood function is less than the given value. This criterion is not used if the value is 0.

Specify square brackets containing a non-negative number followed optionally by the keyword RELATIVE or ABSOLUTE, which indicates the type of change. The default value is 0; the default type is RELATIVE.

**METHOD = FISHER(number) | NEWTON**

*Model parameters estimation method.* The Fisher scoring method is specified by the keyword `FISHER`, the Newton-Raphson method, by the keyword `NEWTON`, and a hybrid method is available by specifying `FISHER(n)`. In the hybrid method,  $n$  is the maximal number of Fisher scoring iterations before switching to the Newton-Raphson method. If convergence is achieved during the Fisher scoring phase of the hybrid method, iterations continue with the Newton-Raphson method.

**MXITER = integer**

*Maximum number of iterations.* Specify a non-negative integer. The default value is 100.

**MXSTEP = integer**

*Maximum step-halving allowed.* Specify a positive integer. The default value is 5.

**PCONVERGE = [number (RELATIVE | ABSOLUTE)]**

*Parameter estimates convergence criterion.* Convergence is assumed if the relative or absolute change in the parameter estimates is less than the given value. This criterion is not used if the value is 0.

Specify square brackets containing a non-negative number followed optionally by the keyword `RELATIVE` or `ABSOLUTE`, which indicates the type of change. The default value is  $10^{-6}$ ; the default type is `RELATIVE`.

**SINGULAR = value**

*Tolerance value used to test for singularity.* Specify a positive value. The default value is  $10^{-12}$ .

## **STATISTICS Subcommand**

The `STATISTICS` subcommand requests various statistics associated with the parameter estimates.

- There are no default keywords on the `STATISTICS` subcommand. If this subcommand is not specified, then none of the statistics listed below are displayed

<b>PARAMETER</b>	<i>Parameter estimates.</i>
<b>EXP</b>	<i>The exponentiated parameter estimates.</i> It is available only for the <code>LOGIT</code> link.
<b>SE</b>	<i>Standard error for each parameter estimate.</i>
<b>TTEST</b>	<i>t test for each parameter estimate.</i>
<b>CINTERVAL</b>	<i>Confidence interval for each parameter estimate and/or exponentiated parameter estimate.</i>
<b>DEFF</b>	<i>Design effect for each parameter estimate.</i>
<b>DEFFSQRT</b>	<i>Square root of design effect for each parameter estimate.</i>

## **NONPARALLEL Subcommand**

The `NONPARALLEL` subcommand requests various statistics associated with a general cumulative link model with non-parallel lines where a separate regression line is fitted for each response category except for the last.

<b>TEST</b>	<i>Test of parallel lines assumption.</i> Test whether regression parameters are equal for all cumulative responses. The general model with non-parallel lines is estimated and the Wald test of equal parameters is applied.
<b>PARAMETER</b>	<i>Parameters of the general model with non-parallel lines.</i> The general model is estimated using the same convergence criteria as for the original model. Both parameters and their standard errors are estimated.
<b>COVB</b>	<i>Covariance matrix for the general model parameters.</i>

## **TEST Subcommand**

The `TEST` subcommand specifies the type of test statistic and the method of adjusting the significance level to be used for hypothesis tests requested on the `MODEL`, `CUSTOM`, and `PRINT` subcommands.

### **TYPE Keyword**

The `TYPE` keyword indicates the type of test statistic.

<b>F</b>	<i>Wald F test.</i> This is the default test statistic if the <code>TYPE</code> keyword is not specified.
<b>ADJF</b>	<i>Adjusted Wald F test.</i>
<b>CHISQUARE</b>	<i>Wald chi-square test.</i>
<b>ADJCHISQUARE</b>	<i>Adjusted Wald chi-square test.</i>

### **PADJUST Keyword**

The `PADJUST` keyword indicates the method of adjusting the significance level.

<b>LSD</b>	<i>Least significant difference.</i> This method does not control the overall probability of rejecting the hypotheses that some linear contrasts are different from the null hypothesis value(s). This is the default.
<b>BONFERRONI</b>	<i>Bonferroni.</i> This method adjusts the observed significance level for the fact that multiple contrasts are being tested.
<b>SEQBONFERRONI</b>	<i>Sequential Bonferroni.</i> This is a sequentially step-down rejective Bonferroni procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.
<b>SIDAK</b>	<i>Sidak.</i> This method provides tighter bounds than the Bonferroni approach.
<b>SEQSIDAK</b>	<i>Sequential Sidak.</i> This is a sequentially step-down rejective Sidak procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.

## ***DOMAIN Subcommand***

The `DOMAIN` subcommand specifies the subpopulation for which the analysis is to be performed.

- The keyword `VARIABLE`, followed by an equals sign, a variable, and a value in parentheses, are required. Put the value inside a pair of quotes if the value is formatted (such as date or currency) or if the factor is of string type.
- The subpopulation is defined by all cases having the given value on the specified variable.
- Analyses are performed only for the specified subpopulation.
- For example, `DOMAIN VARIABLE = myvar (1)` defines the subpopulation by all cases for which variable `MYVAR` has value 1.
- The specified variable may be numeric or string and must exist at the time the `CSORDINAL` procedure is invoked.
- Stratification or cluster variables may be specified, but no other plan file variables are allowed on the `DOMAIN` subcommand.
- Analysis variables may not be specified on the `DOMAIN` subcommand.

## ***MISSING Subcommand***

The `MISSING` subcommand specifies how missing values are handled.

- In general, cases must have valid data for all design variables as well as for the dependent variable and any covariates. Cases with invalid data for any of these variables are excluded from the analysis.
- There is one important exception to the preceding rule. This exception applies when an inclusion probability or population size variable is defined in an analysis plan file. Within a stratum at a given stage, if the inclusion probability or population size values are unequal across cases or missing for a case, then the first valid value found within that stratum is used as the value for the stratum. If strata are not defined, then the first valid value found in the sample is used. If the inclusion probability or population size values are missing for all cases within a stratum (or within the sample if strata are not defined) at a given stage, then an error message is issued.
- The `CLASSMISSING` keyword specifies whether user-missing values are treated as valid. This specification is applied to categorical design variables (that is, strata, cluster, and subpopulation variables), the dependent variable, and any factors.

**EXCLUDE**      *Exclude user-missing values among the strata, cluster, subpopulation, the dependent variable, and factor variables. This is the default.*

**INCLUDE**      *Include user-missing values among the strata, cluster, subpopulation, the dependent variable, and factor variables. Treat user-missing values for these variables as valid data.*

## ***PRINT Subcommand***

The `PRINT` subcommand is used to display optional output.

- If the `PRINT` subcommand is not specified, then the default output includes sample information, variable and factor information, and model summary statistics.
- If the `PRINT` subcommand is specified, then `CSORDINAL` displays output only for those keywords that are specified.

**SAMPLEINFO** *Sample information table.* Displays summary information about the sample, including the unweighted count and the population size. This is default output if the `PRINT` subcommand is not specified.

**VARIABLEINFO**

*Variable information.* Displays summary information about the dependent variable, covariates, and factors. This is default output if the `PRINT` subcommand is not specified.

**SUMMARY** *Model summary statistics.* Displays pseudo- $R^2$  statistics. This is default output if the `PRINT` subcommand is not specified.

**HISTORY(n)** *Iteration history.* Displays coefficient estimates and statistics at every  $n^{\text{th}}$  iteration beginning with the  $0^{\text{th}}$  iteration (the initial estimates). The default is to print every iteration ( $n = 1$ ). The last iteration is always printed if `HISTORY` is specified, regardless of the value of  $n$ .

**GEF** *General estimable function table.*

**LMATRIX** *Set of contrast coefficients (L) matrices.* These are the Type III contrast matrices used in testing model effects.

**COVB** *Covariance matrix for model parameters.*

**CORB** *Correlation matrix for model parameters.*

**CLASSTABLE** *Classification table.* Displays frequencies of observed versus predicted response categories.

**NONE** *No PRINT subcommand output.* None of the `PRINT` subcommand output is displayed. However, if `NONE` is specified with one or more other keywords, then the other keywords override `NONE`.

## **SAVE Subcommand**

The `SAVE` subcommand writes optional model variables to the active dataset.

- Specify one or more temporary variables, each followed by an optional new name in parentheses.
- The optional names must be valid variable names.
- If new names are not specified, `CSORDINAL` uses the default names.
- If a subpopulation is defined on the `DOMAIN` subcommand, then `SAVE` applies only to cases within the subpopulation.

The following rules describe the functionality of the `SAVE` subcommand in relation to the predictor values for each case.

- If all factors and covariates in the model have valid values for the case, then the procedure computes the predicted values. (The `MISSING` subcommand setting is taken into account when defining valid/invalid values for a factor.)
- An additional restriction for factors is that only those values of the factor actually used in building the model are considered valid. For example, suppose factor A takes values 1, 2, and 3 when the procedure builds the model. Also suppose there is a case with a value of 4

on factor A, and valid values on all other factors and covariates. For this case, no predicted values are saved because there is no model coefficient corresponding to factor A = 4.

Computation of predicted values for a given case does not depend on the value of the dependent variable; it could be missing.

**CUMPROB (rootname:n)**

*Cumulative probability.* The user-specified or default name is treated as the root name, and a suffix is added to get new unique variable names. The root name can be followed by a colon and a positive integer giving the number of predicted cumulative probabilities to save. The predicted cumulative probabilities of the first  $n$  response categories are saved. One cumulative predicted probability variable can be saved for each category of the dependent variable. The default root name is *CumulativeProbability*. The default  $n$  is 25. To specify  $n$  without a root name, enter a colon before the number.

**PREDPROB (rootname:n)**

*Predicted probability.* The user-specified or default name is treated as the root name, and a suffix is added to get new unique variable names. The root name can be followed by a colon and a positive integer giving the number of predicted probabilities to save. The predicted probabilities of the first  $n$  response categories are saved. One predicted probability variable can be saved for each category of the dependent variable. The default root name is *PredictedProbability*. The default  $n$  is 25. To specify  $n$  without a root name, enter a colon before the number.

**PREDVAL (varname)**

*Predicted value.* The class or value predicted by the model. The optional variable name must be unique. If the default name is used and it conflicts with existing variable names, then a suffix is added to the default name to make it unique. The default variable name is *PredictedValue*.

**PREDVALPROB (varname)**

*Predicted value probability.* The probability of value predicted by the model. This probability is the maximum probability predicted by the model for a given case. The optional variable name must be unique. If the default name is used and it conflicts with existing variable names, then a suffix is added to the default name to make it unique. The default variable name is *PredictedValueProbability*.

**OBSVALPROB (varname)**

*Observed value probability.* The probability predicted for the observed response value. The optional variable name must be unique. If the default name is used and it conflicts with existing variable names, then a suffix is added to the default name to make it unique. The default variable name is *ObservedValueProbability*.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand saves an external IBM® SPSS® Statistics data file or previously declared dataset containing the parameter covariance or correlation matrix with parameter estimates, standard errors, significance values, and sampling design degrees of freedom. It also saves the parameter estimates and the parameter covariance matrix in XML format.

- At least one keyword and a filename are required.



- The COVB and CORB keywords are mutually exclusive, as are the MODEL and PARAMETER keywords.
- The filename must be specified in full. CSORDINAL does not supply an extension.

**COVB = 'savfile' | 'dataset'**

*Writes the parameter covariance matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**CORB = 'savfile' | 'dataset'**

*Writes the parameter correlation matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**MODEL = 'file'**

*Writes the parameter estimates and the parameter covariance matrix to an XML file.*

**PARAMETER = 'file'**

*Writes the parameter estimates to an XML file.*

# CSPLAN

CSPLAN is available in the Complex Samples option.

```
CSPLAN SAMPLE

/PLAN FILE=file

[/PLANVARS [SAMPLEWEIGHT=varname]]
          [PREVIOUSWEIGHT=varname]

[/PRINT [PLAN**] [MATRIX]]
```

## Design Block: Stage 1

```
/DESIGN [STAGELABEL='label']
        [STRATA=varname [varname [...] ] ]
        [CLUSTER=varname [varname [...] ] ]

/METHOD TYPE={SIMPLE_WOR      } [ESTIMATION={DEFAULT**}]
              {SIMPLE_WR      } {WR          }
              {SIMPLE_SYSTEMATIC}
              {SIMPLE_CHROMY   }
              {PPS_WOR        }
              {PPS_WR         }
              {PPS_SYSTEMATIC }
              {PPS_BREWER     }
              {PPS_MURTHY    }
              {PPS_SAMPFORD  }
              {PPS_CHROMY    }

[/MOS      {VARIABLE=varname} [MIN=value] [MAX=value] ]
          {SOURCE=FROMDATA }

[/SIZE {VALUE=sizevalue           }
      {VARIABLE=varname           }
      {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}

[/RATE {VALUE=ratevalue           }
      {VARIABLE=varname           }
      {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}
      [MINSIZE=value]
      [MAXSIZE=value]

[/STAGEVARS [INCLPROB[(varname)]]]
           [CUMWEIGHT[(varname)]]
           [INDEX[(varname)]]
           [POPSIZE[(varname)]]
           [SAMPSIZE[(varname)]]
           [RATE[(varname)]]
           [WEIGHT[(varname)]]
```

## Design Block: Stages 2 and 3

```
/DESIGN [STAGELABEL='label']
        [STRATA=varname [varname [...] ] ]
        [CLUSTER=varname [varname [...] ] ]

/METHOD TYPE={SIMPLE_WOR      }
              {SIMPLE_WR      }
              {SIMPLE_SYSTEMATIC}
              {SIMPLE_CHROMY   }

[/SIZE {VALUE=sizevalue           }
      {VARIABLE=varname           }
      {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}
      ]]
```

```

        {VARIABLE=varname
        {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}
[/RATE {VALUE=ratevalue
        {VARIABLE=varname
        {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}
        [MINSIZE=value]
        [MAXSIZE=value]

[/STAGEVARS [INCLPROB[ (varname)]]]
           [CUMWEIGHT[ (varname)]]
           [INDEX[ (varname)]]
           [POPSIZE[ (varname)]]
           [SAMPSIZE[ (varname)]]
           [RATE[ (varname)]]
           [WEIGHT[ (varname)]]

```

### Create an Analysis Design

```

CSPLAN ANALYSIS
/PLAN FILE=file
/PLANVARS ANALYSISWEIGHT=varname
[/SRSESTIMATOR TYPE={WOR**}
                {WR   }
[/PRINT [PLAN**] [MATRIX]]

```

### Design Block: Stage 1

```

/DESIGN [STAGELABEL='label']
        [STRATA=varname [varname [...] ] ]
        [CLUSTER=varname [varname [...] ] ]

/ESTIMATOR TYPE= {EQUAL_WOR }
                {UNEQUAL_WOR}
                {WR         }

[/POPSIZE {VALUE=sizevalue
          {VARIABLE=varname
          {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}

[/INCLPROB {VALUE=probvalue
           {VARIABLE=varname
           {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}

```

### Design Block: Stages 2 and 3

```

/DESIGN [STAGELABEL='label']
        [STRATA=varname [varname [...] ] ]
        [CLUSTER=varname [varname [...] ] ]

/ESTIMATOR TYPE= {EQUAL_WOR}
                {WR         }

[/POPSIZE {VALUE=sizevalue
          {VARIABLE=varname
          {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}

[/INCLPROB {VALUE=probvalue
           {VARIABLE=varname
           {MATRIX=varname [varname [...] ] ; catlist value [;catlist value [;...]]}

```

**Display an Existing Plan**

```
CSPLAN VIEW
/PLAN FILE=file
[/PRINT [PLAN**] [MATRIX]]
```

\*\* Default if the subcommand is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

**Example**

```
CSPLAN SAMPLE
/PLAN FILE= '/survey/myfile.csplan'
/DESIGN STRATA=region CLUSTER=school
/METHOD TYPE=PPS_WOR
/MOS VARIABLE=mysizevar
/SIZE VALUE=100.

CSPLAN ANALYSIS
/PLAN FILE= '/survey/myfile.csplan'
/PLANVARS ANALYSISWEIGHT=sampleweight
/DESIGN CLUSTER=district
/ESTIMATOR TYPE=UNEQUAL_WOR
/DESIGN CLUSTER=school
/ESTIMATOR TYPE=EQUAL_WOR
/INCLPROB VARIABLE=sprob.

CSPLAN VIEW
/PLAN FILE= '/survey/myfile.csplan'.
```

**Overview**

CSPLAN creates a complex sample design or analysis specification that is used by companion procedures in the Complex Samples option. CSSELECT uses specifications from a plan file when selecting cases from the active file. Analysis procedures in the Complex Samples option, such as CSDESCRIPTIVES, require a plan file in order to produce summary statistics for a complex sample. You can also use CSPLAN to view sample or analysis specifications within an existing plan file.

The CSPLAN design specification is used only by procedures in the Complex Samples option.

**Options**

**Design Specification.** CSPLAN writes a sample or analysis design to a file. A sample design can be used to extract sampling units from the active file. An analysis design is used to analyze a complex sample. When a sample design is created, the procedure automatically saves an appropriate analysis design to the plan file. Thus, a plan file created for designing a sample can be used for both sample selection and analysis.

Both sample and analysis designs can specify **stratification**, or independent sampling within nonoverlapping groups, as well as **cluster sampling**, in which groups of sampling units are selected. A single or multistage design can be specified with a maximum of three stages.

CSPLAN does not actually execute the plan (that is, it does not extract the sample or analyze data). To sample cases, use a sample design created by CSPLAN as input to CSSELECT. To analyze sample data, use an analysis design created by CSPLAN as input to Complex Samples procedures, such as CSDSCRIPTIVES.

**Sample Design.** A variety of equal- and unequal-probability methods are available for sample selection, including simple and systematic random sampling. CSPLAN offers several methods for sampling with probability proportionate to size (**PPS**), including Brewer's method, Murthy's method, and Sampford's method. Units can be drawn with replacement (**WR**) or without replacement (**WOR**) from the population. At each stage of the design, you can control the number or percentage of units to be drawn. You can also choose output variables, such as stagewise sampling weights, that are created when the sample design is executed.

**Analysis Design.** The following estimation methods are available: with replacement, equal probability without replacement, and unequal probability without replacement. Unequal probability estimation without replacement can be requested in the first stage only. You can specify variables to be used as input to the estimation process, such as overall sample weights and inclusion probabilities.

### **Operations**

- If a sample design is created, the procedure automatically writes a suitable analysis design to the plan file. The default analysis design specifies stratification variables and cluster variables for each stage, as well as an estimation method appropriate for the chosen extraction method.
- CSPLAN writes design specifications in XML format.
- By default, CSPLAN displays output that summarizes the sample or analysis design.

### **Subcommand Order**

- The first DESIGN subcommand must precede all other subcommands except PLAN, PLANVARS, and PRINT.
- PLAN, PLANVARS, and PRINT subcommands can be used in any order.

### **Limitations**

- A maximum of three design blocks can be specified.
- CSPLAN ignores SPLIT FILE and WEIGHT commands with a warning.

## **Basic Specification**

You can specify a sample or analysis design to be created or a plan file to be displayed.

### **Creating a Sample Plan**

- The SAMPLE keyword must be specified on the CSPLAN command.
- A PLAN subcommand is required that specifies a file that will contain the design specification.
- A DESIGN subcommand is required.
- A METHOD subcommand must specify an extraction method.

- Sample size or rate must be specified unless the PPS\_MURTHY or PPS\_BREWER extraction method is chosen.

#### ***Creating an Analysis Plan***

- The ANALYSIS keyword must be specified on the CSPLAN command.
- A PLAN subcommand is required that specifies a file that will contain the analysis specification.
- A PLANVARS subcommand is required that specifies a sample weight variable.
- A DESIGN subcommand is required.
- An ESTIMATOR subcommand must specify an estimator.
- The POPSIZE or INCLPROB subcommand must be specified if the EQUAL\_WOR estimator is selected.

#### ***Displaying an Existing Plan***

- The VIEW keyword must be specified on the CSPLAN command.
- A PLAN subcommand is required that specifies a file whose specifications are to be displayed.

## ***Syntax Rules***

### ***General***

- PLAN, PLANVARS, and PRINT are global. Only a single instance of each global subcommand is allowed.
- Within a subcommand, an error occurs if a keyword or attribute is specified more than once.
- Equals signs shown in the syntax chart are required.
- Subcommand names and keywords (for example, PPS\_WR) must be spelled in full.
- In general, empty subcommands (that is, those that have no specifications) generate an error. DESIGN is the only subcommand that can be empty.
- Any variable names that are specified must be valid IBM® SPSS® Statistics variable names.

### ***Creating a Plan***

- Stages are specified in design blocks. The DESIGN subcommand signals the start of a block. The first block corresponds to stage 1, the second to stage 2, and the third to stage 3. One DESIGN subcommand must be specified per stage.
- The following subcommands are local and apply to the immediately preceding DESIGN subcommand: METHOD, MOS, SIZE, RATE, STAGEVARS, ESTIMATOR, POPSIZE, and INCLPROB. An error occurs if any of these subcommands appears more than once within a block.
- Available METHOD and ESTIMATOR options depend on the stage.
- The following subcommands are honored only if a sample design is requested: METHOD, MOS, SIZE, RATE, and STAGEVARS. An error occurs if any of these subcommands is specified for an analysis design.
- MOS can be specified in stage 1 only.

- The following subcommands can be used only if an analysis design is requested: `ESTIMATOR`, `POPSIZE`, and `INCLPROB`. An error occurs if any of these subcommands is specified for a sample design.
- In general, each variable specified in the design can assume only one role. For example, a weight variable cannot be used as a stratification or cluster variable. Exceptions are listed below.

### **Displaying a Plan**

- If `CSPLAN VIEW` is used, only the `PLAN` and `PRINT` subcommands can be specified.

## **Examples**

### **Simple Sample Design**

```
CSPLAN SAMPLE
  /PLAN FILE='/survey/myfile.csplan'
  /DESIGN
  /METHOD TYPE=SIMPLE_WOR
  /SIZE VALUE=100.
```

- A single-stage sample design is created that is saved in *myfile.csplan*.
- One hundred cases will be selected from the active file when the sample design is executed by the `CSSELECT` procedure.
- The extraction method is simple random sampling without replacement.
- The plan file also includes a default analysis design that uses the `EQUAL_WOR` estimator (the default when units are extracted using the `SIMPLE_WOR` method).

### **Stratified Sample Design**

```
CSPLAN SAMPLE
  /PLAN FILE='/survey/myfile.csplan'
  /DESIGN STRATA=region
  /METHOD TYPE=SIMPLE_WOR
  /RATE MATRIX=REGION; 'East' 0.1 ; 'West' 0.2;
                        'North' 0.1; 'South' 0.3.
```

- A stratified sample design is specified with disproportionate sampling rates for the strata. Sample elements will be drawn independently within each region.
- The extraction method is simple random sampling without replacement.
- `CSPLAN` generates a default analysis design using *region* as a stratification variable and the `EQUAL_WOR` estimator.

### **Stratified Cluster Sample Design**

```
CSPLAN SAMPLE
  /PLAN FILE='/survey/myfile.csplan'
  /DESIGN STRATA=region CLUSTER=school
  /METHOD TYPE=PPS_WOR
  /SIZE VALUE=10
  /MOS VARIABLE=mysizevar.
```

- A stratified cluster sample design is specified.
- Ten schools will be selected within each region with probability proportionate to size.
- Size values for the strata are read from *mysizevar*.
- CSPLAN generates a default analysis design using *region* as a stratification variable and *school* as a cluster variable.
- The UNEQUAL\_WOR estimator will be used for analysis. UNEQUAL\_WOR is the default when units are sampled with probability proportionate to size.

### **Multistage Cluster Sample Design**

```
CSPLAN SAMPLE
  /PLAN FILE='/survey/myfile.csplan'
  /DESIGN STAGELABEL='school districts' CLUSTER=district
  /METHOD TYPE=PPS_WOR
  /RATE VALUE=.2
  /MOS VARIABLE=districtsize
  /DESIGN STAGELABEL='schools' CLUSTER=school
  /METHOD TYPE=SIMPLE_WOR
  /RATE VALUE=0.3.
```

- A multistage cluster sample design is specified.
- Twenty percent of school districts will be drawn with probability proportionate to size.
- Within each selected school district, 30% of schools will be drawn without replacement.
- CSPLAN generates a default analysis design. Since the PPS\_WOR sampling method is specified in stage 1, the UNEQUAL\_WOR estimator will be used for analysis for that stage. The EQUAL\_WOR method will be used to analyze stage 2.

### **Simple Analysis Design**

```
CSPLAN ANALYSIS
  /PLAN FILE='/survey/myfile.csaplan'
  /PLANVARS ANALYSISWEIGHT=sampleweight
  /DESIGN
  /ESTIMATOR TYPE=EQUAL_WOR
  /POPSIZE VALUE=5000.
```

- An analysis design is specified.
- The variable *sampleweight* is specified as the variable containing sample weights for analysis.
- The EQUAL\_WOR estimator will be used for analysis.
- POPSIZE specifies that the sample was drawn from a population of 5,000.

### **Simple Analysis Design**

```
CSPLAN ANALYSIS
  /PLAN FILE='/survey/myfile.csaplan'
  /PLANVARS ANALYSISWEIGHT=sampleweight
  /DESIGN
  /ESTIMATOR TYPE=EQUAL_WOR
  /INCLPROB VALUE=0.10.
```

- An analysis design is specified.



- The variable *sampleweight* is specified as the variable containing sample weights for analysis.
- The EQUAL\_WOR estimator will be used for analysis.
- INCLPROB specifies that 10% of population units were selected for inclusion in the sample.

### **Stratified Analysis Design**

```
CSPLAN ANALYSIS
  /PLAN FILE='/survey/myfile.csaplan'
  /PLANVARS ANALYSISWEIGHT=sampleweight
  /DESIGN STRATA=region
  /ESTIMATOR TYPE=EQUAL_WOR
  /INCLPROB MATRIX=REGION; 'East' 0.1; 'West' 0.2;
                          'North' 0.1; 'South' 0.3.
```

- The analysis design specifies that the sample is stratified by *region*.
- Inclusion probabilities are specified for each stratum.
- The variable *sampleweight* is specified as the variable containing sample weights for analysis.

### **Stratified Clustering Analysis Design**

```
CSPLAN ANALYSIS
  /PLAN FILE='/survey/myfile.csaplan'
  /PLANVARS ANALYSISWEIGHT=sampleweight
  /DESIGN STRATA=district CLUSTER=school
  /ESTIMATOR TYPE=UNEQUAL_WOR.
```

- The analysis design specifies that units were sampled using stratified clustering.
- The variable *sampleweight* is specified as the variable containing sample weights for analysis.
- *District* is defined as a stratification variable and *school* is defined as a cluster variable.
- The UNEQUAL\_WOR estimator will be used for analysis.

### **Multistage Analysis Design**

```
CSPLAN ANALYSIS
  /PLAN FILE='/survey/myfile.csaplan'
  /PLANVARS ANALYSISWEIGHT=sampleweight
  /DESIGN CLUSTER=district
  /ESTIMATOR TYPE=UNEQUAL_WOR
  /DESIGN CLUSTER=school
  /ESTIMATOR TYPE=EQUAL_WOR
  /INCLPROB VARIABLE=sprob.
```

- The analysis design specifies that cases were sampled using multistage clustering. Schools were sampled within districts.
- The UNEQUAL\_WOR estimator will be used in stage 1.
- The EQUAL\_WOR estimator will be used in stage 2.
- The variable *sprob* contains inclusion probabilities, which are required for analysis of the second stage.
- The variable *sampleweight* is specified as the variable containing sample weights for analysis.

**Display Plan**

```
CSPLAN VIEW
  /PLAN FILE='/survey/myfile.csplan'.
```

- The syntax displays the specifications in the plan file *myfile.csplan*.

**CSPLAN Command**

CSPLAN creates a complex sample design or analysis specification.

<b>SAMPLE</b>	<i>Creates a sample design.</i>
<b>ANALYSIS</b>	<i>Creates an analysis design.</i>
<b>VIEW</b>	<i>Displays a sample or analysis design.</i>

**PLAN Subcommand**

The PLAN subcommand specifies the name of a design file to be written or displayed by CSPLAN. The file contains sample and/or analysis design specifications.

<b>FILE</b>	<i>Sampling design file.</i> Specify the filename in full. If you are creating a plan and the file already exists, it is overwritten without warning.
-------------	---

**PLANVARS Subcommand**

PLANVARS is used to name planwise variables to be created when a sample is extracted or used as input to the selection or estimation process.

<b>ANALYSISWEIGHT</b>	<i>Final sample weights for each unit to be used by Complex Samples analysis procedures in the estimation process.</i> ANALYSISWEIGHT is required if an analysis design is specified. It is ignored with a warning if a sample design is specified.
<b>SAMPLEWEIGHT</b>	<i>Overall sample weights that will be generated when the sample design is executed using CSSELECT.</i> A final sampling weight is created automatically when the sample plan is executed. SAMPLEWEIGHT is honored only if a sampling design is specified. It is ignored with a warning if an analysis design is specified. Sample weights are positive for selected units. They take into account all stages of the design as well as previous sampling weights if specified. If SAMPLEWEIGHT is not specified, a default name ( <i>SampleWeight_Final_</i> ) is used for the sample weight variable.
<b>PREVIOUSWEIGHT</b>	<i>Weights to be used in computing final sampling weights in a multistage design.</i> PREVIOUSWEIGHT is honored only if a sampling design is specified. It is ignored with a warning if an analysis design is specified. Typically, the previous weight variable is produced in an earlier stage of a stage-by-stage sample selection process. CSSELECT multiplies previous weights with those for the current stage to obtain final sampling weights.

For example, suppose that you want to sample individuals within cities but only city data are available at the outset of the study. For the first stage of extraction, a design plan is created that specifies that 10 cities are to be sampled from the active file. The `PLANVARS` subcommand specifies that sampling weights are to be saved under the name *CityWeights*:

```
CSPLAN SAMPLE
  /PLAN FILE='/survey/city.csplan'
  /PLANVARS SAMPLEWEIGHT=CityWeights
  /DESIGN CLUSTER=city
  /METHOD TYPE=PPS_WOR
  /MOS VARIABLE=SizeVar
  /SIZE VALUE=10.
```

This plan would be executed using `CSSELECT` on an active file in which each case is a city.

For the next stage of extraction, a design plan is created that specifies that 50 individuals are to be sampled within cities. The design uses the `PREVIOUSWEIGHT` keyword to specify that sample weights generated in the first stage are to be used when computing final sampling weights for selected individuals. Final weights are saved to the variable *FinalWeights*.

```
CSPLAN SAMPLE
  /PLAN FILE='/survey/individuals.csplan'
  /PLANVARS PREVIOUSWEIGHT=CityWeights
              SAMPLEWEIGHT=FinalWeights
  /DESIGN STRATA=city
  /METHOD TYPE=SIMPLE_WOR
  /SIZE VALUE=50.
```

The plan for stage 2 would be executed using `CSSELECT` on an active file in which cases represent individuals and both *city* and *CityWeights* are recorded for each individual. Note that *city* is identified as a stratification variable in this stage, so individuals are sampled within cities.

## ***SRSESTIMATOR Subcommand***

The `SRSESTIMATOR` subcommand specifies the variance estimator used under the simple random sampling assumption. This estimate is needed, for example, in computation of design effects in Complex Samples analysis procedures.

<b>WOR</b>	<i>SRS variance estimator includes the finite population correction.</i> This estimator is the default.
<b>WR</b>	<i>SRS variance estimator does not include the finite population correction.</i> This estimator is recommended when the analysis weights have been scaled so that they do not add up to the population size.

## ***PRINT Subcommand***

<b>PLAN</b>	<i>Displays a summary of plan specifications.</i> The output reflects your specifications at each stage of the design. The plan is shown by default. The <code>PRINT</code> subcommand is used to control output from <code>CSPLAN</code> .
<b>MATRIX</b>	<i>Displays a table of MATRIX specifications.</i> <code>MATRIX</code> is ignored if you do not use the <code>MATRIX</code> form of the <code>SIZE</code> , <code>RATE</code> , <code>POPSIZE</code> , or <code>INCLPROB</code> subcommand. By default, the table is not shown.

## ***DESIGN Subcommand***

The `DESIGN` subcommand signals a stage of the design. It also can be used to define stratification variables, cluster variables, or a descriptive label for a particular stage.

### ***STAGELABEL Keyword***

`STAGELABEL` allows a descriptive label to be entered for the stage that appears in Complex Samples procedure output.

**'Label'**                      *Descriptive stage label.* The label must be specified within quotes. If a label is not provided, a default label is generated that indicates the stage number.

### ***STRATA Keyword***

`STRATA` is used to identify stratification variables whose values represent nonoverlapping subgroups. Stratification is typically done to decrease sampling variation and/or to ensure adequate representation of small groups in a sample.

If `STRATA` is used, `CSSELECT` draws samples independently within each stratum. For example, if *region* is a stratification variable, separate samples are drawn for each region (for example, East, West, North, and South). If multiple `STRATA` variables are specified, sampling is performed within each combination of strata.

**varlist**                      *Stratification variables.*

### ***CLUSTER Keyword***

`CLUSTER` is used to sample groups of sampling units, such as states, counties, or school districts. Cluster sampling is often performed to reduce travel and/or interview costs in social surveys. For example, if census tracts are sampled within a particular city and each interviewer works within a particular tract, he or she would be able to conduct interviews within a small area, thus minimizing time and travel expenses.

- If `CLUSTER` is used, `CSSELECT` samples from values of the cluster variable as opposed to sampling elements (cases).
- If two or more cluster variables are specified, samples are drawn from among all combinations of values of the variables.
- `CLUSTER` is required for nonfinal stages of a sample or analysis plan.
- `CLUSTER` is required if any of the following sampling methods is specified: `PPS_WOR`, `PPS_BREWER`, `PPS_MURTHY`, or `PPS_SAMPFORD`.
- `CLUSTER` is required if the `UNEQUAL_WOR` estimator is specified.

**varlist**                      *Cluster variables.*

## METHOD Subcommand

The METHOD subcommand specifies the sample extraction method. A variety of equal- and unequal-probability methods are available.

The following table lists extraction methods and their availability at each stage of the design. For details on each method, see the CSSELECT algorithms document.

- PPS methods are available only in stage 1. WR methods are available only in the final stage. Other methods are available in any stage.
- If a PPS method is chosen, a measure of size (MOS) must be specified.
- If the PPS\_WOR, PPS\_BREWER, PPS\_SAMPFORD, or PPS\_MURTHY method is selected, first-stage joint inclusion probabilities are written to an external file when the sample plan is executed. Joint probabilities are needed for UNEQUAL\_WOR estimation by Complex Samples analysis procedures.
- By default, CSPLAN chooses an appropriate estimation method for the selected sampling method. If ESTIMATION=WR, Complex Samples analysis procedures use the WR (with replacement) estimator regardless of the sampling method.

Type	Description	Default estimator
SIMPLE_WOR	Selects units with equal probability. Units are extracted without replacement.	EQUAL_WOR
SIMPLE_WR	Selects units with equal probability. Units are extracted with replacement.	WR
SIMPLE_SYSTEMATIC	Selects units at a fixed interval throughout the sampling frame or stratum. A random starting point is chosen within the first interval.	WR
SIMPLE_CHROMY	Selects units sequentially with equal probability. Units are extracted without replacement.	WR
PPS_WOR	Selects units with probability proportional to size. Units are extracted without replacement.	UNEQUAL_WOR
PPS_WR	Selects units with probability proportional to size. Units are extracted with replacement.	WR
PPS_SYSTEMATIC	Selects units by systematic random sampling with probability proportional to size. Units are extracted without replacement.	WR
PPS_CHROMY	Selects units sequentially with probability proportional to size without replacement.	WR
PPS_BREWER	Selects two units from each stratum with probability proportional to size. Units are extracted without replacement.	UNEQUAL_WOR
PPS_MURTHY	Selects two units from each stratum with probability proportional to size. Units are extracted without replacement.	UNEQUAL_WOR
PPS_SAMPFORD	An extension of the Brewer's method that selects more than two units from each stratum with probability proportional to size. Units are extracted without replacement.	UNEQUAL_WOR

## ***ESTIMATION Keyword***

By default, the estimation method used when sample data are analyzed is implied by the specified extraction method. If `ESTIMATION=WR` is specified, the with-replacement estimator is used when summary statistics are produced using Complex Samples analysis procedures.

- The `WR` keyword has no effect if the specified `METHOD` implies `WR` estimation.
- If `ESTIMATION=WR` is specified, the joint probabilities file is not created when the sample plan is executed.
- `ESTIMATION=WR` is available only in the first stage.

## ***SIZE Subcommand***

The `SIZE` subcommand specifies the number of sampling units to draw at the current stage.

- You can specify a single value, a variable name, or a matrix of counts for design strata.
- Size values must be positive integers.
- The `SIZE` subcommand is ignored with a warning if the `PPS_MURTHY` or `PPS_BREWER` method is specified.
- The `SIZE` or `RATE` subcommand must be specified for each stage. An error occurs if both are specified.

**VALUE**            *Apply a single value to all strata.* For example, `VALUE=10` selects 10 units per stratum.

**MATRIX**        *Specify disproportionate sample sizes for different strata.* Specify one or more variables after the `MATRIX` keyword. Then provide one size specification per stratum. A size specification includes a set of category values and a size value. Category values should be listed in the same order as variables to which they apply. Semicolons are used to separate the size specifications.

For example, the following syntax selects 10 units from the North stratum and 20 from the South stratum:

```
/SIZE MATRIX=region; 'North' 10; 'South' 20
```

If there is more than one variable, specify one size per combination of strata. For example, the following syntax specifies rate values for combinations of Region and Sex strata:

```
/SIZE MATRIX=region sex; 'North' 'Male' 10; 'North' 'Female' 15;  
'South' 'Male' 24; 'South' 'Female' 30
```

The variable list must contain all or a subset of stratification variables from the same and previous stages and cluster variables from the previous stages. An error occurs if the list contains variables that are not defined as strata or cluster variables.

Each size specification must contain one category value per variable. If multiple size specifications are provided for the same strata or combination of strata, only the last one is honored.

String and date category values must be quoted.

A semicolon must appear after the variable list and after each size specification. The semicolon is not allowed after the last size specification.

**VARIABLE**       *Specify the name of a single variable that contains the sample sizes.*

## **RATE Subcommand**

The `RATE` subcommand specifies the percentage of units to draw at the current stage—that is, the sampling fraction.

- Specify a single value, a variable name, or a matrix of rates for design strata. In all cases, the value 1 is treated as 100%.
- Rate values must be positive.
- `RATE` is ignored with a warning if the `PPS_MURTHY` or `PPS_BREWER` method is specified.
- Either `SIZE` or `RATE` must be specified for each stage. An error occurs if both are specified.

**VALUE**            *Apply a single value to all strata.* For example, `VALUE=.10` selects 10% of units per stratum.

**MATRIX**        *Specify disproportionate rates for different strata.* Specify one or more variables after the `MATRIX` keyword. Then provide one rate specification per stratum. A rate specification includes a set of category values and a rate value. Category values should be listed in the same order as variables to which they apply. Semicolons are used to separate the rate specifications.

For example, the following syntax selects 10% of units from the North stratum and 20% from the South stratum:

```
/RATE MATRIX=region; 'North' .1; 'South' .2
```

If there is more than one variable, specify one rate per combination of strata. For example, the following syntax specifies rate values for combinations of Region and Sex strata:

```
/RATE MATRIX=region sex; 'North' 'Male' .1; 'North' 'Female' .15;
                          'South' 'Male' .24; 'South' 'Female' .3
```

The variable list must contain all or a subset of stratification variables from the same and previous stages and cluster variables from the previous stages. An error occurs if the list contains variables that are not defined as strata or cluster variables.

Each rate specification must contain one category value per variable. If multiple rate specifications are provided for the same strata or combination of strata, only the last one is honored.

String and date category values must be quoted. A semicolon must appear after the variable list and after each rate specification.

The semicolon is not allowed after the last rate specification.

**VARIABLE**       *Specify the name of a single variable that contains the sample rates.*

## **MINSIZE Keyword**

`MINSIZE` specifies the minimum number of units to draw when `RATE` is specified. `MINSIZE` is useful when the sampling rate for a particular stratum turns out to be very small due to rounding.

**value**            *The value must be a positive integer.* An error occurs if the value exceeds `MAXSIZE`.

### **MAXSIZE Keyword**

MAXSIZE specifies the maximum number of units to draw when RATE is specified. MAXSIZE is useful when the sampling rate for a particular stratum turns out to be larger than desired due to rounding.

**value**                      *The value must be a positive integer. An error occurs if the value is less than MINSIZE.*

### **MOS Subcommand**

The MOS subcommand specifies the measure of size for population units in a PPS design. Specify a variable that contains the sizes or request that sizes be determined when CSSELECT scans the sample frame.

**VARIABLE**                      *Specify a variable containing the sizes.*  
**SOURCE=FROMDATA**              *The CSSELECT procedure counts the number of cases that belong to each cluster to determine the MOS. SOURCE=FROMDATA can be used only if a CLUSTER variable is defined. Otherwise, an error is generated.*

- The MOS subcommand is required for PPS designs. Otherwise, it is ignored with a warning.

### **MIN Keyword**

MIN specifies a minimum MOS for population units that overrides the value specified in the MOS variable or obtained by scanning the data.

**value**                      *The value must be positive. MIN must be less than or equal to MAX.*

- MIN is optional for PPS methods. It is ignored for other methods.

### **MAX Keyword**

MAX specifies a maximum MOS for population units that overrides the value specified in the MOS variable or obtained by scanning the data.

**value**                      *The value must be positive. MAX must be greater than or equal to MIN.*

- MAX is optional for PPS methods. It is ignored for other methods.

### **STAGEVARS Subcommand**

The STAGEVARS subcommand is used to obtain stagewise sample information variables when a sample design is executed. Certain variables are created automatically and cannot be suppressed. The names of both automatic and optional stagewise variables can be user-specified.

- Stagewise inclusion probabilities and cumulative sampling weights are always created.



- A stagewise duplication index is created only when sampling is done with replacement. A warning occurs if index variables are requested when sampling is done without replacement.
- If a keyword is specified without a variable name, a default name is used. The default name indicates the stage to which the variable applies.

### Example

```
/STAGEVARS POPSIZE
           INCLPROB(SelectionProb)
```

- The syntax requests that the population size for the stage be saved using a default name.
- Inclusion probabilities for the stage will be saved using the name *SelectionProb*. (Note that inclusion probabilities are always saved when the sample design is executed. The syntax shown here requests that they be saved using a nondefault name.)

## STAGEVARS Variables

The following table shows available STAGEVARS variables. See the CSSELECT algorithms document for a detailed explanation of each quantity.

If the default variable name is used, a numeric suffix that corresponds to the stage number is added to the root shown below. All names end in an underscore—for example, *InclusionProbability\_1\_*.

Keyword	Default root name	Description	Generated automatically when sample executed?
INCLPROB	<i>InclusionProbability_</i>	Stagewise inclusion (selection) probabilities. The proportion of units drawn from the population at a particular stage.	Yes
CUMWEIGHT	<i>SampleWeightCumulative_</i>	Cumulative sampling weight for a given stage. Takes into account prior stages.	Yes
INDEX	<i>Index_</i>	Duplication index for units selected in a given stage. The index uniquely identifies units selected more than once when sampling is done with replacement.	Yes, when sampling is done with replacement.
POPSIZE	<i>PopulationSize_</i>	Population size for a given stage.	No
SAMPsize	<i>SampleSize_</i>	Number of units drawn at a given stage.	No

Keyword	Default root name	Description	Generated automatically when sample executed?
RATE	<i>SamplingRate_</i>	Stagewise sampling rate.	No
WEIGHT	<i>SampleWeight_</i>	Sampling weight for a given stage. The inverse of the stagewise inclusion probability. Stage weights are positive for each unit selected in a particular stage.	No

## ***ESTIMATOR Subcommand***

The `ESTIMATOR` subcommand is used to choose an estimation method for the current stage. There is no default estimator.

Available estimators depend on the stage:

- `EQUAL_WOR` can be specified in any stage of the design.
- `UNEQUAL_WOR` can be specified in the first stage only. An error occurs if it is used in stage 2 or 3.
- `WR` can be specified in any stage. However, the stage in which it is specified is treated as the last stage. Any subsequent stages are ignored when the data are analyzed.

**EQUAL\_WOR**                      *Equal selection probabilities without replacement.* `POPSIZE` or `INCLPROB` must be specified.

**UNEQUAL\_WOR**                      *Unequal selection probabilities without replacement.* If `POPSIZE` or `INCLPROB` is specified, it is ignored and a warning is issued.

**WR**                                      *Selection with replacement.* If `POPSIZE` or `INCLPROB` is specified, it is ignored and a warning is issued.

## ***POPSIZE Subcommand***

The `POPSIZE` subcommand specifies the population size for each sample element. Specify a single value, a variable name, or a matrix of counts for design strata.

- The `POPSIZE` and `INCLPROB` subcommands are mutually exclusive. An error occurs if both are specified for a particular stage.
- Population size values must be positive integers.

**VALUE**                              *Apply a single value to all strata.* For example, `VALUE=1000` indicates that each stratum has a population size of 1,000.

**MATRIX**                              *Specify disproportionate population sizes for different strata.* Specify one or more variables after the `MATRIX` keyword. Then provide one size specification per stratum. A size specification includes a set of category values and a population size value. Category values should be listed in the same order as variables to which they apply. Semicolons are used to separate the size specifications.

For example, the following syntax specifies that units in the North stratum were sampled from a population of 1,000. The population size for the South stratum is specified as 2,000:

```
/SIZE MATRIX=region; 'North' 1000; 'South' 2000
```

If there is more than one variable, specify one size per combination of strata. For example, the following syntax specifies rate values for combinations of Region and Sex strata:

```
/SIZE MATRIX=region sex; 'North' 'Male' 1000; 'North' 'Female' 1500;
'South' 'Male' 2400; 'South' 'Female' 3000
```

The variable list must contain all or a subset of stratification variables from the same and previous stages and cluster variables from the previous stages. An error occurs if the list contains variables that are not defined as strata or cluster variables.

Each size specification must contain one category value per variable. If multiple size specifications are provided for the same strata or combination of strata, only the last one is honored.

String and date category values must be quoted.

A semicolon must appear after the variable list and after each size specification. The semicolon is not allowed after the last size specification.

#### VARIABLE

*Specify the name of a single variable that contains the population sizes.*

## INCLPROB Subcommand

The INCLPROB subcommand specifies the proportion of units drawn from the population at a given stage. Specify a single value, a variable name, or a matrix of inclusion probabilities for design strata.

- The POPSIZE and INCLPROB subcommands are mutually exclusive. An error occurs if both are specified for a particular stage.
- Proportions must be a positive value less than or equal to 1.

#### VALUE

*Apply a single value to all strata.* For example, VALUE=0.10 indicates that 10% of elements in each stratum were selected.

#### MATRIX

*Specify unequal proportions for different strata.* Specify one or more variables after the MATRIX keyword. Then provide one proportion per stratum. A proportion specification includes a set of category values and a proportion value. Category values should be listed in the same order as variables to which they apply. Semicolons are used to separate the proportion specifications.

For example, the following syntax indicates that 10% of units were selected from the North stratum and 20% were selected from the South stratum:

```
/INCLPROB MATRIX=region; 'North' 0.1; 'South' 0.2
```

If there is more than one variable, specify one proportion per combination of strata. For example, the following syntax specifies proportions for combinations of Region and Sex strata:

```
/INCLPROB MATRIX=region sex; 'North' 'Male' 0.1; 'North' 'Female' 0.15;
'South' 'Male' 0.24; 'South' 'Female' 0.3
```

The variable list must contain all or a subset of stratification variables from the same and previous stages and cluster variables from the previous stages. An error occurs if the list contains variables that are not defined as strata or cluster variables.

Each proportion specification must contain one category value per variable. If multiple proportions are provided for the same strata or combination of strata, only the last one is honored.

String and date category values must be quoted.

A semicolon must appear after the variable list and after each proportion specification. The semicolon is not allowed after the last proportion specification.

**VARIABLE**

*Specify the name of a single variable that contains inclusion probabilities.*

# CSSELECT

CSSELECT is available in the Complex Samples option.

```
CSSELECT
/PLAN FILE='file'
[/CRITERIA [STAGES=n [n [n]]] [SEED={RANDOM**}]
           {value  }]
[/CLASSMISSING {EXCLUDE**}]
           {INCLUDE  }
[/DATA [RENAMEVARS] [PRESORTED]]
[/SAMPLEFILE OUTFILE='savfile'|'dataset' [KEEP=varlist] [DROP=varlist]]
[/JOINTPROB OUTFILE='savfile'|'dataset']
[/SELECTRULE OUTFILE='file']
[/PRINT [SELECTION**] [CPS]]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 16.0

- Added support for SET THREADS.

## Example

```
CSSELECT
  /PLAN FILE='/survey/myfile.csplan'.
```

## Overview

CSSELECT selects complex, probability-based samples from a population. CSSELECT selects units according to a sample design created using the CSPLAN procedure.

## Options

**Scope of Execution.** By default, CSSELECT executes all stages defined in the sampling plan. Optionally, you can execute specific stages of the design. This capability is useful if a full sampling frame is not available at the outset of the sampling process, in which case new stages can be sampled as they become available. For example, CSSELECT might first be used to sample cities, then to sample blocks, and finally to sample individuals. Each time a different stage of the sampling plan would be executed.

**Seed.** By default, a random seed value is used by the CSSELECT random number generator. You can specify a seed to ensure that the same sample will be drawn when CSSELECT is invoked repeatedly using the same sample plan and population frame. The CSSELECT seed value is independent of the global seed specified via the SET command.

**Missing Values.** A case is excluded from the sample frame if it has a system-missing value for any input variable in the plan file. You can control whether user-missing values of stratification and cluster variables are treated as invalid. User-missing values of measure variables are always treated as invalid.

**Input Data.** If the sampling frame is sorted in advance, you can specify that the data are presorted, which may improve performance when stratification and/or clustering is requested for a large sampling frame.

**Sample Data.** CSSELECT writes data to the active dataset (the default) or an external file. Regardless of the data destination, CSSELECT generates final sampling weights, stagewise inclusion probabilities, stagewise cumulative sampling weights, as well as variables requested in the sampling plan.

External files or datasets produced by CSSELECT include selected cases only. By default, all variables in the active dataset are copied to the external file or dataset. Optionally, you can specify that only certain variables are to be copied.

**Joint Probabilities.** First-stage joint inclusion probabilities are automatically saved to an external file when the plan file specifies a PPS without-replacement sampling method. Joint probabilities are used by Complex Samples analysis procedures, such as CSDESCRIPTIVES and CSTABULATE. You can control the name and location of the joint probabilities file.

**Output.** By default, CSSELECT displays the distribution of selected cases by stratum. Optionally, you can display a case-processing summary.

### **Basic Specification**

- The basic specification is a PLAN subcommand that specifies a sample design file.
- By default, CSPLAN writes output data to the active dataset including final sample weights, stagewise cumulative weights, and stagewise inclusion probabilities. See the CSPLAN design for a description of available output variables.

### **Operations**

- CSSELECT selects sampling units according to specifications given in a sample plan. Typically, the plan is created using the CSPLAN procedure.
- In general, elements are selected. If cluster sampling is performed, groups of elements are selected.
- CSSELECT assumes that the active dataset represents the sampling frame. If a multistage sample design is executed, the active dataset should contain data for all stages. For example, if you want to sample individuals within cities and city blocks, then each case should be an individual, and city and block variables should be coded for each individual. When CSSELECT is used to execute particular stages of the sample design, the active dataset should represent the subframe for those stages only.

- A case is excluded from the sample frame if it has a system-missing value for any input variable in the plan.
- You can control whether user-missing values of stratification and cluster variables are treated as valid. By default, they are treated as invalid.
- User-missing values of measure variables are always treated as invalid.
- The CSSELECT procedure has its own seed specification that is independent of the global SET command.
- First-stage joint inclusion probabilities are automatically saved to an external file when the plan file specifies a PPS without-replacement sampling method. By default, the joint probabilities file is given the same name as the plan file (with a different extension) and is written to the same location.
- Output data must be written to an external data file if with-replacement sampling is specified in the plan file.
- This procedure uses the multithreaded options specified by SET THREADS.

### **Syntax Rules**

- The PLAN subcommand is required. All other subcommands are optional.
- Only a single instance of each subcommand is allowed.
- An error occurs if an attribute or keyword is specified more than once within a subcommand.
- An error occurs if the same output file is specified for more than one subcommand.
- Equals signs shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.
- Empty subcommands are not allowed.

### **Limitations**

- WEIGHT and SPLIT FILE settings are ignored with a warning by the CSSELECT procedure.

## **Example**

```
CSSELECT
  /PLAN FILE='/survey/myfile.csplan'
  /CRITERIA SEED=99999
  /SAMPLEFILE OUTFILE='/survey/sample.sav' .
```

- CSSELECT reads the plan file *myfile.csplan*.
- CSSELECT draws cases according to the sampling design specified in the plan file.
- Sampled cases and weights are written to an external file. By default, output data include final sample weights, stagewise inclusion probabilities, stagewise cumulative weights, and any other variables requested in the sample plan.
- The seed value for the random number generator is 99999.

## **PLAN Subcommand**

PLAN identifies the plan file whose specifications are to be used for selecting sampling units. FILE specifies the name of the file. An error occurs if the file does not exist.

## **CRITERIA Subcommand**

CRITERIA is used to control the scope of execution and specify a seed value.

## **STAGES Keyword**

STAGES specifies the scope of execution.

- By default, all stages defined in the sampling plan are executed. STAGES is used to limit execution to specific stages of the design.
- Specify one or more stages. The list can include up to three integer values—for example, STAGES=1 2 3. If two or more values are provided, they must be consecutive. An error occurs if a stage is specified that does not correspond to a stage in the plan file.
- If the sample plan specifies a previous weight variable, it is used in the first stage of the plan.
- When executing latter stages of a multistage sampling design in which the earlier stages have already been sampled, CSSELECT requires the cumulative sampling weights of the last stage sampled, in order to compute the correct final sampling weights for the whole design. For example, if you have executed the first two stages of a three-stage design and saved the second-stage cumulative weights to *SampleWeightCumulative\_2\_*, when you sample the third stage of the design, the active dataset must contain *SampleWeightCumulative\_2\_* to compute the final sampling weights.

## **SEED Keyword**

SEED specifies the random number seed used by the CSSELECT procedure.

- By default, a random seed value is selected. To replicate a particular sample, the same seed, sample plan, and sample frame should be specified when the procedure is executed.
- The CSSELECT seed value is independent of the global seed specified via the SET command.

<b>RANDOM</b>	<i>A seed value is selected at random.</i> This is the default.
<b>value</b>	<i>Specifies a custom seed value.</i> The seed value must be a positive integer.

## **CLASSMISSING Subcommand**

CLASSMISSING is used to control whether user-missing values of classification (stratification and clustering) variables are treated as valid values. By default, they are treated as invalid.

<b>EXCLUDE</b>	<i>User-missing values of stratification and cluster variables are treated as invalid.</i> This is the default.
<b>INCLUDE</b>	<i>User-missing values of stratification and cluster variables are treated as valid values.</i>



CSSELECT always treats user-missing values of measure variables (previous weight, MOS, size, and rate) as invalid.

## ***DATA Subcommand***

DATA specifies general options concerning input and output files.

## ***RENAMEVARS Keyword***

The RENAMEVARS keyword handles name conflicts between existing variables and variables to be created by the CSSELECT procedure.

- If the RENAMEVARS keyword is not specified, conflicting variable names generate an error. This is the default.
- If output data are directed to the active dataset, RENAMEVARS specifies that an existing variable should be renamed with a warning if its name conflicts with that of a variable created by the CSSELECT procedure.
- If output data are directed to an external file or dataset, RENAMEVARS specifies that a variable to be copied from the active dataset should be renamed, with a warning if its name conflicts with that of a variable created by the CSSELECT procedure. See the SAMPLEFILE subcommand for details about copying variables from the active dataset.

## ***PRESORTED Keyword***

By default, CSSELECT assumes that the active dataset is unsorted. The PRESORTED keyword specifies that the data are sorted in advance, which may improve performance when stratification and/or clustering is requested for a large sample frame.

If PRESORTED is used, the data should be sorted first by all stratification variables then by cluster variables consecutively in each stage. The data can be sorted in ascending or descending order. For example, given a sample plan created using the following CSPLAN syntax, the sample frame should be sorted by region, ses, district, type, and school, in that order.

### ***Example***

```
CSPLAN
  /PLAN OUTFILE='/survey/myfile.csplan'
  /DESIGN STRATA=region ses CLUSTER=district type
  /SAMPLE RATE=.2 MOS=districtsize METHOD=PPS_WOR
  /DESIGN CLUSTER=school
  /SAMPLE RATE=.3 METHOD=SIMPLE_WOR.
```

An error occurs if PRESORTED is specified and the data are not sorted in proper order.

## ***SAMPLEFILE Subcommand***

`SAMPLEFILE` is used to write sampled units to an external file or dataset. Datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files.

- The external file or dataset contains sampled cases only. By default, all variables in the active dataset are copied to the external file or dataset.
- If `SAMPLEFILE` is specified, data are not written to the active dataset.
- `SAMPLEFILE` must be used if with-replacement sampling is specified in the plan file. Otherwise, an error is generated.
- `KEEP` and `DROP` can be used simultaneously; the effect is cumulative. An error occurs if you specify a variable already named on a previous `DROP` or one not named on a previous `KEEP`.

## ***OUTFILE Keyword***

The `OUTFILE` keyword specifies the name of the external file or the name of a dataset. An external file, a file handle, or a dataset name must be specified. If the file or dataset exists, it is overwritten without warning.

## ***KEEP Keyword***

The `KEEP` keyword lists variables to be copied from the active dataset to the file or dataset specified on the `OUTFILE` keyword. `KEEP` has no bearing on the active dataset.

- At least one variable must be specified.
- Variables not listed are not copied.
- An error occurs if a specified variable does not exist in the active dataset.
- Variables are copied in the order in which they are listed.

## ***DROP Keyword***

The `DROP` keyword excludes variables from the file or dataset specified on the `OUTFILE` keyword. `DROP` has no bearing on the active dataset.

- At least one variable must be specified.
- Variables not listed are copied.
- The `ALL` keyword can be used to drop all variables.
- An error occurs if a specified variable does not exist in the active dataset.

## ***JOINTPROB Subcommand***

First-stage joint inclusion probabilities are automatically saved to an external IBM® SPSS® Statistics data file when the plan file specifies a PPS without-replacement sampling method. By default, the joint probabilities file is given the same name as the plan file (with a different

extension), and it is written to the same location. JOINTPROB is used to override the default name and location of the file.

- OUTFILE specifies the name of the file. In general, if the file exists, it is overwritten without warning.
- The joint probabilities file is generated only when the plan file specifies PPS\_WOR, PPS\_BREWER, PPS\_SAMPFORD, or PPS\_MURTHY as the sampling method. A warning is generated if JOINTPROB is used when any other sampling method is requested in the plan file.

### ***Structure of the Joint Probabilities File***

Complex Samples analysis procedures will expect the following variables in the joint probability file in the order listed below. If there are other variables beyond the joint probability variables, they will be silently ignored.

1. **Stratification variables.** These are the stratification variables used in the first stage of sampling. If there is no stratification in first stage, no stratification variables are included in the file.
2. **Cluster variables.** These are variables used to identify each primary sampling unit (PSU) within a stratum. At least one cluster variable is always included, since it is required for all selection methods that generate the joint probabilities as well as for the estimation method using them.
3. **System PSU id.** This variable labels PSU's within a stratum. The variable name used is *Unit\_No\_*.
4. **Joint probability variables.** These variables store the joint inclusion probabilities for each pair of units. The default names of these variables will have the form *Joint\_Prob\_n\_*; for example, the joint inclusion probabilities of the 2nd and 3rd units will be the values located at case 2 of *Joint\_Prob\_3\_* or case 3 of *Joint\_Prob\_2\_*. Since the analysis procedures extract joint probabilities by location, it is safe to rename these variables at your convenience. Within each stratum, these joint inclusion probabilities will form a square symmetric matrix. Since the joint inclusion probabilities only vary for the off diagonal entries, the diagonal elements correspond to the first stage inclusion probabilities. The maximum number of joint inclusion probability variables will be equal to the maximum sample size across all strata.

**Example**

Figure 48-1  
Joint probabilities file

	county	town	Unit_No_	Joint_Prob_1	Joint_Prob_2	Joint_Prob_3	Joint_Prob_4	Joint_Prob_5
1	1	10	1	.31	.10	.11	.12	.
2	1	11	2	.10	.39	.15	.16	.
3	1	9	3	.11	.15	.44	.21	.
4	1	12	4	.12	.16	.21	.48	.
5	2	12	1	.22	.04	.07	.08	.
6	2	6	2	.04	.23	.07	.08	.
7	2	7	3	.07	.07	.41	.19	.
8	2	2	4	.08	.08	.19	.45	.
9	3	5	1	.58	.31	.32	.	.
10	3	3	2	.31	.61	.36	.	.
11	3	4	3	.32	.36	.63	.	.
12	4	14	1	.26	.06	.06	.07	.09
13	4	8	2	.06	.29	.07	.08	.10
14	4	4	3	.06	.07	.29	.08	.10
15	4	2	4	.07	.08	.08	.33	.12
16	4	13	5	.09	.10	.10	.12	.43
17	5	3	1	.74	.25	.27	.	.
18	5	6	2	.25	.41	.13	.	.
19	5	4	3	.27	.13	.43	.	.

The file *poll\_jointprob.sav* contains first-stage joint probabilities for selected townships within counties. *County* is a first-stage stratification variable, and *Township* is a cluster variable. Combinations of these variables identify all first-stage PSUs uniquely. *Unit\_No\_* labels PSUs within each stratum and is used to match up with *Joint\_Prob\_1\_*, *Joint\_Prob\_2\_*, *Joint\_Prob\_3\_*, *Joint\_Prob\_4\_*, and *Joint\_Prob\_5\_*. The first two strata each have 4 PSUs; therefore, the joint inclusion probability matrices are 4×4 for these strata, and the *Joint\_Prob\_5\_* column is left empty for these rows. Similarly, strata 3 and 5 have 3×3 joint inclusion probability matrices, and stratum 4 has a 5×5 joint inclusion probability matrix.

The need for a joint probabilities file is seen by perusing the values of the joint inclusion probability matrices. When the sampling method is not a PPS WOR method, the selection of a PSU is independent of the selection of another PSU, and their joint inclusion probability is simply the product of their inclusion probabilities. In contrast, the joint inclusion probability for Townships 9 and 10 of County 1 is approximately 0.11 (see the first case of *Joint\_Prob\_3\_* or the third case of *Joint\_Prob\_1\_*), or less than the product of their individual inclusion probabilities (the product of the first case of *Joint\_Prob\_1\_* and the third case of *Joint\_Prob\_3\_* is  $0.31 \times 0.44 = 0.1364$ ).

**SELETRULE Subcommand**

SELETRULE generates a text file containing a rule that describes characteristics of selected units.

- The selection rule is not generated by default.

- `OUTFILE` specifies the name of the file. If the file exists, it is overwritten without warning.
- The selection rule is written in generic notation, for example—`(a EQ 1) AND (b EQ 2) '`. You can transform the selection rule into SQL code or command syntax that can be used to extract a subframe for the next stage of a multistage extraction.

## ***PRINT Subcommand***

`PRINT` controls output display.

- |                  |  |
|------------------|--|
| <b>SELECTION</b> | <i>Summarizes the distribution of selected cases across strata.</i> The information is reported per design stage. The table is shown by default. |
| <b>CPS</b>       | <i>Displays a case processing summary.</i>   |

# CSTABULATE

CSTABULATE is available in the Complex Samples option.

```
CSTABULATE
/PLAN FILE = file
[/JOINTPROB FILE = file]
/TABLES VARIABLES = varlist [BY varname]
[/CELLS [POPSIZE] [ROWPCT] [COLPCT] [TABLEPCT]]
[/STATISTICS [SE] [CV] [DEFF] [DEFFSQRT] [CIN [{95** }]] [COUNT]
           {value}
           --- options for one-way frequency tables ---
           [CUMULATIVE]
           --- options for two-way crosstabulations ---
           [EXPECTED] [RESID] [ASRESID]]
[/TEST --- options for one-way frequency tables ---
       [HOMOGENEITY]
       --- options for two-way crosstabulations ---
       [INDEPENDENCE]
       --- options for two-by-two crosstabulations ---
       [ODDSRATIO] [RELISK] [RISKDIFF]]
[/SUBPOP TABLE = varname [BY varname [BY ...]] [DISPLAY = {LAYERED }]]
           {SEPARATE }
[/MISSING [SCOPE = {TABLE }]] [CLASSMISSING = {EXCLUDE }]]
           {LISTWISE}           {INCLUDE }
```

\*\* Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
CSTABULATE
/PLAN FILE = '/survey/myfile.xml'
/TABLES VARIABLES = a.
```

## Overview

CSTABULATE displays one-way frequency tables or two-way crosstabulations, and associated standard errors, design effects, confidence intervals, and hypothesis tests, for samples drawn by complex sampling methods. The procedure estimates variances by taking into account the sample design used to select the sample, including equal probability and probability proportional to

size (PPS) methods, and with-replacement (WR) and without-replacement (WOR) sampling procedures. Optionally, CSTABULATE creates tables for subpopulations.

### **Basic Specification**

- The basic specification is a PLAN subcommand and the name of a complex sample analysis specification file, which may be generated by CSPLAN, and a TABLES subcommand with at least one variable specified.
- This specification displays a population size estimate and its standard error for each cell in the defined table, as well as for all marginals.

### **Operations**

- CSTABULATE computes table statistics for sampling designs supported by CSPLAN and CSSELECT.
- The input dataset must contain the variables to be analyzed and variables related to the sampling design.
- The complex sample analysis specification file provides an analysis plan based on the sampling design.
- For each cell and marginal in the defined table, the default output is the population size estimate and its standard error.
- WEIGHT and SPLIT FILE settings are ignored by CSTABULATE.

### **Syntax Rules**

- The PLAN and TABLES subcommands are required. All other subcommands are optional.
- Each subcommand may be specified only once.
- Subcommands can be specified in any order.
- All subcommand names and keywords must be spelled in full.
- Equals signs (=) shown in the syntax chart are required.
- Empty subcommands are not allowed.

## **Examples**

### **Example: Frequency tables**

```
* Complex Samples Frequencies.
CSTABULATE
/PLAN FILE = 'samplesDirectory\nhis2000_subset.csaplan'
/TABLES VARIABLES = VITANY
/CELLS POPSIZE TABLEPCT
/STATISTICS SE CIN(95)
/SUBPOP TABLE = AGE_CAT
/MISSING SCOPE = TABLE CLASSMISSING = EXCLUDE.
```

- The procedure will compute estimates based on the complex sample analysis specification given in *nhis2000\_subset.csaplan*.

- One-way frequency tables are produced for variable *VITANY*. Estimates, standard errors, and 95% confidence intervals are displayed for the population size and table percent for each category.
- In addition, a separate table is produced for these statistics by levels of *AGE\_CAT*.
- All other options are set to their default values.

**Example: Crosstabulation table**

```
* Complex Samples Crosstabs.
CSTABULATE
  /PLAN FILE = 'samplesDirectory\demo.csplan'
  /TABLES VARIABLES = news BY response
  /SUBPOP TABLE = inccat DISPLAY=LAYERED
  /CELLS ROWPCT
  /STATISTICS SE
  /TEST ODDSRATIO RELRISK
  /MISSING SCOPE = LISTWISE CLASSMISSING = INCLUDE.
```

- The procedure will compute estimates based on the complex sampling plan in *demo.csplan*.
- The crosstabulation of *news* by *response* is produced overall and again by levels of *inccat*.
- The estimates and standard errors of the row percentages are reported in the cells of the crosstabulation tables.
- In addition, the odds ratio and relative risk for *news* by *response* is computed for the overall population and separately for levels of *inccat*.
- All other options are set to their default values.

## **PLAN Subcommand**

The `PLAN` subcommand specifies the name of an XML file containing analysis design specifications. This file is written by `CSPLAN`.

- The `PLAN` subcommand is required.

**FILE**                      *Specifies the name of an external file.*

## **JOINTPROB Subcommand**

The `JOINTPROB` subcommand is used to specify the file or dataset containing the first stage joint inclusion probabilities for the `UNEQUAL_WOR` estimation. `CSSELECT` writes this file in the same location and with the same name (but different extension) as the plan file. When the `UNEQUAL_WOR` estimation is specified, `CSTABULATE` will use the default location and name of the file unless the `JOINTPROB` subcommand is used to override them.

**FILE**                      *Specifies the name of the file or dataset containing the joint inclusion probabilities.*

## **TABLES Subcommand**

The `TABLES` subcommand specifies the tabulation variables.



- If a single variable list is specified, then a one-way frequency table is displayed for each variable in the list.
- If the variable list is followed by the BY keyword and a variable, then two-way crosstabulations are displayed for each pair of variables. Pairs of variables are defined by crossing the variable list to the left of the BY keyword with the variable to the right. Each variable on the left defines the row dimension in a two-way crosstabulation, and the variable to the right defines the column dimension. For example, TABLES VARIABLES = A B BY C displays two tables: *A by C* and *B by C*.
- Numeric or string variables may be specified.
- Plan file and subpopulation variables may not be specified on the TABLES subcommand.
- Within the variable list, all specified variables must be unique. Also, if a variable is specified after the BY keyword, then it must be different from all variables preceding the BY keyword.

**VARIABLES**            *Specifies the tabulation variables.*

## **CELLS Subcommand**

The CELLS subcommand requests various summary value estimates associated with the table cells.

If the CELLS subcommand is not specified, then CSTABULATE displays the population size estimate for each cell in the defined table(s), as well as for all marginals. However, if the CELLS subcommand is specified, then only those summary values that are requested are displayed.

**POPSIZE**            *The population size estimate for each cell and marginal in a table. This is the default output if the CELLS subcommand is not specified.*

**ROWPCT**            *Row percentages. The population size estimate in each cell in a row is expressed as a percentage of the population size estimate for that row. Available for two-way crosstabulations. For one-way frequency tables, specifying this keyword gives the same output as the TABLEPCT keyword.*

**COLPCT**            *Column percentages. The population size estimate in each cell in a column is expressed as a percentage of the population size estimate for that column. Available for two-way crosstabulations. For one-way frequency tables, specifying this keyword gives the same output as the TABLEPCT keyword.*

**TABLEPCT**          *Table percentages. The population size estimate in each cell of a table is expressed as a percentage of the population size estimate for that table.*

## **STATISTICS Subcommand**

The STATISTICS subcommand requests various statistics associated with the summary value estimates in the table cells.

If the STATISTICS subcommand is not specified, then CSTABULATE displays the standard error for each summary value estimate in the defined table(s) cells. However, if the STATISTICS subcommand is specified, then only those statistics that are requested are displayed.

**SE**                    *The standard error for each summary value estimate. This is the default output if the STATISTICS subcommand is not specified.*

**CV**                    *Coefficient of variation.*

**DEFF**                *Design effects.*

<b>DEFFSQRT</b>	<i>Square root of the design effects.</i>
<b>CIN [(value)]</b>	<i>Confidence interval.</i> If the <b>CIN</b> keyword is specified alone, then the default 95% confidence interval is computed. Optionally, <b>CIN</b> may be followed by a value in parentheses, where $0 \leq \text{value} < 100$ .
<b>COUNT</b>	<i>Unweighted counts.</i> The number of valid observations in the dataset for each summary value estimate.
<b>CUMULATIVE</b>	<i>Cumulative summary value estimates.</i> Available for one-way frequency tables only.
<b>EXPECTED</b>	<i>Expected summary value estimates.</i> The summary value estimate in each cell if the two variables in a crosstabulation are statistically independent. Available for two-way crosstabulations only and displayed only if the <b>TABLEPCT</b> keyword is specified on the <b>CELLS</b> subcommand.
<b>RESID</b>	<i>Residuals.</i> The difference between the observed and expected summary value estimates in each cell. Available for two-way crosstabulations only and displayed only if the <b>TABLEPCT</b> keyword is specified on the <b>CELLS</b> subcommand.
<b>ASRESID</b>	<i>Adjusted Pearson residuals.</i> Available for two-way crosstabulations only and displayed only if the <b>TABLEPCT</b> keyword is specified on the <b>CELLS</b> subcommand.

## TEST Subcommand

The **TEST** subcommand requests statistics or tests for summarizing the entire table.

Furthermore, if subpopulations are defined on the **SUBPOP** subcommand using only first-stage stratification variables (or a subset of them), then tests are performed for each subpopulation also.

<b>HOMOGENEITY</b>	<i>Test of homogeneous proportions.</i> Available for one-way frequency tables only.
<b>INDEPENDENCE</b>	<i>Test of independence.</i> Available for two-way crosstabulations only.
<b>ODDSRATIO</b>	<i>Odds ratio.</i> Available for two-by-two crosstabulations only.
<b>RELRISK</b>	<i>Relative risk.</i> Available for two-by-two crosstabulations only.
<b>RISKDIFF</b>	<i>Risk difference.</i> Available for two-by-two crosstabulations only.

## SUBPOP Subcommand

The **SUBPOP** subcommand specifies subpopulations for which analyses are to be performed.

- The set of subpopulations is defined by specifying a single categorical variable, or two or more categorical variables, separated by the **BY** keyword, whose values are crossed.
- For example, `/SUBPOP TABLE = A` defines subpopulations based on the levels of variable *A*.
- For example, `/SUBPOP TABLE = A BY B` defines subpopulations based on crossing the levels of variables *A* and *B*.
- A maximum of 16 variables may be specified.
- Numeric or string variables may be specified.
- All specified variables must be unique.
- Stratification or cluster variables may be specified, but no other plan file variables are allowed on the **SUBPOP** subcommand.

- Tabulation variables may not be specified on the SUBPOP subcommand.
- The BY keyword is used to separate variables.

The DISPLAY keyword specifies the layout of results for subpopulations.

**LAYERED**            *Results for all subpopulations are displayed in the same table. This is the default.*  
**SEPARATE**         *Results for different subpopulations are displayed in different tables.*

## ***MISSING Subcommand***

The MISSING subcommand specifies how missing values are handled.

- All design variables must have valid data. Cases with invalid data for any design variable are deleted from the analysis.

The SCOPE keyword specifies which cases are used in the analyses. This specification is applied to tabulation variables but not design variables.

**TABLE**             *Each table is based on all valid data for the tabulation variable(s) used in creating the table. Tables for different variables may be based on different sample sizes. This is the default.*  
**LISTWISE**         *Only cases with valid data for all tabulation variables are used in creating the tables. Tables for different variables are always based on the same sample size.*

The CLASSMISSING keyword specifies whether user-missing values are treated as valid. This specification is applied to tabulation variables and categorical design variables (that is, strata, cluster, and subpopulation variables).

**EXCLUDE**           *Exclude user-missing values. This is the default.*  
**INCLUDE**           *Include user-missing values. Treat user-missing values as valid data.*

# CTABLES

CTABLES is available in the Custom Tables option.

*Note:* Square brackets that are used in the CTABLES syntax chart are required parts of the syntax and are not used to indicate optional elements. All subcommands except /TABLE are optional.

CTABLES

```
/FORMAT MINCOLWIDTH={DEFAULT} MAXCOLWIDTH={DEFAULT}
      {value } {value }

      UNITS={POINTS} EMPTY= {ZERO } MISSING= { '.' }
      {INCHES} {BLANK } {'chars'}
      {CM } {'chars'}
```

/VLABELS VARIABLES= varlist

```
      DISPLAY= {DEFAULT}
      {NAME }
      {LABEL }
      {BOTH }
      {NONE }
```

/MRSETS COUNTDUPLICATES= {**NO** }
 {YES}

/SMISSING {**VARIABLE**}
 {LISTWISE}

/PCOMPUTE &catname= EXPR(expression)

/PPROPERTIES &catname &catname...
 LABEL= "label"
 FORMAT= summary format ... summary format
 HIDESOURCECATS= {**NO\*\***}
 {YES }

/HIDESSMALLCOUNTS COUNT = {**5** }
 {integer}

/TABLE rows BY columns BY layers

/SLABELS POSITION= {**COLUMN**} VISIBLE= {**YES**}
 {ROW } {NO }
 {LAYER }

/CLABELS {**AUTO** }
 {ROWLABELS= {**OPPOSITE**} }
 {LAYER }
 {COLLABELS= {**OPPOSITE**} }
 {LAYER }

/CATEGORIES VARIABLES= varlist

```
      { [value, value, value...] }
      { ORDER= {A} KEY= {VALUE } MISSING= {EXCLUDE} }
      {D } {LABEL } {INCLUDE}
      {summary(varname)}
```

TOTAL= {**NO** } LABEL= "label" POSITION= {**AFTER**} EMPTY= {**INCLUDE**}
 {YES } {BEFORE} {EXCLUDE}

Explicit value lists can include SUBTOTAL='label', HSUBTOTAL='label', MISSING, OTHERNM.

/TITLES CAPTION= ['text' 'text'...]
 CORNER= ['text' 'text'...]

```

TITLE=  ['text' 'text'...]
Text can contain the symbols )DATE )TIME )TABLE

/SIGTEST TYPE= CHISQUARE  ALPHA= {0.05          }
                               {significance level}

INCLUDEMRSETS={YES**}
              {NO   }

CATEGORIES={ALLVISIBLE**}
          {SUBTOTALS  }

/COMPARETEST TYPE= {PROP}  ALPHA= {0.05          }
                 {MEAN}    {significance level}

ADJUST= {BONFERRONI}  ORIGIN=COLUMN
        {NONE        }

INCLUDEMRSETS={YES**}  MEANSVARIANCE={ALLCATS   }
              {NO   }    {TESTEDCATS}

CATEGORIES={ALLVISIBLE**}
          {SUBTOTALS  }

MERGE={NO**}
      {YES}

```

Row, column, and layer elements each have the general form

```

varname {[C]} [summary 'label' format...] {+} varname ...
        {[S]}                               {>}

```

When nesting (>) and concatenation (+) are combined, as in  $a + b > c$ , nesting occurs before concatenation; parentheses can be used to change precedence, as in  $(a + b) > c$ .

Summary functions available for all variables: COUNT ROWPCT.COUNT  
 COLPCT.COUNT TABLEPCT.COUNT SUBTABLEPCT.COUNT LAYERPCT.COUNT  
 LAYERROWPCT.COUNT LAYERCOLPCT.COUNT ROWPCT.VALIDN COLPCT.VALIDN  
 TABLEPCT.VALIDN SUBTABLEPCT.VALIDN LAYERPCT.VALIDN LAYERROWPCT.VALIDN  
 LAYERCOLPCT.VALIDN ROWPCT.TOTALN COLPCT.TOTALN TABLEPCT.TOTALN  
 SUBTABLEPCT.TOTALN LAYERPCT.TOTALN LAYERROWPCT.TOTALN  
 LAYERCOLPCT.TOTALN

Summary functions available for scale variables and for totals and subtotals of numeric variables: MAXIMUM MEAN MEDIAN MINIMUM MISSING MODE PTILE RANGE SEMEAN  
 STDDEV SUM TOTALN VALIDN VARIANCE ROWPCT.SUM COLPCT.SUM TABLEPCT.SUM  
 SUBTABLEPCT.SUM LAYERPCT.SUM LAYERROWPCT.SUM LAYERCOLPCT.SUM

Summary functions available for multiple response variables and their totals:  
 RESPONSES ROWPCT.RESPONSES COLPCT.RESPONSES TABLEPCT.RESPONSES  
 SUBTABLEPCT.RESPONSES LAYERPCT.RESPONSES LAYERROWPCT.RESPONSES  
 LAYERCOLPCT.RESPONSES ROWPCT.RESPONSES.COUNT COLPCT.RESPONSES.COUNT  
 TABLEPCT.RESPONSES.COUNT SUBTABLEPCT.RESPONSES.COUNT  
 LAYERPCT.RESPONSES.COUNT LAYERROWPCT.RESPONSES.COUNT  
 LAYERCOLPCT.RESPONSES.COUNT ROWPCT.COUNT.RESPONSES  
 COLPCT.COUNT.RESPONSES TABLEPCT.COUNT.RESPONSES  
 SUBTABLEPCT.COUNT.RESPONSES LAYERPCT.COUNT.RESPONSES LAYERROWPCT.  
 COUNT.RESPONSES LAYERCOLPCT.COUNT.RESPONSES

For unweighted summaries, prefix `U` to a function name, as in `UCOUNT`.

Formats for summaries: `COMMAw.d` `DOLLARw.d` `Fw.d` `NEGPARENw.d` `NEQUALw.d` `PARENw.d` `PCTw.d` `PCTPARENw.d` `DOTw.d` `CCA...` `CCEw.d` `Nw.d` `Ew.d` and all `DATE` formats

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

#### Release 13.0

- `HSUBTOTAL` keyword introduced on the `CATEGORIES` subcommand.

#### Release 14.0

- `INCLUDEMRSETS` keyword introduced on the `SIGTEST` and `COMPARETEST` subcommands.
- `CATEGORIES` keyword introduced on the `SIGTEST` and `COMPARETEST` subcommands.
- `MEANSVARIANCE` keyword introduced on the `COMPARETEST` subcommand.

#### Release 18.0

- `MERGE` keyword introduced on the `COMPARETEST` subcommand.
- `PCOMPUTE` and `PPROPERTIES` subcommands introduced.

#### Release 19.0

- `HIDESMALLCOUNTS` subcommand introduced.

### **Examples**

```
CTABLES /TABLE POLVIEWS [COLPCT] BY AGECAT.
```

```
CTABLES /TABLE $MLTNEWS [COUNT COLPCT] BY SEX
  /SLABELS VISIBLE=NO
  /CATEGORIES VARIABLES=SEX TOTAL=YES.
```

```
CTABLES /TABLE (CONFINAN + CONBUS + CONBUS + CONEDUC
  + CONPRESS + CONMEDIC) [COUNT ROWPCT]
  /CLABELS ROWLABELS=OPPOSITE.
```

## **Overview**

The Custom Tables procedure produces tables in one, two, or three dimensions and provides a great deal of flexibility for organizing and displaying the contents.

- In each dimension (row, column, and layer), you can stack multiple variables to concatenate tables and nest variables to create subtables. See the `TABLE` subcommand.
- You can let Custom Tables determine summary statistics according to the measurement level in the dictionary, or you can assign one or more summaries to specific variables and override the measurement level without altering the dictionary. See the `TABLE` subcommand.

- You can create multiple response sets with the `MRSETS` command and use them like ordinary categorical variables in a table expression. You can control the percentage base by choosing an appropriate summary function, and you can control with the `MRSETS` subcommand whether duplicate responses from a single respondent are counted.
- You can assign totals to categorical variables at different nesting levels to create subtable and table totals, and you can assign subtotals across subsets of the values of a variable. See the `CATEGORIES` subcommand.
- You can determine, on a per-variable basis, which categories to display in the table, including whether to display missing values and empty categories for which variable labels exist. You can also sort categories by name, label, or the value of a summary function. See the `CATEGORIES` subcommand.
- You can specify whether to show or hide summary and category labels and where to position the labels. For variable labels, you can specify whether to show labels, names, both, or neither. See the `SLABELS`, `CLABELS`, and `VLABELS` subcommands.
- You can request chi-square tests and pairwise comparisons of column proportions and means. See the `SIGTEST` and `COMPARETEST` subcommands.
- You can assign custom titles and captions (see the `TITLES` subcommand) and control what is displayed for empty cells and those for which a summary function cannot be computed. See the `FORMAT` subcommand.
- `CTABLES` ignores `SPLIT FILE` requests if layered splits (compare groups in the graphical user interface) are requested. You can compare groups by using the split variables at the highest nesting level for row variables. See the `TABLE` subcommand for nesting variables.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

## **Syntax Conventions**

- The basic specification is a `TABLE` subcommand with at least one variable in one dimension. Multiple `TABLE` subcommands can be included in one `CTABLES` command.
- The global subcommands `FORMAT`, `VLABELS`, `MRSETS`, and `SMISSING` must precede the first `TABLE` subcommand and can be named in any order.
- The local subcommands `SLABELS`, `CLABELS`, `CATEGORIES`, `TITLES`, `SIGTEST`, and `COMPARETEST` follow the `TABLE` subcommand in any order and refer to the immediately preceding table expression.
- The `HIDESMALLCOUNTS` subcommand applies to all `TABLE` subcommands and must precede the first `TABLE` subcommand.
- In general, if subcommands are repeated, their specifications are merged. The last value of each specified attribute is honored.
- Equals signs that are shown in the syntax charts are required.
- Square brackets that are shown in the syntax charts are required.

- All keywords except summary function names, attribute values, and explicit category list keywords can be truncated to as few as three characters. Function names must be spelled in full.
- The slash before all subcommands, including the first subcommand, is required.

## Examples

### Example: Column Percentages

```
CTABLES /TABLE POLVIEWS [COLPCT] BY AGECAT.
```

Figure 50-1

		Age category					
		Less than 25	25 to 34	35 to 44	45 to 54	55 to 64	65 or older
		Column N %	Column N %	Column N %	Column N %	Column N %	Column N %
Think of self as liberal or conservative	Extremely liberal	4.5%	2.5%	2.1%	2.4%	1.3%	2.2%
	Liberal	18.8%	15.7%	14.6%	11.3%	10.5%	9.4%
	Slightly liberal	13.5%	14.2%	13.2%	15.4%	10.5%	10.5%
	Moderate	36.8%	37.1%	32.7%	37.2%	39.3%	38.8%
	Slightly conservative	14.3%	14.9%	19.3%	15.0%	18.4%	13.4%
	Conservative	11.7%	13.0%	14.6%	15.4%	16.4%	21.2%
	Extremely conservative	4%	2.7%	3.5%	3.3%	3.6%	4.5%

- *POLVIEWS* defines the rows, and *AGECAT* defines the columns. Column percentages are requested, overriding the default *COUNT* function.

### Example: Using a Multiple Response Set

```
CTABLES /TABLE $MLTNEWS [COUNT COLPCT] BY SEX
/SLABELS VISIBLE=NO
/CATEGORIES VARIABLES=SEX TOTAL=YES.
```

Figure 50-2

		Gender					
		Male		Female		Total	
News sources	Get news from newspapers	375	41.9%	430	36.3%	805	38.7%
	Get news from news magazines	121	13.5%	173	14.6%	294	14.1%
	Get news from television	451	50.3%	626	52.8%	1077	51.8%
	Get news from radio	233	26.0%	318	26.8%	551	26.5%
	Get news from internet	359	40.1%	508	42.9%	867	41.7%

- *\$MLTNEWS* is a multiple response set.
- The *COLPCT* function uses the number of respondents as the percentage base, so each cell shows the percentage of males or females who gave each response, and the sum of percentage for each column is greater than 100.
- Summary labels are hidden.
- The *CATEGORIES* subcommand creates a total for both sexes.

### Example: Concatenation

```
CTABLES /TABLE (CONFINAN + CONBUS + CONBUS + CONEDUC
+ CONPRESS + CONMEDIC) [COUNT ROWPCT]
```



```
/CLABELS ROWLABELS=OPPOSITE.
```

Figure 50-3

	A great deal		Only some		Hardly any	
	Count	Row N %	Count	Row N %	Count	Row N %
Confidence in banks & financial institutions	490	26.3%	1068	57.3%	306	16.4%
Confidence in major companies	500	27.5%	1078	59.2%	243	13.3%
Confidence in major companies	500	27.5%	1078	59.2%	243	13.3%
Confidence in education	511	27.2%	1055	56.1%	315	16.7%
Confidence in press	176	9.5%	878	47.2%	808	43.4%
Confidence in medicine	844	45.0%	864	46.1%	167	8.9%

- The six confidence variables all have the same categories with the same value labels for each category.
- The CLABELS subcommand moves the category labels to the columns.

## TABLE Subcommand

The TABLE subcommand specifies the structure of the table, including the variables and summary functions that define each dimension. The TABLE subcommand has the general form

```
/TABLE rows BY columns BY layers
```

The minimum specification for a row, column, or layer is a variable name. You can specify one or more dimensions.

See [SET TABLERENDER](#) for information on how multiple layer dimensions are rendered in lightweight tables.

## Variable Types

The variables that are used in a table expression can be category variables, scale variables, or multiple response sets. Multiple response sets are defined by the MRSETS command and always begin with a \$. Custom Tables uses the measurement level in the dictionary for the active data file to identify category and scale variables. You can override the default variable type for numeric variables by placing [C] or [S] after the variable name. Thus, to treat the category variable HAPPY as a scale variable and obtain a mean, you would specify

```
/TABLE HAPPY [S].
```

## Category Variables and Multiple Response Sets

Category variables define one cell per value. See the CATEGORIES subcommand for ways of controlling how categories are displayed. Multiple response sets also define one cell per value.

### Example

```
CTABLES /TABLE HAPPY.
```

Figure 50-4

		Count
General happiness	Very happy	891
	Pretty happy	1575
	Not too happy	340

- The counts for *HAPPY* are in the rows.

**Example**

CTABLES /TABLE BY HAPPY.

Figure 50-5

General happiness		
Very happy	Pretty happy	Not too happy
Count	Count	Count
891	1575	340

- The counts for *HAPPY* are in the columns.

**Example**

CTABLES /TABLE BY BY HAPPY

Figure 50-6

General happiness Very happy	
Count	
891	

- The counts for *HAPPY* are in layers.

**Stacking and Nesting**

Stacking (or concatenating) variables creates multiple logical tables within a single table structure.

**Example**

CTABLES /TABLE HAPPY + HAPMAR BY CHILDCAT.

Figure 50-7

		Number of children (grouped categories)			
		None	1-2	3-4	5 or more
		Count	Count	Count	Count
General happiness	Very happy	197	412	221	59
	Pretty happy	499	662	314	97
	Not too happy	98	136	79	27
Happiness of marriage	Very happy	111	462	232	49
	Pretty happy	51	238	133	22
	Not too happy	5	18	10	4

- The output contains two tables: one table for general happiness by number of children and one table for happiness in marriage by number of children. Except for missing values, all of the cases in the data appear in both tables.

Nesting variables creates hierarchical tables.

### Example

```
CTABLES /TABLE SEX > HAPMAR BY CHILDCAT.
```

Figure 50-8

				Number of children (grouped categories)			
				None	1-2	3-4	5 or more
				Count	Count	Count	Count
Gender	Male	Happiness	Very happy	48	216	102	30
		of marriage	Pretty happy	25	110	58	11
			Not too happy	3	7	4	1
Female	Happiness	of marriage	Very happy	63	246	130	19
			Pretty happy	26	128	75	11
			Not too happy	2	11	6	3

- The output contains one table with a subtable for each value of *SEX*. The same subtables would result from the table expression `HAPMAR BY CHILDCAT BY SEX`, but the subtables would appear in separate layers.

Stacking and nesting can be combined. When they are combined, by default, nesting takes precedence over stacking. You can use parentheses to alter the order of operations.

### Example

```
CTABLES /TABLE (HAPPY + HAPMAR) > SEX.
```

Figure 50-9

				Count
General happiness	Very happy	Gender	Male	373
			Female	518
	Pretty happy	Gender	Male	712
			Female	863
	Not too happy	Gender	Male	133
			Female	207
Happiness of marriage	Very happy	Gender	Male	396
			Female	459
	Pretty happy	Gender	Male	205
			Female	240
	Not too happy	Gender	Male	15
			Female	22

- The output contains two tables. Without the parentheses, the first table, for general happiness, would not have separate rows for male and female.

## Scale Variables

Scale variables, such as *age in years* or *population of towns*, do not define multiple cells within a table. The table expression `/TABLE AGE` creates a table with one cell containing the mean of *AGE* across all cases in the data. You can use nesting and/or dimensions to display summary

statistics for scale variables within categories. The nature of scale variables prevents their being arranged hierarchically. Therefore:

- A scale variable cannot be nested under another scale variable.
- Scale variables can be used in only one dimension.

### Example

```
CTABLES /TABLE AGE > HAPPY BY SEX.
```

Figure 50-10

			Gender	
			Male	Female
			Mean	Mean
Age of respondent	General happiness	Very happy	47	47
		Pretty happy	44	45
		Not too happy	43	47

## Specifying Summaries

You can specify one or more summary functions for variables in any one dimension. For category variables, summaries can be specified only for the variables at the lowest nesting level. Thus, in the table expression

```
/TABLE SEX > (HAPPY + HAPMAR) BY AGECAT
```

you can assign summaries to *HAPPY* and *HAPMAR* or to *AGECAT*, but not to both and not to *SEX*.

If a scale variable appears in a dimension, that dimension becomes the statistics dimension, and all statistics must be specified for that dimension. A scale variable need not be at the lowest level of nesting. Thus, the following is a valid specification:

```
CTABLES /TABLE AGE [MINIMUM, MAXIMUM, MEAN] > SEX > HAPPY.
```

A multiple response variable also need not be at the lowest level of nesting. The following specification is a valid specification:

```
CTABLES /TABLE $MLTCARS [COUNT, RESPONSES] > SEX.
```

However, if two multiple response variables are nested, as in *\$MULTCARS* > *\$MULTNEWS*, summaries can be requested only for the variable at the innermost nesting level (in this case, *\$MULTNEWS*).

The general form for a summary specification is

```
[summary 'label' format, ..., summary 'label' format]
```

- The specification follows the variable name in the table expression. You can apply a summary specification to multiple variables by enclosing the variables in parentheses. The following specifications are equivalent:

```
/TABLE SEX [COUNT] + HAPPY [COUNT, COLPCT]
/TABLE (SEX + HAPPY [COLPCT])[COUNT]
```

- The brackets are required even if only one summary is specified.
- Commas are optional.
- Label and format are both optional; defaults are used if label and format are not specified.
- If totals or subtotals are defined for a variable (on the CATEGORIES subcommand), by default, the same functions that are specified for the variable are used for the totals. You can use the keyword TOTALS within the summary specification to specify different summary functions for the totals and subtotals. The specification then has the form [summary 'label' format ... TOTALS [summary 'label' format...]]. You must still specify TOTAL=YES on the CATEGORIES subcommand to see the totals.
- Summaries that are available for category variables are also available for scale variables and multiple response sets. Functions that are specific to scale variables and to multiple response sets are also available.
- If case weighting is in effect, summaries are calculated taking into account the current WEIGHT value. To obtain unweighted summaries, prefix a U to the function name, as in UCOUNT. Unweighted functions are not available where weighting would not apply, as in the MINIMUM and MAXIMUM functions.

### Example

```
CTABLES /TABLE SEX > HAPMAR [COLPCT] BY CHILDCAT.
```

Figure 50-11

				Number of children (grouped categories)			
				None	1-2	3-4	5 or more
				Column N %	Column N %	Column N %	Column N %
Gender	Male	Happiness of marriage	Very happy	63.2%	64.9%	62.2%	71.4%
			Pretty happy	32.9%	33.0%	35.4%	26.2%
		Not too happy	3.9%	2.1%	2.4%	2.4%	
Female	Happiness of marriage	Very happy	69.2%	63.9%	61.6%	57.6%	
		Pretty happy	28.6%	33.2%	35.5%	33.3%	
		Not too happy	2.2%	2.9%	2.8%	9.1%	

### Example

```
CTABLES /TABLE AGE CAT > TVHOURS [MEAN F5.2,
STDDEV 'Standard Deviation' F5.2, PTILE 90 '90th Percentile'].
```

Figure 50-12

			Mean	Standard Deviation	90th Percentile
Age category	Less than 25	Hours per day watching TV	2.85	2.03	5
	25 to 34	Hours per day watching TV	2.78	2.37	5
	35 to 44	Hours per day watching TV	2.56	2.11	5
	45 to 54	Hours per day watching TV	2.58	1.97	5
	55 to 64	Hours per day watching TV	3.02	2.22	6
	65 or older	Hours per day watching TV	3.58	2.50	6

- Each summary function for the row variable appears by default in a column.

- Labels for standard deviation and the 90th percentile override the defaults.
- Because *TVHOURS* is recorded in whole hours and has an integer print format, the default general print formats for mean and standard deviation would also be integer, so overrides are specified.

Table 50-1  
Summary functions: all variables

Function	Description	Default Label*	Default Format
COUNT	Number of cases in each category. This is the default for categorical and multiple response variables.	Count	Count
ROWPCT . COUNT	Row percentage based on cell counts. Computed within subtable.	Row %	Percent
COLPCT . COUNT	Column percentage based on cell counts. Computed within subtable.	Column %	Percent
TABLEPCT . COUNT	Table percentage based on cell counts.	Table %	Percent
SUBTABLEPCT . COUNT	Subtable percentage based on cell counts.	Subtable %	Percent
LAYERPCT . COUNT	Layer percentage based on cell counts. Same as table percentage if no layers are defined.	Layer %	Percent
LAYERROWPCT . COUNT	Row percentage based on cell counts. Percentages sum to 100% across the entire row (that is, across subtables).	Layer Row %	Percent
LAYERCOLPCT . COUNT	Column percentage based on cell counts. Percentages sum to 100% across the entire column (that is, across subtables).	Layer Column %	Percent
ROWPCT . VALIDN	Row percentage based on valid count.	Row Valid N %	Percent
COLPCT . VALIDN	Column percentage based on valid count.	Column Valid N %	Percent
TABLEPCT . VALIDN	Table percentage based on valid count.	Table Valid N %	Percent
SUBTABLEPCT . VALIDN	Subtable percentage based on valid count.	Subtable Valid N %	Percent
LAYERPCT . VALIDN	Layer percentage based on valid count.	Layer Valid N %	Percent
LAYERROWPCT . VALIDN	Row percentage based on valid count. Percentages sum to 100% across the entire row.	Layer Row Valid N %	Percent
LAYERCOLPCT . VALIDN	Column percentage based on valid count. Percentages sum to 100% across the entire column.	Layer Column Valid N %	Percent
ROWPCT . TOTALN	Row percentage based on total count, including user-missing and system-missing values.	Row Total N %	Percent
COLPCT . TOTALN	Column percentage based on total count, including user-missing and system-missing values.	Column Total N %	Percent

Function	Description	Default Label*	Default Format
TABLEPCT . TOTALN	Table percentage based on total count, including user-missing and system-missing values.	Table Total N %	Percent
SUBTABLEPCT . TOTALN	Subtable percentage based on total count, including user-missing and system-missing values.	Subtable Total N %	Percent
LAYERPCT . TOTALN	Layer percentage based on total count, including user-missing and system-missing values.	Layer Total N %	Percent
LAYERROWPCT . TOTALN	Row percentage based on total count, including user-missing and system-missing values. Percentages sum to 100% across the entire row.	Layer Row Total N %	Percent
LAYERCOLPCT . TOTALN	Column percentage based on total count, including user-missing and system-missing values. Percentages sum to 100% across the entire column.	Layer Column Total N %	Percent

\* This is the default on a U.S.-English system.

The .COUNT suffix can be omitted from percentages that are based on cell counts. Thus, ROWPCT is equivalent to ROWPCT .COUNT.

**Table 50-2**  
Summary functions: scale variables, totals, and subtotals

Function	Description	Default Label	Default Format
MAXIMUM	Largest value.	Maximum	General
MEAN	Arithmetic mean. The default for scale variables.	Mean	General
MEDIAN	50th percentile.	Median	General
MINIMUM	Smallest value.	Minimum	General
MISSING	Count of missing values (both user-missing and system-missing).	Missing	General
MODE	Most frequent value. If there is a tie, the smallest value is shown.	Mode	General
PTILE	Percentile. Takes a numeric value between 0 and 100 as a required parameter.  PTILE is computed the same way as APTILE in the TABLES command. Note that in the TABLES command, the default percentile method was HPTILE.	Percentile #####.##	General
RANGE	Difference between maximum and minimum values.	Range	General
SEMEAN	Standard error of the mean.	Std Error of Mean	General
STDDEV	Standard deviation.	Std Deviation	General
SUM	Sum of values.	Sum	General

## CTABLES

Function	Description	Default Label	Default Format
TOTALN	Count of nonmissing, user-missing, and system-missing values. The count excludes valid values hidden via the CATEGORIES subcommand.	Total N	Count
VALIDN	Count of nonmissing values.	Valid N	Count
VARIANCE	Variance.	Variance	General
ROWPCT . SUM	Row percentage based on sums.	Row Sum %	Percent
COLPCT . SUM	Column percentage based on sums.	Column Sum %	Percent
TABLEPCT . SUM	Table percentage based on sums.	Table Sum %	Percent
SUBTABLEPCT . SUM	Subtable percentage based on sums.	Subtable Sum %	Percent
LAYERPCT . SUM	Layer percentage based on sums.	Layer Sum %	Percent
LAYERROWPCT . SUM	Row percentage based on sums. Percentages sum to 100% across the entire row.	Layer Row Sum %	Percent
LAYERCOLPCT . SUM	Column percentage based on sums. Percentages sum to 100% across the entire column.	Layer Column Sum %	Percent

Table 50-3  
Summary functions: multiple response sets

Function	Description	Default Label	Default Format
RESPONSES	Count of responses.	Responses	Count
ROWPCT . RESPONSES	Row percentage based on responses. Total number of responses is the denominator.	Row Responses %	Percent
COLPCT . RESPONSES	Column percentage based on responses. Total number of responses is the denominator.	Column Responses %	Percent
TABLEPCT . RESPONSES	Table percentage based on responses. Total number of responses is the denominator.	Table Responses %	Percent
SUBTABLEPCT . RESPONSES	Subtable percentage based on responses. Total number of responses is the denominator.	Subtable Responses %	Percent
LAYERPCT . RESPONSES	Layer percentage based on responses. Total number of responses is the denominator.	Layer Responses %	Percent
LAYERROWPCT . RESPONSES	Row percentage based on responses. Total number of responses is the denominator.  Percentages sum to 100% across the entire row (that is, across subtables).	Layer Row Responses %	Percent
LAYERCOLPCT . RESPONSES	Column percentage based on responses. Total number of responses is the denominator.  Percentages sum to 100% across the entire column (that is, across subtables).	Layer Column Responses %	Percent



<b>Function</b>	<b>Description</b>	<b>Default Label</b>	<b>Default Format</b>
ROWPCT . RESPONSES . COUNT	Row percentage: Responses are the numerator, and total count is the denominator.	Row Responses % (Base: Count)	Percent
COLPCT . RESPONSES . COUNT	Column percentage: Responses are the numerator, and total count is the denominator.	Column Responses % (Base: Count)	Percent
TABLEPCT . RESPONSES . COUNT	Table percentage: Responses are the numerator, and total count is the denominator.	Table Responses % (Base: Count)	Percent
SUBTABLEPCT . RE- SPONSES . COUNT	Subtable percentage: Responses are the numerator, and total count is the denominator.	Subtable Responses % (Base: Count)	Percent
LAYERPCT . RESPONSES . COUNT	Layer percentage: Responses are the numerator, and total count is the denominator.	Layer Responses % (Base: Count)	Percent
LAYERROWPCT . RE- SPONSES . COUNT	Row percentage: Responses are the numerator, and total count is the denominator.  Percentages sum to 100% across the entire row (that is, across subtables).	Layer Row Responses % (Base: Count)	Percent
LAYERCOLPCT . RE- SPONSES . COUNT	Column percentage: Responses are the numerator, and total count is the denominator.  Percentages sum to 100% across the entire column (that is, across subtables).	Layer Column Responses % (Base: Count)	Percent
ROWPCT . COUNT . RESPONSES	Row percentage: Count is the numerator, and total responses are the denominator.	Row Count % (Base: Responses)	Percent
COLPCT . COUNT . RESPONSES	Column percentage: Count is the numerator, and total responses are the denominator.	Column Count % (Base: Responses)	Percent
TABLEPCT . COUNT . RESPONSES	Table percentage: Count is the numerator, and total responses are the denominator.	Table Count % (Base: Responses)	Percent
SUBTABLEPCT . COUNT . RESPONSES	Subtable percentage: Count is the numerator, and total responses are the denominator.	Subtable Count % (Base: Responses)	Percent
LAYERPCT . COUNT . RESPONSES	Layer percentage: Count is the numerator, and total responses are the denominator.	Layer Count % (Base: Responses)	Percent

## CTABLES

Function	Description	Default Label	Default Format
LAYERROWPCT.COUNT.RE- SPONSES	Row percentage: Count is the numerator, and total responses are the denominator.  Percentages sum to 100% across the entire row (that is, across subtables).	Layer Row Count % (Base: Responses)	Percent
LAYERCOLPCT.COUNT.RE- SPONSES	Row percentage: Count is the numerator, and total responses are the denominator.  Percentages sum to 100% across the entire column (that is, across subtables).	Layer Column Count % (Base: Responses)	Percent

### Formats for Summaries

A default format is assigned to each summary function:

<b>Count</b>	The value is expressed in F (standard numeric) format with 0 decimal places. If you have fractional weights and want a count that reflects those weights, use F format with appropriate decimal places.
<b>Percent</b>	The value is expressed with one decimal place and a percent symbol.
<b>General</b>	The value is expressed in the variable's print format.

These default formats are internal to CTABLES and cannot be used in table expressions. To override the default formats, use any of the print formats that are available in the Core system except Z, PBHEX, and HEX, or use the additional formats that are described in the following table.

Table 50-4  
Additional formats for summaries

Format	Description	Example
NEGPARENw.d	Parentheses appear around negative numbers.	-1234.567 formatted as NEGPAREN9.2 yields (1234.57).
NEQUALw.d	"N=" precedes the number.	1234.567 formatted as NEQUAL9.2 yields N=1234.57.
PARENw.d	The number is parenthesized.	1234.567 formatted as PAREN8.2 yields (1234.57).
PCTPARENw.d	A percent symbol follows the parenthesized value.	1234.567 formatted as PCTPAREN10.2 yields (1234.57%).

### Missing Values in Summaries

The following table presents the rules for including cases in a table for VALIDN, COUNT, and TOTALN functions when values are included or excluded explicitly through an explicit category list or implicitly through inclusion or exclusion of user-missing values.

Table 50-5  
Inclusion/exclusion of values in summaries

Variable and Value Type	VALIDN	COUNT	TOTALN
Categorical Variable: shown valid value Multiple Dichotomy Set: at least one “true” value Multiple Category Set: at least one shown valid value Scale Variable: valid value	Include	Include	Include
Categorical Variable: included user-missing value Multiple Category Set: all values are included user-missing Scale Variable: user-missing or system-missing	Exclude	Include	Include
Categorical Variable: excluded user-missing or system-missing value Multiple Dichotomy Set: all values are “false” Multiple Category Set: all values are excluded user-missing, system-missing, or excluded valid, but at least one value is not excluded valid	Exclude	Exclude	Include
Categorical Variable: excluded valid value Multiple Dichotomy Set: all values are excluded valid values	Exclude	Exclude	Exclude

## SLABELS Subcommand

The SLABELS subcommand controls the position of summary statistics in the table and controls whether summary labels are shown.

```
/SLABELS POSITION= {COLUMN} VISIBLE= {YES}
                  {ROW   }          {NO  }
                  {LAYER }
```

By default, summaries appear in the columns and labels are visible.

### Example: Summary Label Positioning

```
CTABLES /TABLE NEWS [COUNT COLPCT].
```

Figure 50-13

		Count	Column N %
How often does respondent read newspaper	Every day	805	43.0%
	Few times a week	420	22.5%
	Once a week	294	15.7%
	Less than once a week	202	10.8%
	Never	149	8.0%

```
CTABLES /TABLE NEWS [COUNT COLPCT]
/SLABELS POSITION=ROW VISIBLE=NO.
```

Figure 50-14

How often does respondent read newspaper	Every day	805 43.0%
	Few times a week	420 22.5%
	Once a week	294 15.7%
	Less than once a week	202 10.8%
	Never	149 8.0%

## CLABELS Subcommand

The CLABELS subcommand controls the location of category labels.

```
/CLABELS {AUTO           }
          {ROWLABELS= {OPPOSITE} }
          {LAYER       }
          {COLLABELS= {OPPOSITE} }
          {LAYER       }
```

By default, category labels are nested under the variables to which they belong. Category labels for row and column variables can be moved to the opposite dimension or to the layers. If labels exist in both dimensions, only one dimension, row labels or column labels, can be moved; they cannot be swapped.

### Example

```
CTABLES
  /TABLE (CONFINAN + CONEDUC + CONBUS + CONMEDIC + CONPRESS + CONTV )
```

Figure 50-15

		Count
Confidence in banks & financial institutions	A great deal	490
	Only some	1068
	Hardly any	306
Confidence in education	A great deal	511
	Only some	1055
	Hardly any	315
Confidence in major companies	A great deal	500
	Only some	1078
	Hardly any	243
Confidence in medicine	A great deal	844
	Only some	864
	Hardly any	167
Confidence in press	A great deal	176
	Only some	878
	Hardly any	808
Confidence in television	A great deal	196
	Only some	936
	Hardly any	744

- Six variables are stacked in the rows, and their category labels are stacked under them.

```
CTABLES
  /TABLE (CONFINAN + CONEDUC + CONBUS + CONMEDIC + CONPRESS + CONTV )
  /SLABELS VISIBLE=NO /CLABELS ROWLABELS=OPPOSITE
```

Figure 50-16

	A great	Only some	Hardly any
Confidence in banks & financial institutions	490	1068	306
Confidence in education	511	1055	315
Confidence in major companies	500	1078	243
Confidence in medicine	844	864	167
Confidence in press	176	878	808
Confidence in television	196	936	744

- The category labels are moved to the columns. Where variables are stacked, as in this example, the value labels for all of the variables must be exactly the same to allow for this format. Additionally, all must have the same category specifications, and data-dependent sorting is not allowed.

## CATEGORIES Subcommand

The CATEGORIES subcommand controls the order of categories in the rows and columns of the table, controls the showing and hiding of ordinary and user-missing values, and controls the computation of totals and subtotals.

```
/CATEGORIES VARIABLES= varlist

{ [value, value, value...]
{ ORDER= {A} KEY= {VALUE           } MISSING= {EXCLUDE} }
  {D}      {LABEL           }      {INCLUDE}
          {summary (varname)}

TOTAL= {NO  } LABEL= "label" POSITION= {AFTER } EMPTY= {INCLUDE}
       {YES }                       {BEFORE} {EXCLUDE}
```

The minimum specification is a variable list and one of the following specifications: a category specification, TOTAL specification, or EMPTY specification. The variable list can be a list of variables or the keyword ALL, which refers to all category variables in the table expression. ALL cannot be used with the explicit category list.

## Explicit Category Specification

The explicit category specification is a bracketed list of data values or value ranges in the order in which they are to be displayed in the table. Values not included in the list are excluded from the table. This form allows for subtotals and showing or hiding of specific values (both ordinary and user-missing).

- The list can include both ordinary and user-missing values but not the system-missing value (.).
- Values are optionally separated by commas.
- String and date values must be quoted. Date values must be consistent with the variable's print format.
- The LO, THRU, and HI keywords can be used in the value list to refer to a range of categories. LO and HI can be used only as part of a range specification.
- The MISSING keyword can be used to refer to all user-missing values.

- The `OTHERNM` keyword can be used to refer to all nonmissing values that are not explicitly named in the list. The keyword can be placed anywhere within the list. The values to which it refers appear in ascending order.
- If a value is repeated in the list, the last instance is honored. Thus, for a variable `RATING` with integer values 1 through 5, the following specifications are equal:

```
/CATEGORIES VARIABLES = RATING [1,2,4,5,3]
/CATEGORIES VARIABLES = RATING [1 THRU 5,3]
/CATEGORIES VARIABLES = RATING [OTHERNM,3]
```

- For a multiple dichotomy set, you can order the variables in the set by using the names of the variables in the set. The variable names are not enclosed in quotation marks.
- The `SUBTOTAL` keyword is used within a category list to request subtotals for a variable. The position of a subtotal within the list determines where it will appear in the table and the categories to which it applies. By default, a subtotal applies to all values that precede it up to the next subtotal. If `POSITION=BEFORE` is specified (For more information, see the topic [Totals on p. 531.](#)), subtotals apply to the categories that follow them in the list. Hierarchical and overlapping subtotals are not supported. You can specify a label for a subtotal by placing the label in quotation marks immediately following the `SUBTOTAL` keyword and an equals sign, as illustrated in the following example:

### Example

```
CTABLES /TABLE AGE CAT
/CATEGORIES VARIABLES=AGE CAT [1, 2, 3, SUBTOTAL='Subtotal < 45',
4, 5, 6, SUBTOTAL='Subtotal 45+'].
```

Figure 50-17

		Count
Age	Less than 25	242
category	25 to 34	627
	35 to 44	679
	Subtotal < 45	1548
	45 to 54	481
	55 to 64	320
	65 or older	479
	Subtotal 45+	1280

- The `HSUBTOTAL` keyword functions just like the `SUBTOTAL` keyword, except that only the subtotal is displayed in the table; the categories that define the subtotal are not included in the table. So you can use `HSUBTOTAL` to collapse categories in a table without recoding the original variables.

### Example

```
CTABLES /TABLE AGE CAT
/CATEGORIES VARIABLES=AGE CAT [1, 2, 3, HSUBTOTAL='Under 45',
4, 5, 6, HSUBTOTAL='45 or older']..
```

Figure 50-18

		Count
Age category	Under 45	1548
	45 or older	1280

## Implicit Category Specification

The implicit list allows you to sort the categories and to show or hide user-missing values without having to enumerate the values. The implicit list also provides for data-dependent sorting. If you do not supply an explicit value list, you can use the following keywords:

<b>ORDER</b>	<i>The sorting order.</i> You can select <code>A</code> (the default) for ascending order, or <code>D</code> for descending order.
<b>KEY</b>	<i>The sort key.</i> You can specify <code>VALUE</code> (the default) to sort by the values or <code>LABEL</code> to sort by the value labels. When values are sorted by label, any unlabeled values appear after the labeled values in the table. You can also specify a summary function for data-dependent sorting.
<b>MISSING</b>	<i>Whether user-missing values are included.</i> You can specify <code>EXCLUDE</code> (the default) or <code>INCLUDE</code> . System-missing values are never included.

**Data-Dependent Sorting.** The following conventions and limitations apply to sorting by using a summary function as the key:

- The sort function must be a summary function that is supported in `CTABLES`.
- The sort function must be used in the table. The exception to this rule is `COUNT`. You can sort by `COUNT` even if counts do not appear in the table.
- Data-dependent sorting is not available if category labels are repositioned by using the `CLABELS` subcommand.
- Summary functions that are available only for scale variables require that you give the variable name in parentheses, as in `MEAN(age)`. For percentiles, the variable name must be followed by a comma and an integer value between 0 and 100, as in `PTILE(age, 75)`. Other functions, such as `COUNT`, do not require a variable name, but you can supply a variable name to restrict the sort.
- When a variable name is given, and multiple logical tables are created through stacking, the entire table is sorted based on the first logical table that includes the categorical variable that is being sorted and the variable that is specified in the key.
- When a table contains more than one dimension, the sort is based on the distribution of the key within the categories of the sorted variable, without regard to the contents of the other dimensions. Thus, given the table

```
CTABLES /TABLE A BY B + C /CAT VAR=A ORDER=A KEY=COUNT(A) ,
```

the rows are sorted according to the counts for the categories of *A*, without regard to the values of *B* and *C*. If there are no missing values in the other dimension, the result is the same as sorting on the totals for that dimension (in this case, *B* or *C*). If the other dimension has an unbalanced pattern of missing values, the sorting may give unexpected results; however, the result is unaffected by differences in the pattern for *B* and *C*.

- If the sort variable is crossed with stacked category variables, the first table in the stack determines the sort order.
- To ensure that the categories are sorted the same way in each layer of the pivot table, layer variables are ignored for the purpose of sorting.

**Example**

```

CTABLES
  /TABLE CAR1 BY AGECAT
  /CATEGORIES VARIABLES=AGECAT TOTAL=YES
  /CATEGORIES VARIABLES=CAR1 ORDER=D KEY=COUNT.

```

Figure 50-19

		Age category						Total
		Less than 25	25 to 34	35 to 44	45 to 54	55 to 64	65 or older	
		Count	Count	Count	Count	Count	Count	
Car maker, most recent car	American	99	267	293	214	140	215	1228
	Japanese	73	136	140	107	66	104	626
	German	18	91	69	63	36	61	338
	Korean	23	77	88	45	35	50	318
	Swedish	18	32	46	20	24	25	165
	Other	11	24	43	32	19	24	153

- The first CATEGORIES subcommand requests a total across all age categories.
- The second CATEGORIES subcommand requests a sort of the categories of *CAR1* in descending order (using COUNT as the key). The categories of *CAR1* are sorted according to the total counts.

**Example**

```

CTABLES
  /TABLE AGE [MEAN F5.1] > CAR1 BY SEX
  /CATEGORIES VARIABLES=SEX TOTAL=YES
  /CATEGORIES VARIABLES=CAR1 KEY=MEAN(AGE) .

```

Figure 50-20

			Gender		
			Male	Female	Total
			Mean	Mean	Mean
Age of respondent	Car maker, most recent car	Swedish	42.6	45.6	44.3
		Japanese	43.5	45.5	44.7
		Korean	43.4	46.2	45.0
		American	45.3	46.5	45.9
		German	44.3	47.6	46.2
		Other	48.6	46.4	47.3

- The first CATEGORIES subcommand requests a total across the values of *SEX*.
- The second CATEGORIES subcommand requests that the categories of *CAR1* be sorted according to the mean of *AGE*. The categories are sorted according to the total means for both sexes, and that would be the case if the totals were not shown in the table.



## Totals

A total can be specified for any category variable regardless of its level of nesting within a dimension. Totals can be requested in more than one dimension. The following options are available:

<b>TOTAL</b>	<i>Whether to display a total for a variable.</i> You can specify <code>TOTAL=NO</code> (the default) or <code>TOTAL=YES</code> .
<b>LABEL</b>	<i>The label for the total.</i> The specification is a quoted string.
<b>POSITION</b>	<i>Whether a total comes after or before the categories of the variable being totaled.</i> You can specify <code>AFTER</code> (the default) or <code>BEFORE</code> . <code>POSITION</code> also determines whether subtotals that are specified in an explicit list of categories apply to the categories that precede them ( <code>AFTER</code> ) or follow them ( <code>BEFORE</code> ).

Scale variables cannot be totaled directly. To obtain a total or subtotals for a scale variable, request the total or subtotals for the category variable within whose categories the summaries for the scale variable appear.

### Example

```
CTABLES /TABLE AGECAT
  /CATEGORIES VARIABLES=AGECAT TOTAL=YES LABEL='Total Respondents'.
```

Figure 50-21

		Count
Age	Less than 25	242
category	25 to 34	627
	35 to 44	679
	45 to 54	481
	55 to 64	320
	65 or older	479
Total Respondents		2828

### Example

```
CTABLES /TABLE AGE [MEAN 'Average' F5.1] > SEX
  /CATEGORIES VARIABLES=SEX TOTAL=YES LABEL='Combined'.
```

Figure 50-22

			Average
Age of respondent	Gender	Male	44.6
		Female	46.3
	Combined		45.6

- The summary function for *AGE* appears in cells that are determined by the values of *SEX*. The total is requested for *SEX* to obtain the average age across both sexes.

## Empty Categories

Empty categories are those categories for which no cases appear in the data. For an explicit category list, this includes all explicitly named values and all labeled values that are implied by THRU, OTHERNM, or MISSING. For an implicit category list, this includes all values for which value labels exist.

**EMPTY**            *Whether to show categories whose count is zero.* You can specify `EMPTY=INCLUDE` (the default) or `EMPTY=EXCLUDE`.

## TITLES Subcommand: Titles, Captions, and Corner Text

The TITLES subcommand specifies table annotations. If the subcommand is used, a title, caption, or corner text must be specified. No caption, title, or corner text is displayed by default.

```
/TITLES  CAPTION= ['text' 'text'...]
         CORNER=  ['text' 'text'...]
         TITLE=   ['text' 'text'...]
```

**CAPTION**            *Caption lines.* The caption appears below the table. Multiple lines can be specified. Each line must be quoted.

**CORNER**            *Corner text.* Corner text appears in the corner cell of the table, above row titles and next to column titles. Multiple lines can be specified. Each line must be quoted. Pivot tables show all corner text that fits in the corner cell. The specified text is ignored if the table has no corner cell.

The system default TableLook uses the corner area for display of row dimension labels. To display CTABLES corner text, the Row Dimension Labels setting in Table Properties should be set to Nested. This choice can be preset in the default TableLook.

**TITLE**              *Title text.* The title appears above the table. Multiple lines can be specified. Each line must be quoted.

The following symbols can be used within any caption, corner text, or title line. Each symbol must be specified by using an opening right parenthesis and all uppercase letters.

**)DATE**              *Current date.* Displays a locale-appropriate date stamp that includes the year, month, and day.

**)TIME**              *Current time.* Displays a locale-appropriate time stamp.

**)TABLE**            *Table description.* Inserts a description of the table, which consists of the table expression stripped of measurement levels, statistics specifications, and /TABLE. If variable labels are available, they are used instead of variable names in the table expression.

### Example

```
CTABLES /VLABELS VARIABLES=SEX HAPMAR DISPLAY=NONE
/TABLE SEX > HAPMAR BY CHILDCAT [COLPCT]
/SLABELS VISIBLE=NO
/TITLE TITLE = 'Marital Happiness for Men and Women '+
              'by Number of Children'
CAPTION= 'Report created at )TIME on )DATE' ')TABLE'.
```

Figure 50-23

**Marital Happiness for Men and Women by Number of Children**

		Number of children (grouped categories)			
		None	1-2	3-4	5 or more
Male	Very happy	63.2%	64.9%	62.2%	71.4%
	Pretty happy	32.9%	33.0%	35.4%	26.2%
	Not too happy	3.9%	2.1%	2.4%	2.4%
Female	Very happy	69.2%	63.9%	61.6%	57.6%
	Pretty happy	28.6%	33.2%	35.5%	33.3%
	Not too happy	2.2%	2.9%	2.8%	9.1%

Report created at 8:48:27 AM on 11/22/2004

Gender &gt; Happiness of marriage BY Number of children (grouped categories)

- The VLABELS subcommand suppresses the display of variable labels for *SEX* and *HAPMAR*.
- The SLABELS subcommand suppresses the default label for the summary function.
- The TITLE specification on the TITLE subcommand uses the standard IBM® SPSS® Statistics convention to break a single string across input lines.
- The CAPTION specification uses the ) DATE, ) TIME, and ) TABLE keywords to print the date, time, and a description of the table structure.

## Significance Testing

Custom Tables can perform the chi-square test of independence and pairwise comparisons of column proportions for tables that contain at least one category variable in both the rows and the columns. Custom Tables can perform pairwise comparisons of column means for tables that contain at least one summary variable in the rows and one category variable in the columns.

### Chi-Square Tests: SIGTEST Subcommand

```

/SIGTEST TYPE= CHISQUARE ALPHA= {0.05          }
                                {significance level}

INCLUDEMRSETS= {YES**}
                {NO   }

CATEGORIES= {ALLVISIBLE**}
             {SUBTOTALS  }

```

The SIGTEST subcommand has the following specifications:

<b>TYPE</b>	<i>Type of significance test.</i> The specification is required. The only current choice is CHISQUARE.
<b>ALPHA</b>	<i>Significance level for the test.</i> The specification must be greater than 0 and less than 1. The default is 0.05.
<b>INCLUDEMRSETS</b>	<i>Include multiple response variables in tests.</i> If there are no multiple response sets, this keyword is ignored. If INCLUDEMRSETS=YES and COUNTDUPLICATES=YES on the MRSETS subcommand, multiple response sets are suppressed with a warning.
<b>CATEGORIES</b>	<i>Replacing categories with subtotals for testing.</i> If SUBTOTALS is specified, each subtotal replaces its categories for significance testing. If ALLVISIBLE is specified, only subtotals that are specified by using the HSUBTOTAL keyword replace their categories for testing.

**Example**

```
CTABLES /TABLE AGECAT BY MARITAL
/CATEGORIES VARIABLES=AGECAT MARITAL TOTAL=YES
/SIGTEST TYPE=CHISQUARE.
```

Figure 50-24

		Marital status					
		Married	Widowed	Divorced	Separated	Never married	Total
		Count	Count	Count	Count	Count	Count
Age category	Less than 25	37	1	5	5	194	242
	25 to 34	271	13	63	16	263	626
	35 to 44	379	11	129	44	116	679
	45 to 54	275	18	123	13	52	481
	55 to 64	186	31	76	7	20	320
	65 or older	197	209	48	8	17	479
	Total	1345	283	444	93	662	2827

Figure 50-25

**Pearson Chi-Square Tests**

		Marital status
Age category	Chi-square	1473.381
	df	20
	Sig.	.000*

Results are based on nonempty rows and columns in each innermost subtable.

\*. The Chi-square statistic is significant at the 0.05 level.

**Pairwise Comparisons of Proportions and Means: COMPARETEST Subcommand**

```
/COMPARETEST TYPE= {PROP} ALPHA= {0.05}
                   {MEAN}          {significance level}

ADJUST= {BONFERRONI} ORIGIN=COLUMN
        {NONE}        }

INCLUDEMRSETS={YES**} MEANSVARIANCE={ALLCATS }
              {NO }                {TESTEDCATS}

CATEGORIES={ALLVISIBLE**}
           {SUBTOTALS }

MERGE={NO}
      {YES}
```

The COMPARETEST subcommand has the following specifications:

- TYPE** *The type of pairwise comparison.* The specification is required. To compare proportions when the test variable in the rows is categorical, choose PROP. The table must include counts or simple column percentages. To compare means when the test variable in the rows is scale and the column variable is categorical, choose MEAN. The table must include the mean as a summary statistic.
- ALPHA** *The significance level for the test.* The specification must be greater than 0 and less than 1. The default is 0.05.
- ADJUST** *The method for adjusting p values for multiple comparisons.* Valid options are NONE and BONFERRONI. If ADJUST is not specified, the Bonferroni correction is used.

<b>ORIGIN</b>	<i>The direction of the comparison.</i> This specification will determine whether column means (proportions) or row means (proportions) are being compared. Currently, only COLUMN is supported.
<b>INCLUDEMRSETS</b>	<i>Include multiple response variables in tests.</i> If there are no multiple response sets, this keyword is ignored. If INCLUDEMRSETS=YES and COUNTDUPLICATES=YES on the MRSETS subcommand, multiple response sets are suppressed with a warning.
<b>MEANSVARIANCE</b>	<i>Computation of variance for means test.</i> The variance for the means test is always based on the categories that are compared for multiple response tests, but for ordinary categorical variables, the variance can be estimated from just the categories that are compared or all categories. This keyword is ignored unless TYPE=MEAN.
<b>CATEGORIES</b>	<i>Replacing categories with subtotals for testing.</i> If SUBTOTALS is specified, each subtotal replaces its categories for significance testing. If ALLVISIBLE is specified, only subtotals that are specified by using the HSUBTOTAL keyword replace their categories for testing.
<b>MERGE</b>	<i>Merge significance indicators into the main table.</i> If NO is specified, significant differences are indicated in a separate table that follows the main table. If YES is specified, significance differences indicators are merged with the main table. The main table identifies significant differences with APA-style formatting using subscript letters. ( <i>Note:</i> The APA-style table includes a caption that explains how to interpret the subscripts in the table. If you specify a CAPTION on the TITLES subcommand, that caption will be displayed instead of the caption that explains the subscripts.)

**Example**

```
CTABLES /TABLE AGE CAT BY MARITAL
/CATEGORIES VARIABLES=AGE CAT MARITAL TOTAL=YES
/COMPARETEST TYPE=PROP ALPHA=.01.
```

Figure 50-26

Comparisons of Column Proportions<sup>a</sup>

		Marital status				
		Married	Widowed	Divorced	Separated	Never married
		(A)	(B)	(C)	(D)	(E)
Age category	Less than 25				B	A B C D
	25 to 34	B		B	B	A B C D
	35 to 44	B E		B E	A B C E	B
	45 to 54	B E		B E		
	55 to 64	E	E	E		
	65 or older	E	A C D E	E		

Results are based on two-sided tests with significance level .01. For each significant pair, the key of the category with the smaller column proportion appears under the category with the larger column proportion.

a. Tests are adjusted for all pairwise comparisons within a row of each innermost subtable using the Bonferroni correction.

- The table of counts is identical to that shown in the example for chi-square above.
- The comparison output shows a number of predictable pairs for marital status among different age groups that are significant at the 0.01 level that is specified with ALPHA in the command.

**Example**

```
CTABLES /TABLE AGE > SEX BY MARITAL
/CATEGORIES VARIABLES=SEX TOTAL=YES
```

```
/COMPARETEST TYPE=MEAN.
```

Figure 50-27

			Marital status				
			Married	Widowed	Divorced	Separated	Never married
			Mean	Mean	Mean	Mean	Mean
Age of respondent	Gender	Male	49	66	48	44	32
		Female	45	70	48	41	32
		Total	47	70	48	42	32

Figure 50-28

**Comparisons of Column Means<sup>a</sup>**

			Marital status				
			Married	Widowed	Divorced	Separated	Never married
			(A)	(B)	(C)	(D)	(E)
Age of respondent	Gender	Male	E	A C D E	E	E	
		Female	E	A C D E	D E	E	

Results are based on two-sided tests assuming equal variances with significance level 0.05. For each significant pair, the key of the smaller category appears under the category with larger mean.

<sup>a</sup>. Tests are adjusted for all pairwise comparisons within a row of each innermost subtable using the Bonferroni correction.

## FORMAT Subcommand

```
/FORMAT MINCOLWIDTH={DEFAULT} MAXCOLWIDTH={DEFAULT}
      {value } {value }
      UNITS={POINTS} EMPTY={ZERO } MISSING= {'.' }
      {INCHES} {BLANK } {'chars' }
      {CM } {'chars' }
```

The **FORMAT** subcommand controls the appearance of the table. At least one of the following attributes must be specified: **MINCOLWIDTH**, **MAXCOLWIDTH**, **UNITS**, **EMPTY**, or **MISSING**.

**MINCOLWIDTH** *The minimum width of columns in the table.* This setting includes the main tables as well as any tables of significance tests. **DEFAULT** honors the column labels setting in the current TableLook. The value must be less than or equal to the setting for **MAXCOLWIDTH**.

**MAXCOLWIDTH** *The maximum width of columns in the table.* This setting includes the main tables as well as any tables of significance tests. **DEFAULT** honors the column labels setting in the current TableLook. The value must be greater than or equal to the setting for **MINCOLWIDTH**.

**UNITS** *The measurement system for column width values.* The default is **POINTS**. You can also specify **INCHES** or **CM** (centimeters). **UNITS** is ignored unless **MINCOLWIDTH** or **MAXCOLWIDTH** is specified.

**EMPTY** *Fill characters used when a count or percentage is zero.* **ZERO** (the default) displays a 0 using the format for the cell statistic. **BLANK** leaves the statistic blank. You can also specify a quoted character string. If the string is too wide for the cell, the text is truncated.

**MISSING**

If `FORMAT EMPTY=BLANK`, there will be no visible difference between cells that have a count of 0 and cells for which no statistics are defined.

*Fill characters used when a cell statistic cannot be computed.* This specification applies to non-empty cells for which a statistic, such as standard deviation, cannot be computed. The default is a period (.). You can specify a quoted string. If the string is too wide for the cell, the text is truncated.

**VLABELS Subcommand**

```
/VLABELS VARIABLES=varlist
      DISPLAY={DEFAULT}
              {NAME   }
              {LABEL  }
              {BOTH   }
              {NONE   }
```

By default, the display of variable labels is controlled by the `TVARS` specification on the `SET` command in the Core system. The `VLABELS` subcommand allows you to show a name, label, or both for each table variable. The minimum specification is a variable list and a `DISPLAY` specification. To give different specifications for different variables, use multiple `VLABELS` subcommands.

- VARIABLES**      *The variables to which the subcommand applies.* You can use `ALL` or `VARNAME` TO `VARNAME`, which refers to the order of variables in the current active data file. If a specified variable does not appear in a table, `VLABELS` is ignored for that variable.
- DISPLAY**        *Whether the variable's name, label, both, or neither is shown in the table.* `DEFAULT` honors the `SET TVARS` setting. `NAME` shows the variable name only. `LABEL` shows the variable label only. `BOTH` shows the variable name and label. `NONE` hides the name and label.

**SMISSING Subcommand**

```
/SMISSING {VARIABLE}
          {LISTWISE}
```

If more than one scale variable is included in a table, you can control whether cases that are missing on one variable are included in summaries for which they have valid values.

- VARIABLE**      *Exclude cases variable by variable.* A case is included in summaries for each scale variable for which the case has a valid value regardless of whether the case has missing values for other scale variables in the table.
- LISTWISE**      *Exclude cases that are missing on any scale variable in the table.* This process ensures that summaries for all scale variables in the table are based on the same set of cases.

Listwise deletion applies on a per-table basis. Thus, given the specification

```
/TABLE (AGE [MEAN,COUNT]>SEX) + (AGE+CHILDS) [MEAN,COUNT] > HAPPY
```

all cases with valid values for `AGE` will be used in the `AGE > SEX` table, regardless of whether they have missing values for `CHILDS` (assuming that they also have valid values for `SEX`).

## MRSETS Subcommand

```
/MRSETS COUNTDUPLICATES= {NO }
                        {YES}
```

For multiple response sets that combine multiple category variables, a respondent can select the same response for more than one of the variables. Typically, only one response is desired. For example, \$MAGS can combine MAG1 to MAG5 to record which magazines a respondent reads regularly. If a respondent indicated the same magazine for MAG1 and MAG2, you would not want to count that magazine twice. However, if \$CARS combines CAR1 to CAR5 to indicate which cars a respondent owns now, and a respondent owns two cars of the same make, you might want to count both responses. The MRSETS subcommand allows you to specify whether duplicates are counted. By default, duplicates are not counted.

The MRSETS specification applies only to RESPONSES and percentages based on RESPONSES. MRSETS does not affect counts, which always ignore duplicates.

## PCOMPUTE Subcommand

In addition to displaying the aggregated results of summary statistics, a table can display one or more categories computed from those aggregated results and/or constant values. These are known as computed categories or postcomputes and are specified by the PCOMPUTE subcommand.

```
/PCOMPUTE &catname= EXPR(expression)
```

The left side of the equation specifies the computed category name, which must be a valid name preceded by an ampersand (&). If two or more PCOMPUTE subcommands specify the same category name, the last one takes precedence. Computed categories *cannot* be referenced in a table expression or in the expression for another computed category.

The right side of the equation defines the expression used to calculate the postcomputed category. The expression is calculated for each summary statistic and can consist of:

- **References to category values.** These are enclosed in brackets and must be quoted if they are strings or dates (for example, [1] or ['Male']). All categories in an expression must be of the same type (numeric, string, or date). Both valid and user-missing values are allowed.
- **Arithmetic operators.** The supported operators are - (subtraction), + (addition), \* (multiplication), / (division), and \*\* (exponentiation).
- **Parentheses.** These are used to control the order of evaluation.
- **Numeric constants.** Constants can be used by themselves (for example, &x=EXPR(50) or as part of a larger expression (for example, &x=EXPR(([1] + [2]) \* 0.8)).
- **Keywords.** The following keywords are supported in the expression: LO, THRU, HIGH, MISSING, and OTHERNM. If a range is specified, it must appear in square brackets (for example, &x=EXPR([LO THRU HIGH]) or &x=EXPR([1 THRU 5])). The values in the range are totaled.
- **References to totals and subtotals.** If only one SUBTOTAL or HSUBTOTAL is used in a category list, the PCOMPUTE expression may refer to it by the SUBTOTAL keyword only or by keyword and position. (Both SUBTOTAL and HSUBTOTAL are referenced by the SUBTOTAL keyword in the computed category's expression.) If multiple subtotals are used, the expression must refer



to them by position. The position is specified within square brackets following the keyword and must be an integer greater than or equal to 1 (for example, SUBTOTAL [2]). Totals are always referenced by keyword only.

All categories referenced in the expression must be also be explicitly specified in the CATEGORIES subcommand. As with other items in the table, the computed category's position in the CATEGORIES subcommand determines its position in the table.

By default, the values of computed categories use the same formatting for summary statistics as the other categories in the variable. The PPROPERTIES subcommand allows you to specify different formatting. This subcommand also allows you to specify a label for the computed category and hide the categories used in the computed category's expression. [For more information, see the topic PPROPERTIES Subcommand on p. 539.](#)

### Examples

```
CTABLES
  /PCOMPUTE &x=EXPR([2]-[1])
  /PPROPERTIES &x LABEL = "Female - Male"
  /TABLE SEX
  /CATEGORIES VARIABLES=SEX [1, 2, &x].
```

```
CTABLES
  /PCOMPUTE &x=EXPR([2]/[1])
  /PPROPERTIES &x LABEL = "Female / Male" FORMAT = COUNT F8.3
  /TABLE SEX
  /CATEGORIES VARIABLES=SEX [1, 2, &x].
```

```
CTABLES
  /PCOMPUTE &x = EXPR(SUBTOTAL[1] + SUBTOTAL[2])
  /PPROPERTIES &x LABEL="0-12"
  /TABLE educ[C]
  /CATEGORIES VARIABLES=educ [0 THRU 8, HSUBTOTAL="0-8", 9 THRU 12, HSUBTOTAL="9-12", &x,
  OTHERNM, HSUBTOTAL=">12"] TOTAL=YES.
```

## PPROPERTIES Subcommand

```
/PPROPERTIES &catname &catname...
  LABEL= "label"
  FORMAT= summary format ... summary format
  HIDESOURCECATS= {NO**}
                  {YES }
```

The PPROPERTIES subcommand defines properties for one or more computed categories. (For more information about creating computed categories, see the topic [PCOMPUTE Subcommand](#) on p. 538.) One PPROPERTIES subcommand can be used to define properties for multiple computed categories. If multiple PPROPERTIES subcommands are specified for a single computed category, the results are cumulative.

The PPROPERTIES subcommand has the following keywords:

<b>LABEL</b>	<i>Label for the computed category.</i> The quoted text defines the label that appears in the table. If no label is specified, the computed category's expression is used for the label. Like other labels, the maximum length is 120 bytes. The label can also include the symbol ) LABEL [n] to display the label for the <i>n</i> th category specified in the CATEGORIES subcommand.
<b>FORMAT</b>	<i>One or more pairs of summary statistics and display formats for the computed category.</i> Each summary statistic is followed by a display format for that statistic. For a list of summary statistics, see the topic <a href="#">Specifying Summaries</a> on p. 518. For a list of formats, see the topic <a href="#">Formats for Summaries</a> on p. 524.
<b>HIDESOURCECATS</b>	<i>Hide categories, totals, or subtotals used in the computed category's expression.</i> Use YES to hide the categories, totals, or subtotals used to evaluate the computed category. This setting ignores the EMPTY keyword. If HIDESOURCECAT=YES, categories are hidden, even if EMPTY is set to INCLUDE. You cannot use YES if you request significance tests.

## ***HIDESMALLCOUNTS Subcommand***

HIDESMALLCOUNTS allows you to hide counts displayed in tables for count values that are less than a specified integer. Hidden values are displayed as <N, where N is the specified integer.

### ***COUNT Keyword***

The COUNT keyword specifies the cutoff, below which count values are hidden. Values equal to the cutoff are not hidden. The value of COUNT must be an integer and greater than or equal to 2. The default is 5.

# CURVEFIT

CURVEFIT is available in the Statistics Base option

```
CURVEFIT VARIABLES= varname [WITH varname]
  [/MODEL= [LINEAR**] [LOGARITHMIC] [INVERSE]
           [QUADRATIC] [CUBIC] [COMPOUND]
           [POWER] [S] [GROWTH] [EXPONENTIAL]
           [LGSTIC] [ALL]]
  [/CIN={95** }
        {value}]
  [/UPPERBOUND={NO**}]
                {n }
  [/{CONSTANT† }
   {NOCONSTANT}]
  [/PLOT={FIT**}]
        {NONE }
  [/ID = varname]
  [/PRINT=ANOVA]
  [/SAVE=[PRED] [RESID] [CIN]]
  [/APPLY [= 'model name'] [/{SPECIFICATIONS}]
          {FIT }
  [/TEMPLATE= 'filename']
```

\*\*Default if the subcommand is omitted.

†Default if the subcommand is omitted and there is no corresponding specification on the TSET command.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 19.0

- TEMPLATE subcommand introduced.

## Example

```
CURVEFIT VARIABLES = VARY
  /MODEL=CUBIC.
```

## Overview

CURVEFIT fits selected curves to a line plot, allowing you to examine the relationship between one or more dependent variables and one independent variable. CURVEFIT also fits curves to time series and produces forecasts, forecast errors, lower confidence limits, and upper confidence limits. You can choose curves from a variety of regression models.

### Options

**Model Specification.** There are 11 regression models available on the MODEL subcommand. You can fit any or all of these to the data. The keyword ALL is available to fit all 11 models. You can control whether the regression equation includes a constant term using the CONSTANT or NOCONSTANT subcommand.

**Upperbound Value.** You can specify the upperbound value for the logistic model using the UPPERBOUND subcommand.

**Output.** You can produce an analysis-of-variance summary table using the PRINT subcommand. You can suppress the display of the curve-fitting plot using the PLOT subcommand.

**New Variables.** To evaluate the regression statistics without saving predicted and residual variables, specify TSET NEWVAR=NONE prior to CURVEFIT. To save the new variables and replace the variables saved earlier, use TSET NEWVAR=CURRENT (the default). To save the new variables without erasing variables saved earlier, use TSET NEWVAR=ALL or the SAVE subcommand on CURVEFIT.

**Forecasting.** When used with the PREDICT command, CURVEFIT can produce forecasts and confidence limits beyond the end of the series. [For more information, see the topic PREDICT on p. 1556.](#)

**Templates.** You can specify a template, using the TEMPLATE subcommand, to override the default chart attribute settings on your system.

### Basic Specification

The basic specification is one or more dependent variables. If the variables are not time series, you must also specify the keyword WITH and an independent variable.

- By default, the LINEAR model is fit.
- A 95% confidence interval is used unless it is changed by a TSET CIN command prior to the procedure.
- CURVEFIT produces a plot of the curve, a regression summary table displaying the type of curve used, the  $R^2$  coefficient, degrees of freedom, overall  $F$  test and significance level, and the regression coefficients.
- For each variable and model combination, CURVEFIT creates four variables: fit/forecast values, residuals, lower confidence limits, and upper confidence limits. These variables are automatically labeled and added to the active dataset unless TSET NEWVAR=NONE is specified prior to CURVEFIT. [For more information, see the topic SAVE Subcommand on p. 547.](#)

**Subcommand Order**

- Subcommands can be specified in any order.

**Syntax Rules**

- `VARIABLES` can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each one is executed.

**Operations**

- When `CURVEFIT` is used with the `PREDICT` command to forecast values beyond the end of a time series, the original and residual series are assigned the system-missing value after the last case in the original series.
- If a model requiring a log transformation (`COMPOUND`, `POWER`, `S`, `GROWTH`, `EXPONENTIAL`, or `LGSTIC`) is requested and there are values in the dependent variable(s) less than or equal to 0, the model cannot be fit because nonpositive values cannot be log-transformed.
- `CURVEFIT` uses listwise deletion of missing values. Whenever one dependent variable is missing a value for a particular case or observation, that case or observation will not be included in any computations.
- For the models `QUADRATIC` and `CUBIC`, a message is issued if the tolerance criterion is not met. (See `TSET` for information on changing the tolerance criterion.)
- Since `CURVEFIT` automatically generates four variables for each dependent variable and model combination, the `ALL` specification after `MODEL` should be used cautiously to avoid creating and adding to the active dataset many more variables than are necessary.
- The residual variable is always reported in the original metric. To compute the logged residual (which should be used for diagnostic checks) for the models `COMPOUND`, `POWER`, `S`, `GROWTH`, and `EXPONENTIAL`, specify

```
COMPUTE NEWVAR = LN(VAR) - LN(FIT#n) .
```

where *NEWVAR* is the logged residual, *VAR* is the name of the dependent variable or observed series, and *FIT#n* is the name of the fitted variable generated by `CURVEFIT`.

For the `LGSTIC` (logistic) model, the logged residual can be obtained by

```
COMPUTE NEWERR = LN(VAR) - LN(1/FIT#n) .
```

or, if upperbound value *u* is specified on the `UPPERBOUND` subcommand, by

```
COMPUTE NEWVAR = LN(1/VAR - 1/u) - LN(1/FIT#n) .
```

- `CURVEFIT` obeys the `WEIGHT` command when there is an independent variable. The `WEIGHT` specification is ignored if no independent variable is specified.

**Limitations**

- A maximum of 1 `VARIABLES` subcommand. There is no limit on the number of dependent variables or series named on the subcommand.
- A maximum of 1 independent variable can be specified after the keyword `WITH`.

## Examples

### **Comparing linear and quadratic models**

```
* Curve Estimation.
CURVEFIT /VARIABLES=sales WITH advert
/CONSTANT
/MODEL=LINEAR QUADRATIC
/PLOT FIT
/SAVE=PRED RESID .
```

- The procedure fits a curve to *sales* using *advert*.
- The `MODEL` subcommand requests linear and quadratic models; the `CONSTANT` subcommand specifies that a constant term should be included in each model.
- The `PLOT` subcommand requests a plot of *sales* by *advert* for the observed values and each fit model.
- The `SAVE` subcommand requests that fitted values and residuals be saved to the active dataset.

### **A cubic model**

```
* Curve Estimation.
CURVEFIT /VARIABLES=infected WITH time
/CONSTANT
/MODEL= CUBIC
/PRINT ANOVA
/PLOT FIT
/SAVE=PRED RESID .
```

- The procedure fits a curve to *infected* using *time*.
- The `MODEL` subcommand requests a cubic model; the `CONSTANT` subcommand specifies that a constant term should be included in the model.
- The `PRINT` subcommand requests an ANOVA table.
- The `PLOT` subcommand requests a plot of *infected* by *time* for the observed values and the fit model.
- The `SAVE` subcommand requests that fitted values and residuals be saved to the active dataset.

## **VARIABLES Subcommand**

`VARIABLES` specifies the variables and is the only required subcommand.

- If the dependent variables specified are not time series, you must also specify the keyword `WITH` and an independent variable.

## **MODEL Subcommand**

`MODEL` specifies the model or models to be fit to the data. The default model is `LINEAR`.

- You can fit any or all of the 11 available models.
- Model name keywords can be abbreviated to the first three characters.

- You can use the keyword `ALL` to fit all models.
- When the `LGSTIC` model is specified, the upperbound value is included in the output.

The following table lists the available models and their regression equations. The linear transformations for the last six models are also shown.

<b>Keyword</b>	<b>Equation</b>	<b>Linear equation</b>
LINEAR	$Y = b_0 + b_1 t$	
LOGARITHMIC	$Y = b_0 + b_1 \ln(t)$	
INVERSE	$Y = b_0 + b_1/t$	
QUADRATIC	$Y = b_0 + b_1 t + b_2 t^2$	
CUBIC	$Y = b_0 + b_1 t + b_2 t^2 + b_3 t^3$	
COMPOUND	$Y = b_0 b_1^t$	$\ln(Y) = \ln(b_0) + t \ln(b_1)$
POWER	$Y = b_0 t^{b_1}$	$\ln(Y) = \ln(b_0) + b_1 \ln(t)$
S	$Y = e^{b_0 + b_1/t}$	$\ln(Y) = b_0 + b_1/t$
GROWTH	$Y = e^{b_0 + b_1 t}$	$\ln(Y) = b_0 + b_1 t$
EXPONENTIAL	$Y = b_0 e^{b_1 t}$	$\ln(Y) = \ln(b_0) + b_1 t$
LGSTIC (logistic)	$Y = (1/u + b_0 b_1^t)_{-1}$	$\ln(1/Y - 1/u) = \ln(b_0) + t \ln(b_1)$

where

$b_0$  = a constant

$b_n$  = regression coefficient

$t$  = independent variable or time value

$\ln$  = the natural logarithm

$u$  = upperbound value for `LGSTIC`

### **Example**

```
CURVEFIT VARIABLES = VARX.
```

- This command fits a curve to `VARX` using the linear regression model (the default).

### **Example**

```
CURVEFIT VARIABLES = VARY
/MODEL=GROWTH EXPONENTIAL.
```

- This command fits two curves to `VARY`, one using the growth model and the other using the exponential model.

## **UPPERBOUND Subcommand**

UPPERBOUND is used with the logistic model (keyword LGSTIC) to specify an upper boundary value to be used in the regression equation.

- The specification on UPPERBOUND must be a positive number and must be greater than the largest data value in any of the specified dependent variables.
- The default UPPERBOUND value is infinity, so that  $1/u = 0$  and is dropped from the equation.
- You can specify UPPERBOUND NO to reset the value to infinity when applying a previous model.
- If you specify UPPERBOUND without LGSTIC, it is ignored.
- Note that UPPERBOUND is a subcommand and cannot be used within a MODEL subcommand. For example, the following specification is *not* valid:

```
/MODEL=CUBIC LGSTIC /UPPER=99 LINEAR
```

The correct specification is:

```
/MODEL=CUBIC LGSTIC LINEAR
/UPPER=99
```

## **CONSTANT and NOCONSTANT Subcommands**

CONSTANT and NOCONSTANT indicate whether a constant term should be estimated in the regression equation. The specification overrides the corresponding setting on the TSET command.

- CONSTANT indicates that a constant should be estimated. It is the default unless changed by TSET NOCONSTANT prior to the current procedure.
- NOCONSTANT eliminates the constant term from the model.

### **Example**

```
CURVEFIT VARIABLES = Y1
/MODEL=COMPOUND
/NOCONSTANT.
```

- In this example, a compound curve is fit to *Y1* with no constant term in the model.

## **CIN Subcommand**

CIN controls the size of the confidence interval for confidence interval variables saved with the SAVE subcommand. It has no effect on the model results displayed in output tables.

- The specification on CIN must be greater than 0 and less than 100.
- The default confidence interval is 95.
- The CIN subcommand overrides the TSET CIN setting.



## ***PLOT Subcommand***

PLOT specifies whether the curve-fitting plot is displayed. If PLOT is not specified, the default is FIT. The curve-fitting plot is displayed. PLOT=FIT is generally used with an APPLY subcommand to turn off a PLOT=NONE specification in the applied model.

<b>FIT</b>	<i>Display the curve-fitting plot.</i>
<b>NONE</b>	<i>Do not display the plot.</i>

## ***ID Subcommand***

ID specifies an identification variable. When in point selection mode, you can click on an individual chart point to display the value of the ID variable for the selected case.

## ***SAVE Subcommand***

SAVE saves the values of predicted, residual, and/or confidence interval variables generated during the current session in the active dataset.

- SAVE saves the specified variables with default names: *FIT<sub>n</sub>* for predicted values, *ERR<sub>n</sub>* for residuals, *LCL<sub>n</sub>* for the lower confidence limit, and *UCL<sub>n</sub>* for the upper confidence limit, where *n* increments each time any variable is saved for a model.
- SAVE overrides the CURRENT or NONE setting on TSET NEWVARS (see TSET).

<b>PRED</b>	<i>Predicted variable.</i>
<b>RESID</b>	<i>Residual variable.</i>
<b>CIN</b>	<i>Confidence interval.</i> The default is a 95% confidence interval. Use the CIN subcommand to specify a different confidence interval.

## ***PRINT Subcommand***

PRINT is used to produce an additional analysis-of-variance table for each model and variable.

- The only specification on PRINT is the keyword ANOVA.

## ***APPLY Subcommand***

APPLY allows you to use a previously defined CURVEFIT model without having to repeat the specifications.

- The specifications on APPLY can include the name of a previous model in quotes and one of two keywords. All of these specifications are optional.
- If a model name is not specified, the model specified on the previous CURVEFIT command is used.
- To change one or more of the specifications of the model, specify the subcommands of only those portions you want to change after the subcommand APPLY.

- If no variables or series are specified on the CURVEFIT command, the dependent variables that were originally specified with the model being reapplied are used.
- To change the dependent variables used with the model, enter new variable names before or after the APPLY subcommand.

The keywords available for APPLY on CURVEFIT are:

<b>SPECIFICATIONS</b>	<i>Use only the specifications from the original model. This is the default.</i>
<b>FIT</b>	<i>Use the coefficients estimated for the original model in the equation.</i>

### Example

```
CURVEFIT VARIABLES = X1
/MODEL=QUADRATIC.
CURVEFIT VARIABLES = Z1
/APPLY.
```

- The first command fits a quadratic curve to *X1*.
- The second command fits the same type of curve to *Z1*.

### Example

```
CURVEFIT VARIABLES = X1 Y1 Z1
/MODEL=QUADRATIC.
CURVEFIT APPLY
/MODEL=CUBIC.
```

- The first command fits quadratic curves to *X1*, *Y1*, and *Z1*.
- The second command fits curves to the same three series using the cubic model.

## TEMPLATE Subcommand

TEMPLATE uses an existing file as a template and applies it to any charts generated by the current CURVEFIT command. The template overrides the default settings that are used to create any charts, and the specifications on the current CURVEFIT command override the template. Templates are created in the Chart Editor by saving an existing chart as a template.

### Example

```
CURVEFIT VARIABLES = vary
/MODEL=CUBIC
/TEMPLATE=' /templates/mytemplate.sgt'.
```

## References

- Abraham, B., and J. Ledolter. 1983. *Statistical methods of forecasting*. New York: John Wiley and Sons.
- Draper, N. R., and H. Smith. 1981. *Applied regression analysis*, 2nd ed. New York: John Wiley and Sons.

Montgomery, D. C., and E. A. Peck. 1982. *Introduction to linear regression analysis*. New York: John Wiley and Sons.

# DATA LIST

```
DATA LIST [FILE='file'] [ENCODING='encoding specification']
[{FIXED}]
  {FREE } [{"delimiter", "delimiter",..., TAB}]
  {LIST }

[RECORDS={1}] [SKIP={n}] [{TABLE }]
  {n}                {NOTABLE}

/{1 } varname {col location [(format)]} [varname ...]
  {rec #}                {(FORTRAN-like format) }

[/ {2 } ...] [/ ...]
  {rec #}
```

## Numeric and string input formats:

Type	Column-style format	FORTRAN-like format
Numeric (default)	d or F, d	Fw.d
Restricted numeric	N, d	Nw.d
Scientific notation	E, d	Ew.d
Numeric with commas	COMMA, d	COMMAw.d
Numeric with dots	DOT, d	DOTw.d
Numeric with commas and dollar sign	DOLLAR, d	DOLLARw.d
Numeric with percent sign	PCT, d	PCTw.d
Zoned decimal	Z, d	Zw.d
String	A	Aw

## Format elements to skip columns:

Type	Column-style format	FORTRAN-like format
Tab to column <i>n</i>		Tn
Skip <i>n</i> columns		nX

## Date and time input formats:

Type	Data input	Format	FORTRAN-like format
International date	dd-mmm-yyyy	DATE	DATEw
American date	mm/dd/yyyy	ADATE	ADATEw
European date	dd/mm/yy	EDATE	EDATEw
Julian date	yyddd	JDATE	JDATEw
Sorted date	yy/mm/dd	SDATE	SDATEw
Quarter and year	qQyyyy	QYR	QYRw
Month and year	mm/yyyy	MOYR	MOYRw
Week and year	wkWKyyyy	WKYR	WKYRw
Date and time	dd-mmm-yyyy hh:mm:ss.ss	DATETIME	DATETIMEw.d
Time	hh:mm:ss.ss	TIME	TIMEw.d

Type	Data input	Format	FORTRAN-like format
Days and time	ddd hh:mm:ss.ss	DTIME	DTIMEw . d
Day of the week	string	WKDAY	WKDAYw
Month	string	MONTH	MONTHw

*Note:* For default numeric (F) format and scientific notation (E) format, the decimal indicator of the input data must match the IBM® SPSS® Statistics locale decimal indicator (period or comma). Use `SHOW DECIMAL` to display the current decimal indicator and `SET DECIMAL` to set the decimal indicator. (Comma and Dollar formats only recognize a period as the decimal indicator, and Dot format only recognizes the comma as the decimal indicator.)

### Release History

Release 16.0

- `ENCODING` subcommand added for Unicode support.

### Example

```
DATA LIST /ID 1-3 SEX 5 (A) AGE 7-8 OPINION1 TO OPINION5 10-14.
```

## Overview

`DATA LIST` defines a raw data file (a raw data file contains numbers and other alphanumeric characters) by assigning names and formats to each variable in the file. Raw data can be inline (entered with your commands between `BEGIN DATA` and `END DATA`) or stored in an external file. They can be in fixed format (values for the same variable are always entered in the same location on the same record for each case) or in freefield format (values for consecutive variables are not in particular columns but are entered one after the other, separated by blanks or commas).

For information on defining matrix materials, see `MATRIX DATA`. For information on defining complex data files that cannot be defined with `DATA LIST`, see `FILE TYPE` and `REPEATING DATA`. For information on reading IBM® SPSS® Statistics data files and portable files, see `GET` and `IMPORT`.

The program can also read data files created by other software applications. Commands that read these files include `GET CAPTURE` and `GET TRANSLATE`.

### Options

**Data Source.** You can use inline data or data from an external file.

**Data Formats.** You can define numeric (with or without decimal places) and string variables using an array of input formats (percent, dollar, date and time, and so forth). You can also specify column binary and unaligned positive integer binary formats (available only if used with the `MODE=MULTIPUNCH` setting on the `FILE HANDLE` command).

**Data Organization.** You can define data that are in fixed format (values in the same location on the same record for each case), in freefield format with multiple cases per record, or in freefield format with one case on each record using the `FIXED`, `FREE`, and `LIST` keywords.

**Multiple Records.** For fixed-format data, you can indicate the number of records per case on the RECORDS subcommand. You can specify which records to read in the variable definition portion of DATA LIST.

**Summary Table.** For fixed-format data, you can display a table that summarizes the variable definitions using the TABLE subcommand. You can suppress this table using NOTABLE.

**Value Delimiter.** For freefield-format data (keywords FREE and LIST), you can specify the character(s) that separate data values, or you can use the keyword TAB to specify the tab character as the delimiter. Any delimiter other than the TAB keyword must be enclosed in quotation marks, and the specification must be enclosed in parentheses, as in DATA LIST FREE(" , ").

**End-of-File Processing.** You can specify a logical variable that indicates the end of the data using the END subcommand. This logical variable can be used to invoke special processing after all the cases from the data file have been read.

### **Basic Specification**

- The basic specification is the FIXED, LIST, or FREE keyword followed by a slash that signals the beginning of variable definition.
- FIXED is the default.
- If the data are in an external file, the FILE subcommand must be used.
- If the data are inline, the FILE subcommand is omitted and the data are specified between the BEGIN DATA and END DATA commands.
- Variable definition for fixed-format data includes a variable name, a column location, and a format (unless the default numeric format is used). The column location is not specified if FORTRAN-like formats are used, since these formats include the variable width.
- Variable definition for freefield data includes a variable name and, optionally, a delimiter specification and a FORTRAN-like format specification. If format specifications include a width and number of decimal positions (for example, F8.2), the width and decimal specifications are not used to read the data but are assigned as print and write formats for the variables.

### **Subcommand Order**

Subcommands can be named in any order. However, all subcommands must precede the first slash, which signals the beginning of variable definition.

### **Syntax Rules**

Subcommands on DATA LIST are separated by spaces or commas, not by slashes.

## **Examples**

\* Column-style format specifications.

```
DATA LIST /ID 1-3 SEX 5 (A) AGE 7-8 OPINION1 TO OPINION5 10-14.  
BEGIN DATA  
001 m 28 12212  
002 f 29 21212
```

```
003 f 45 32145
...
128 m 17 11194
END DATA.
```

- The data are inline between the `BEGIN DATA` and `END DATA` commands, so the `FILE` subcommand is not specified. The data are in fixed format. The keyword `FIXED` is not specified because it is the default.
- Variable definition begins after the slash. Variable *ID* is in columns 1 through 3. Because no format is specified, numeric format is assumed. Variable *ID* is therefore a numeric variable that is three digits wide.
- Variable *SEX* is a short string variable in column 5. Variable *SEX* is one byte wide.
- *AGE* is a two-column numeric variable in columns 7 and 8.
- Variables *OPINION1*, *OPINION2*, *OPINION3*, *OPINION4*, and *OPINION5* are named using the `TO` keyword. Each is a one-column numeric variable, with *OPINION1* located in column 10 and *OPINION5* located in column 14.
- The `BEGIN DATA` and `END DATA` commands enclose the inline data. Note that the values of *SEX* are lowercase letters and must be specified as such on subsequent commands.

## Operations

- `DATA LIST` creates a new active dataset.
- Variable names are stored in the active dataset dictionary.
- Formats are stored in the active dataset dictionary and are used to display and write the values. To change output formats of numeric variables defined on `DATA LIST`, use the `FORMATS` command.
- For default numeric (F) format and scientific notation (E) format, the decimal indicator of the input data must match the IBM® SPSS® Statistics locale decimal indicator (period or comma). Use `SHOW DECIMAL` to display the current decimal indicator and `SET DECIMAL` to set the decimal indicator. (Comma and Dollar formats only recognize a period as the decimal indicator, and Dot format only recognizes the comma as the decimal indicator.)

## Fixed-Format Data

- The order of the variables in the active dataset dictionary is the order in which they are defined on `DATA LIST`, not their sequence in the input data file. This order is important if you later use the `TO` keyword to refer to variables on subsequent commands.
- In numeric format, blanks to the left or right of a number are ignored; embedded blanks are invalid. When the program encounters a field that contains one or more blanks interspersed among the numbers, it issues a warning message and assigns the system-missing value to that case.
- Alphabetical and special characters, except the decimal point and leading plus and minus signs, are not valid in numeric variables and are set to system-missing if encountered in the data.
- See [String Formats](#) for information on column width specifications for string variables in UTF-8 data files.

- The system-missing value is assigned to a completely blank field for numeric variables. The value assigned to blanks can be changed using the `BLANKS` specification on the `SET` command.
- The program ignores data contained in columns and records that are not specified in the variable definition.

## ***Freefield Data***

`FREE` can read freefield data with multiple cases recorded on one record or with one case recorded on more than one record. `LIST` can read freefield data with one case on each record.

- Line endings are read as delimiters between values.
- If you use FORTRAN-like format specifications (for example, `DOLLAR12.2`), width and decimal specifications are not used to read the data but are assigned as print and write formats for the variable.

For freefield data *without* explicitly specified value delimiters:

- Commas and blanks are interpreted as delimiters between values.
- Extra blanks are ignored.
- Multiple commas with or without blank space between them can be used to specify missing data.
- If a valid value contains commas or blank spaces, enclose the values in quotes.

For data with explicitly specified value delimiters (for example, `DATA LIST FREE (",")`):

- Multiple delimiters without any intervening space can be used to specify missing data.
- The specified delimiters cannot occur within a data value, even if you enclose the value in quotes.

*Note:* Freefield format with specified value delimiters is typically used to read data in text format written by a computer program, not for data manually entered in a text editor.

## ***FILE Subcommand***

`FILE` specifies the raw data file. `FILE` is required when data are stored in an external data file. `FILE` must not be used when the data are stored in a file that is included with the `INCLUDE` command or when the data are inline (see `INCLUDE` and `BEGIN DATA-END DATA`).

- `FILE` must be separated from other `DATA LIST` subcommands by at least one blank or comma.
- `FILE` must precede the first slash, which signals the beginning of variable definition.

## ***ENCODING Subcommand***

`ENCODING` specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is UTF8. For more information, see [SET command](#), [UNICODE subcommand](#).



- In code page mode, the default is the current locale setting. For more information, see [SET command](#), [LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), `UTF8`, `UTF16`, `UTF16BE` (big endian), `UTF16LE` (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- In Unicode mode, the defined width of string variables is tripled for code page and UTF-16 text data files. Use [ALTER TYPE](#) to automatically adjust the defined width of string variables.
- See [String Formats](#) for information on column width specifications for string variables in UTF-8 data files.
- If there is no `FILE` subcommand, the `ENCODING` subcommand is ignored.

## ***FIXED, FREE, and LIST Keywords***

`FIXED`, `FREE`, or `LIST` indicates the format of the data. Only one of these keywords can be used on each `DATA LIST`. The default is `FIXED`.

<b>FIXED</b>	<i>Fixed-format data.</i> Each variable is recorded in the same column location on the same record for each case in the data. <code>FIXED</code> is the default.
<b>FREE</b>	<i>Freefield data.</i> The variables are recorded in the same order for each case but not necessarily in the same column locations. More than one case can be entered on the same record. By default, values are separated by blanks or commas. You can also specify different value delimiters.
<b>LIST</b>	<i>Freefield data with one case on each record.</i> The variables are recorded in freefield format as described for the keyword <code>FREE</code> except that the variables for each case must be recorded on one record.

- `FIXED`, `FREE`, or `LIST` must be separated from other `DATA LIST` subcommands by at least one blank or comma.
- `FIXED`, `FREE`, or `LIST` must precede the first slash, which signals the beginning of data definition.
- For fixed-format data, you can use column-style or FORTRAN-like formats, or a combination of both. For freefield data, you can use only FORTRAN-like formats.
- For fixed-format data, the program reads values according to the column locations specified or implied by the FORTRAN-like format. Values in the data do *not* have to be in the same order as the variables named on `DATA LIST` and do *not* have to be separated by a space or column.
- See [String Formats](#) for information on column width specifications for string variables in UTF-8 format data files.
- For freefield data, the program reads values sequentially in the order in which the variables are named on `DATA LIST`. Values in the data *must* be in the order in which the variables are named on `DATA LIST` and *must* be separated by at least one valid delimiter.
- For freefield data, multiple blank spaces can be used to indicate missing information only if a blank space is explicitly specified as the delimiter. In general, it is better to use multiple nonblank delimiters (for example, two commas with no intervening space) to specify missing data.
- In freefield format, a value cannot be split across records.

**Example**

\* Data in fixed format.

```
DATA LIST FILE="/data/hubdata.txt" FIXED RECORDS=3
  /1 YRHIRED 14-15 DEPT 19 SEX 20.
```

- FIXED indicates explicitly that the *hubdata.txt* file is in fixed format. Because FIXED is the default, the keyword FIXED could have been omitted.
- Variable definition begins after the slash. Column locations are specified after each variable. Since formats are not specified, the default numeric format is used. Variable widths are determined by the column specifications: *YRHIRED* is two digits wide, and *DEPT* and *SEX* are each one digit wide.

**Example**

\* Data in freefield format.

```
DATA LIST FREE / POSTPOS NWINS.
BEGIN DATA
2, 19, 7, 5, 10, 25, 5, 17, 8, 11, 3,, 6, 8, 1, 29
END DATA.
```

- Data are inline, so FILE is omitted. The keyword FREE is used because data are in freefield format with multiple cases on a single record. Two variables, *POSTPOS* and *NWINS*, are defined. Since formats are not specified, both variables receive the default F8.2 format.
- All of the data are recorded on one record. The first two values build the first case in the active dataset. For the first case, *POSTPOS* has value 2 and *NWINS* has value 19. For the second case, *POSTPOS* has value 7 and *NWINS* has value 5, and so on. The active dataset will contain eight cases.
- The two commas without intervening space after the data value 3 indicate a missing data value.

**Example**

\* Data in list format.

```
DATA LIST LIST (" ")/ POSTPOS NWINS.
BEGIN DATA
2,19
7,5
10,25
5,17
8,11
3,
6,8
1,29
END DATA.
```

- This example defines the same data as the previous example, but LIST is used because each case is recorded on a separate record. FREE could also be used. However, LIST is less prone to errors in data entry. If you leave out a value in the data with FREE format, all values after the missing value are assigned to the wrong variable. Since LIST format reads a case from each record, a missing value will affect only one case.

- A comma is specified as the delimiter between values.
- Since line endings are interpreted as delimiters between values, the second comma after the value 3 (in the sixth line of data) is not necessary to indicate that the value of *NWINS* is missing for that case.

## **TABLE and NOTABLE Subcommands**

TABLE displays a table summarizing the variable definitions supplied on DATA LIST. NOTABLE suppresses the summary table. TABLE is the default.

- TABLE and NOTABLE can be used only for fixed-format data.
- TABLE and NOTABLE must be separated from other DATA LIST subcommands by at least one blank or comma.
- TABLE and NOTABLE must precede the first slash, which signals the beginning of variable definition.

## **RECORDS Subcommand**

RECORDS indicates the number of records per case for fixed-format data. In the variable definition portion of DATA LIST, each record is preceded by a slash. By default, DATA LIST reads one record per case.

- The only specification on RECORDS is a single integer indicating the *total* number of records for each case (even if the DATA LIST command does not define all the records).
- RECORDS can be used only for fixed-format data and must be separated from other DATA LIST subcommands by at least one blank or comma. RECORDS must precede the first slash, which signals the beginning of variable definition.
- Each slash in the variable definition portion of DATA LIST indicates the beginning of a new record. The first slash indicates the first (or only) record. The second and any subsequent slashes tell the program to go to a new record.
- To skip a record, specify a slash without any variables for that record.
- The number of slashes in the variable definition cannot exceed the value of the integer specified on RECORDS.
- The sequence number of the record being defined can be specified after each slash. DATA LIST reads the number to determine which record to read. If the sequence number is used, you *do not* have to use a slash for any skipped records. However, the records to be read must be in their sequential order.
- The slashes for the second and subsequent records can be specified within the variable list, or they can be specified on a format list following the variable list (see the example below).
- All variables to be read from one record should be defined before you proceed to the next record.
- Since RECORDS can be used only with fixed format, it is not necessary to define all the variables on a given record or to follow their order in the input data file.

**Example**

```
DATA LIST FILE="/data/hubdata.txt" RECORDS=3
  /2 YRHIRED 14-15 DEPT 19 SEX 20.
```

- DATA LIST defines fixed-format data. RECORDS can be used only for fixed-format data.
- RECORDS indicates that there are three records per case in the data. Only one record per case is defined in the data definition.
- The sequence number (2) before the first variable definition indicates that the variables being defined are on the second record. Because the sequence number is provided, a slash is not required for the first record, which is skipped.
- The variables *YRHIRED*, *DEPT*, and *SEX* are defined and will be included in the active dataset. Any other variables on the second record or on the other records are not defined and are not included in the active dataset.

**Example**

```
DATA LIST FILE="/data/hubdata.txt" RECORDS=3
  / /YRHIRED 14-15 DEPT 19 SEX 20.
```

- This command is equivalent to the one in the previous example. Because the record sequence number is omitted, a slash is required to skip the first record.

**Example**

```
DATA LIST FILE="/data/hubdata.txt" RECORDS=3
  /YRHIRED (T14,F2.0) / /NAME (T25,A24).
```

- RECORDS indicates there are three records for each case in the data.
- *YRHIRED* is the only variable defined on the first record. The FORTRAN-like format specification T14 means tab over 14 columns. Thus, *YRHIRED* begins in column 14 and has format F2.0.
- The second record is skipped. Because the record sequence numbers are not specified, a slash must be used to skip the second record.
- *NAME* is the only variable defined for the third record. *NAME* begins in column 25 and is a string variable with a width of 24 bytes (format A24).

**Example**

```
DATA LIST FILE="/data/hubdata.txt" RECORDS=3
  /YRHIRED NAME (T14,F2.0 / / T25,A24).
```

- This command is equivalent to the one in the previous example. *YRHIRED* is located on the first record, and *NAME* is located on the third record.
- The slashes that indicate the second and third records are specified within the format specifications. The format specifications follow the complete variable list.

## SKIP Subcommand

SKIP skips the first  $n$  records of the data file.

### Example

```
DATA LIST LIST SKIP=2 /numvar.
BEGIN DATA
Some text describing the file
followed by some more text
1
2
3
END DATA.
```

## END Subcommand

END provides control of end-of-file processing by specifying a variable that is set to a value of 0 until the end of the data file is encountered, at which point the variable is set to 1. The values of all variables named on DATA LIST are left unchanged. The logical variable created with END can then be used on DO IF and LOOP commands to invoke special processing after all of the cases from a particular input file have been built.

- DATA LIST and the entire set of commands used to define the cases must be enclosed within an INPUT PROGRAM–END INPUT PROGRAM structure. The END FILE command must also be used to signal the end of case generation.
- END can be used only with fixed-format data. An error is generated if the END subcommand is used with FREE or LIST.

### Example

```
INPUT PROGRAM.
NUMERIC      TINCOME (DOLLAR8.0).          /* Total income
LEAVE        TINCOME.
DO IF        $CASENUM EQ 1.
+ PRINT      EJECT.
+ PRINT      / 'Name          Income'.
END IF
DATA LIST    FILE=INCOME END=#EOF NOTABLE / NAME 1-10(A)
                                                    INCOME 16-20(F).

DO IF        #EOF.
+ PRINT      / 'TOTAL          ', TINCOME.
+ END FILE.
ELSE.
+ PRINT      / NAME, INCOME (A10,COMMA8).
+ COMPUTE    TINCOME = TINCOME+INCOME. /* Accumulate total income
END IF.
END INPUT PROGRAM.

EXECUTE.
```

- The data definition commands are enclosed within an INPUT PROGRAM–END INPUT PROGRAM structure.
- NUMERIC indicates that a new numeric variable, *TINCOME*, will be created.

- LEAVE tells the program to leave variable *TINCOME* at its value for the previous case as each new case is read, so that it can be used to accumulate totals across cases.
- The first DO IF structure, enclosing the PRINT EJECT and PRINT commands, tells the program to display the headings *Name* and *Income* at the top of the display (when *\$CASENUM* equals 1).
- DATA LIST defines variables *NAME* and *INCOME*, and it specifies the scratch variable *#EOF* on the END subcommand.
- The second DO IF prints the values for *NAME* and *INCOME* and accumulates the variable *INCOME* into *TINCOME* by passing control to ELSE as long as *#EOF* is not equal to 1. At the end of the file, *#EOF* equals 1, and the expression on DO IF is true. The label *TOTAL* and the value for *TINCOME* are displayed, and control is passed to END FILE.

### Example

```
* Concatenate three raw data files.

INPUT PROGRAM.
NUMERIC #EOF1 TO #EOF3. /*These will be used as the END variables.

DO IF #EOF1 & #EOF2 & #EOF3.
+   END FILE.
ELSE IF #EOF1 & #EOF2.
+   DATA LIST FILE=THREE END=#EOF3 NOTABLE / NAME 1-20(A)
+       AGE 25-26 SEX 29(A).
+   DO IF NOT #EOF3.
+       END CASE.
+   END IF.
ELSE IF #EOF1.
+   DATA LIST FILE=TWO END=#EOF2 NOTABLE / NAME 1-20(A)
+       AGE 21-22 SEX 24(A).
+   DO IF NOT #EOF2.
+       END CASE.
+   END IF.
ELSE.
+   DATA LIST FILE=ONE END=#EOF1 NOTABLE /1 NAME 1-20(A)
+       AGE 21-22 SEX 24 (A).
+   DO IF NOT #EOF1.
+       END CASE.
+   END IF.
END IF.
END INPUT PROGRAM.

REPORT FORMAT AUTOMATIC LIST /VARS=NAME AGE SEX.
```

- The input program contains a DO IF–ELSE IF–END IF structure.
- Scratch variables are used on each END subcommand so the value will not be reinitialized to the system-missing value after each case is built.
- Three data files are read, two of which contain data in the same format. The third requires a slightly different format for the data items. All three DATA LIST commands are placed within the DO IF structure.
- END CASE builds cases from each record of the three files. END FILE is used to trigger end-of-file processing once all data records have been read.

## Variable Definition

The variable definition portion of `DATA LIST` assigns names and formats to the variables in the data. Depending on the format of the file, you may also need to specify record and column location. The following sections describe variable names, location, and formats.

### Variable Names

- Variable names must conform to variable-naming rules. System variables (beginning with a `$`) cannot be defined on `DATA LIST`. For more information, see the topic [Variable Names on p. 49](#).
- The keyword `TO` can be used to generate names for consecutive variables in the data. Leading zeros in the number are preserved in the name. `X1 TO X100` and `X001 TO X100` both generate 100 variable names, but the first 99 names are not the same in the two lists. `X01 TO X9` is not a valid specification.
- The order in which variables are named on `DATA LIST` determines their order in the active dataset.

#### Example

```
DATA LIST FREE / ID SALARY #V1 TO #V4.
```

- The `FREE` keyword indicates that the data are in freefield format. Six variables are defined: `ID`, `SALARY`, `#V1`, `#V2`, `#V3`, and `#V4`. `#V1` to `#V4` are scratch variables that are not stored in the active dataset. Their values can be used in transformations but not in procedure commands.

### Variable Location

For fixed-format data, variable locations are specified either explicitly using column locations or implicitly using FORTRAN-like formats. For freefield data, variable locations are not specified. Values are read sequentially in the order in which variables are named on the variable list.

#### Fixed-Format Data

- If column-style formats are used, you must specify the column location of each variable after the variable name. If the variable is one column wide, specify the column number. Otherwise, specify the first column number followed by a dash (–) and the last column number.
- If several adjacent variables on the same record have the same width and format type, you can use one column specification after the last variable name. Specify the beginning column location of the first variable, a dash, and the ending column location of the last variable. The program divides the total number of columns specified equally among the variables. If the number of columns does not divide equally, an error message is issued.
- The same column locations can be used to define multiple variables.
- For FORTRAN-like formats, column locations are implied by the width specified on the formats. For more information, see the topic [Variable Formats on p. 563](#). To skip columns, use the `Tn` or `nX` format specifications.

- With fixed format, column-style and FORTRAN-like specifications can be mixed on the same `DATA LIST` command.
- Record location is indicated by a slash or a slash and record number before the names of the variables on that record. [For more information, see the topic RECORDS Subcommand on p. 557.](#)
- The program ignores data in columns and on records that are not specified on `DATA LIST`.
- In the data, values do not have to be separated by a space or comma.
- See [String Formats](#) for information on column width specifications for string variables in UTF-8 format data files.

**Example**

```
DATA LIST FILE="/data/hubdata.txt" RECORDS=3
  /1 YRHIRED 14-15 DEPT 19 SEX 20
  /2 SALARY 21-25.
```

- The data are in fixed format (the default) and are read from the file *HUBDATA*.
- Three variables, *YRHIRED*, *DEPT*, and *SEX*, are defined on the first record of the *HUBDATA* file. One variable, *SALARY*, is read from columns 21 through 25 on the second record. The total number of records per case is specified as 3 even though no variables are defined on the third record. The third record is simply skipped in data definition.

**Example**

```
DATA LIST FILE="/data/hubdata.txt" RECORDS=3
  /1 DEPT 19 SEX 20 YRHIRED 14-15 MOHIRED 12-13 HIRED 12-15
  /2 SALARY 21-25.
```

- The first two defined variables are *DEPT* and *SEX*, located in columns 19 and 20 on record 1. The next three variables, *YRHIRED*, *MOHIRED*, and *HIRED*, are also located on the first record.
- *YRHIRED* is read from columns 14 and 15, *MOHIRED* from columns 12 and 13, and *HIRED* from columns 12 through 15. The variable *HIRED* is a four-column variable with the first two columns representing the month when an employee was hired (the same as *MOHIRED*) and the last two columns representing the year of employment (the same as *YRHIRED*).
- The order of the variables in the dictionary is the order in which they are defined on `DATA LIST`, not their sequence in the *HUBDATA* file.

**Example**

```
DATA LIST FILE="/data/hubdata.txt" RECORDS=3
  /1 DEPT 19 SEX 20 MOHIRED YRHIRED 12-15
  /2 SALARY 21-25.
```

- A single column specification follows *MOHIRED* and *YRHIRED*. `DATA LIST` divides the total number of columns specified equally between the two variables. Thus, each variable has a width of two columns.



### Example

\* Mixing column-style and FORTRAN-like format specifications.

```
DATA LIST FILE=PRSNL / LNAME M_INIT STREET (A20,A1,1X,A10)
      AGE 35-36.
```

- FORTRAN-like format specifications are used for string variables *LNAME*, *M\_INIT*, and *STREET*. These variables must be adjacent in the data file. *LNAME* is 20 bytes wide and is located in columns 1–20. *M\_INIT* is one byte wide and is located in column 21. The `1X` specification defines a blank column between *M\_INIT* and *STREET*. *STREET* is 10 bytes wide and is located in columns 23–32.
- A column-style format is used for the variable *AGE*. *AGE* begins in column 35, ends in column 36, and by default has numeric format.

### Freefield Data

- In freefield data, column location is irrelevant since values are not in fixed column positions. Instead, values are simply separated from each other by blanks, commas, or a specified delimiter. Any number of consecutive blanks are interpreted as one delimiter unless a blank space is explicitly specified as the value delimiter. A value cannot be split across records.
- If there are not enough values to complete the last case, a warning is issued and the incomplete case is dropped.
- The specified delimiter can only be used within data values if the value is enclosed in quotes.
- To include a single quote (apostrophe) in a string value, enclose the value in double quotes. To include double quotes in a string value, enclose the value in single quotes. [For more information, see the topic String Values in Command Specifications on p. 40.](#)
- See [String Formats](#) for information on width specifications for string variables in UTF-8 data files.

### Variable Formats

Two types of format specifications are available: column-style and FORTRAN-like. With each type, you can specify both numeric and string formats. The difference between the two types is that FORTRAN-like formats include the width of the variable and column-style formats do not.

- Column-style formats are available only for fixed-format data.
- Column-style and FORTRAN-like formats can be mixed on the same `DATA LIST` to define fixed-format data.
- A value that cannot be read according to the format type specified is assigned the system-missing value and a warning message is issued.

The following sections discuss the rules for specifying column-style and FORTRAN-like formats, followed by additional considerations for numeric and string formats.

### ***Column-Style Format Specifications***

The following rules apply to column-style formats:

- Data must be in a fixed format.
- Column locations must be specified after variable names. The width of a variable is determined by the number of specified columns. [For more information, see the topic Fixed-Format Data on p. 561.](#)
- Following the column location, specify the format type in parentheses. The format type applies only to the variable or the list of variables associated with the column location specification immediately before it. If no format type is specified, numeric (F) format is used.
- To include decimal positions in the format, specify the format type followed by a comma and the number of decimal positions. For example, (DOLLAR) specifies only whole dollar amounts, and (DOLLAR, 2) specifies DOLLAR format with two decimal positions.
- Since column positions are explicitly specified, the variables can be named in any order.

### ***FORTRAN-like Format Specifications***

The following rules apply to FORTRAN-like formats:

- Data can be in either fixed or freefield format.
- Column locations cannot be specified. The width of a variable is determined by the width portion (w) of the format specification. The width must specify the number of bytes in the widest value.
- One format specification applies to only one variable. The format is specified in parentheses after the variable to which it applies. Alternatively, a variable list can be followed by an equal number of format specifications contained in one set of parentheses. When a number of consecutive variables have the same format, the number can be used as a multiplying factor preceding the format. For example, (3F5.2) assigns the format F5.2 to three consecutive variables.
- For fixed data, the number of formats specified (either explicitly or implied by the multiplication factor) must be the same as the number of variables. Otherwise, the program issues an error message. If no formats are specified, all variables have the default format F8.2.
- For freefield data, variables with no specified formats take the default F8.2 format. However, an asterisk (\*) must be used to indicate where the default format stops. Otherwise, the program tries to apply the next specified format to every variable before it and issues an error message if the number of formats specified is less than the number of variables.
- For freefield data, width and decimal specifications are not used to read the data but are assigned as print and write formats for the variable.
- For fixed data, T<sub>n</sub> can be used before a format to indicate that the variable begins at the *n*th column, and nX can be used to skip *n* columns before reading the variable. When T<sub>n</sub> is specified, variables named do not have to follow the order of the variables in the data.
- For freefield data, variables are located according to the sequence in which they are named on DATA LIST. The order of variables on DATA LIST must correspond to the order of variables in the data.

- To include decimal positions in the format for fixed-format data, specify the total width followed by a decimal point and the number of decimal positions. For example, (DOLLAR5) specifies a five-column DOLLAR format without decimal positions, and (DOLLAR5.2) specifies a five-column DOLLAR format, two columns of which are decimal positions.

## Numeric Formats

- Format specifications on DATA LIST are input formats. Based on the width specification and format type, the program generates output (print and write) formats for each variable. The program automatically expands the output format to accommodate punctuation characters such as decimal points, commas, dollar signs, or date and time delimiters. (The program does not automatically expand the output formats you assign on the FORMATS, PRINT FORMATS, and WRITE FORMATS commands. For information on assigning output formats, refer to these commands.)
- Scientific notation is accepted in input data with F, COMMA, DOLLAR, DOT, and PCT formats. The same rules apply to these formats as to E format. The values 1.234E3, 1.234+3, and 1.234E 3 are all legitimate. The last value (with a blank space) will cause freefield data to be misread and therefore should be avoided when LIST or FREE is specified.

## Implied Decimal Positions

- For fixed-format data, decimal positions can be coded in the data or implied by the format. If decimal positions are implied but are not entered in the data, the program interprets the rightmost digits in each value as the decimal digits. A coded decimal point in a value overrides the number of implied decimal places. For example, (DOLLAR, 2) specifies two decimal positions. The value 123 is interpreted as 1.23; however, the value 12.3 is interpreted as 12.3 because the coded decimal position overrides the number of implied decimal positions.
- For freefield data, decimal positions cannot be implied but must be coded in the data. If decimal positions are specified in the format but a data value does not include a decimal point, the program fills the decimal places with zeros. For example, with F3.1 format (three columns with one decimal place), the value 22 is displayed as 22.0. If a value in the data has more decimal digits than are specified in the format, the additional decimals are truncated in displayed output (but not in calculations). For example, with F3.1 format, the value 2.22 is displayed as 2.2 even though in calculations it remains 2.22.

The table below compares how values are interpreted for fixed and freefield formats. Values in the table are for a four-column numeric variable.

Table 52-1  
Interpretation of values in fixed and freefield format

Values	Fixed		Freefield	
	Default	Two defined decimal places	Default	Two defined decimal places
2001	2001	20.01	2001.00	2001.00
201	201	2.01	201.00	201.00
-201	-201	-2.01	-201.00	-201.00
2	2	.02	2.00	2.00
20	20	.20	20.00	20.00

Values	Fixed		Freefield	
	Default	Two defined decimal places	Default	Two defined decimal places
2.2	2.2	2.2	2.20	2.20
.201	.201	.201	.201	.201
2 01	Undefined	Undefined	Two values	Two values

**Example**

```
DATA LIST
  /MODEL 1 RATE 2-6(PCT,2) COST 7-11(DOLLAR) READY 12-21(ADATE) .
BEGIN DATA
1935 7878811-07-1988
2 16754654606-08-1989
3 17684783612-09-1989
END DATA.
```

- Data are inline and in fixed format (the default).
- Each variable is followed by its column location. After the column location, a column-style format is specified in parentheses.
- *MODEL* begins in column 1, is one column wide, and receives the default numeric F format.
- *RATE* begins in column 2 and ends in column 6. The PCT format is specified with two decimal places. A comma is used to separate the format type from the number of decimal places. Decimal points are not coded in the data. Thus, the program reads the rightmost digits of each value as decimal digits. The value 935 for the first case in the data is interpreted as 9.35. Note that it does not matter where numbers are entered within the column width.
- *COST* begins in column 7 and ends in column 11. DOLLAR format is specified.
- *READY* begins in column 12 and ends in column 21. ADATE format is specified.

**Example**

```
DATA LIST FILE="/data/data1.txt"
  /MODEL (F1) RATE (PCT5.2) COST (DOLLAR5) READY (ADATE10) .
```

- In this example, the FILE subcommand is used because the data are in an external file.
- The variable definition is the same as in the preceding example except that FORTRAN-like format specifications are used rather than column-style. Column locations are not specified. Instead, the format specifications include a width for each format type.
- The width (w) portion of each format must specify the total number of bytes in the widest value. DOLLAR5 format for *COST* accepts the five-digit value 78788, which displays as \$78,788. Thus, the specified input format DOLLAR5 generates an output format DOLLAR7. The program automatically expands the width of the output format to accommodate the dollar sign and comma in displayed output.

## String Formats

String (alphanumeric) variables can contain any numbers, letters, or characters, including special characters and embedded blanks. Numbers entered as values for string variables cannot be used in calculations unless you convert them to numeric format (see `RECODE`). On `DATA LIST`, a string variable is defined with an `A` format if data are in standard character form or an `AHEX` format if data are in hexadecimal form.

- For fixed-format data, the width of a string variable is either implied by the column location specification or specified by the `w` on the FORTRAN-like format. For freefield data, the width must be specified on the FORTRAN-like format.
- `AHEX` format is available only for fixed-format data. Since each set of two hexadecimal characters represents one standard character, the width specification must be an even number. The output format for a variable in `AHEX` format is `A` format with half the specified width.
- If a string in the data is longer than its specified width, the string is truncated and a warning message is displayed. If the string in the data is shorter, it is right-padded with blanks and no warning message is displayed.
- For fixed-format data, all characters within the specified or implied columns, including leading, trailing, and embedded blanks and punctuation marks, are read as the value of the string.
- For freefield data without a specified delimiter, string values in the data must be enclosed in quotes if the string contains a blank or a comma. Otherwise, the blank or comma is treated as a delimiter between values. [For more information, see the topic String Values in Command Specifications on p. 40.](#)

### Example

```
DATA LIST FILE="/data/wins.txt" FREE /POSTPOS NWINS * POSNAME (A24).
```

- `POSNAME` is specified as a 24-byte string. The asterisk preceding `POSNAME` indicates that `POSTPOS` and `NWINS` are read with the default format. If the asterisk was not specified, the program would apply the `A24` format to `POSNAME` and then issue an error message indicating that there are more variables than specified formats.

### Example

```
DATA LIST FILE="/data/wins.txt" FREE /POSTPOS * NWINS (A5) POSWINS.
```

- Both `POSTPOS` and `POSWINS` receive the default numeric format `F8.2`.
- `NWINS` receives the specified format of `A5`.

### String Width in Code Page, UTF-8, and UTF-16 Files

“Column” and width specifications for numbers are the same for all encodings. Each digit represents a single column. For example, a column specification of `VAR1 1-4 (F)` or a width specification of `VAR1 (F4)` defines a width wide enough for a 4-digit number.

For string variables, however, column and width specifications can be different for different encodings.

- For code page and UTF-8 files, column and width specifications for alphabetic characters represent bytes.
- In code page files, Latin (Roman) characters are each one byte. This includes characters with diacritical marks, such as accents. For example, resume and résumé both have a width of 6 bytes (or columns) in a code page file. Asian characters, however, require two bytes for each character.
- In UTF-8 files, Latin characters without diacritical marks are each one byte, but characters with diacritical marks require two bytes. For example, resume requires a width of 6 bytes while résumé requires a width of 8 bytes.
- Characters from Greek, Cyrillic, Coptic, Armenian, Hebrew, Arabic, Syriac and Tana alphabets each require two bytes in UTF-8 files.
- Characters from Asian alphabets require 3 or more bytes for each character in UTF-8 files.
- For UTF-16 files, column and width specifications represent characters. Therefore, each character requires one column for all alphabets.

For more information on Unicode encoding, go to <http://www.unicode.org>.

# DATAFILE ATTRIBUTE

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=name('value') name('value')...
           arrayname[1]('value') arrayname[2]('value')...
DELETE=name name...
           arrayname[n] arrayname...
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 14.0

- Command introduced.

## Example

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=OriginalVersion ('1').
```

## Overview

DATAFILE ATTRIBUTE provides the ability for you to define your own data file attributes and assign attribute values to the active dataset.

- User-defined data file attributes are saved with the data file in the data dictionary.
- The DATAFILE ATTRIBUTE command takes effect immediately, updating the data dictionary without requiring a data pass.
- You can display a list of data file and variable attributes with DISPLAY ATTRIBUTES. [For more information, see the topic DISPLAY on p. 648.](#)

## Basic Specification

The basic specification is:

- ATTRIBUTE keyword followed by an equals sign (=) and one or more attribute names that follow variable naming rules, with each attribute name followed by a quoted attribute value, enclosed in parentheses.

or

- DELETE keyword followed by an equals sign (=) and a list of defined attribute names or attribute arrays.

## Syntax Rules

- The keywords ATTRIBUTE and DELETE must each be followed by an equals sign (=).

- Each `ATTRIBUTE` keyword must be followed by a name that follows variable naming rules and a single, quoted attribute value, enclosed in parentheses. [For more information, see the topic Variable Names on p. 49.](#)
- Attribute names that begin with `@` are not displayed by `DISPLAY DICTIONARY` or `DISPLAY ATTRIBUTES`. They can only be displayed with `DISPLAY @ATTRIBUTES`.
- Attribute names that begin with a dollar sign (\$) are reserved for internal use.
- All attribute values must be quoted (single or double quotes), even if the values are numbers.
- Attribute values can be up to 32,767 bytes in length.

**Example**

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=OriginalVersion ('1')
             CreationDate('10/28/2004')
             RevisionDate('10/29/2004').
```

**Attribute Arrays**

If you append an integer enclosed in square brackets to the end of an attribute name, the attribute is interpreted as an array of attributes. For example:

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=FileAttribute[99]('not quite 100').
```

will create 99 attributes—*FileAttribute[01]* through *FileAttribute[99]*—and will assign the value “not quite 100” to the last one.

- Array subscripts (the value enclosed in square brackets) must be integers greater than 0. (Array subscript numbering starts with 1, not 0.)
- If the root name of an attribute array is the same as an existing attribute name, the attribute array replaces the existing attribute. If no value is assigned to the first element in the array (subscript [1]), the original attribute value is used for that element value.

With the `DELETE` keyword, the following rules apply to attribute arrays:

- If you specify `DELETE` followed by an array root name and no value in square brackets, all attributes in the array are deleted.
- If you specify `DELETE` with an array name followed by an integer value in square brackets, the specified array element is deleted and the integer values for all subsequent attributes in the array (in numeric order) are changed to reflect the new order of array elements.

**Example**

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=RevisionDate('10/29/2004').
DATAFILE ATTRIBUTE
  ATTRIBUTE=RevisionDate[2] ('10/21/2005').
DATAFILE ATTRIBUTE
  DELETE=RevisionDate[1].
DATAFILE ATTRIBUTE
  DELETE=RevisionDate.
```



- The first `DATAFILE ATTRIBUTE` command creates the attribute *RevisionDate* with a value of 10/29/2004.
- The second `DATAFILE ATTRIBUTE` command creates an array attribute named *RevisionDate*, which replaces the original attribute of the same name. Two array elements are created: *RevisionDate[1]* retains the original value of *RevisionDate*, and *RevisionDate[2]* has a value of 10/21/2005.
- The third `DATAFILE ATTRIBUTE` command deletes *RevisionDate[1]*, and the array element formerly known as *RevisionDate[2]* becomes the new *RevisionDate[1]* (with a value of 10/21/2005).
- The last `DATAFILE ATTRIBUTE` command deletes all attributes in the *RevisionDate* array, since it specifies the array root name without an integer value in brackets.

# ***DATASET ACTIVATE***

```
DATASET ACTIVATE name  
    [WINDOW={ASIS }]  
    {FRONT}
```

## ***Release History***

Release 14.0

- Command introduced.

## ***Example***

```
GET FILE='/data/mydata.sav'.  
DATASET NAME file1.  
COMPUTE AvgIncome=income/famsize.  
GET DATA /TYPE=XLS  
    /FILE='/data/exceldata.xls'.  
COMPUTE TotIncome=SUM(income1, income2, income3).  
DATASET NAME file2.  
DATASET ACTIVATE file1.
```

## ***Overview***

The `DATASET` commands (`DATASET NAME`, `DATASET ACTIVATE`, `DATASET DECLARE`, `DATASET COPY`, `DATASET CLOSE`) provide the ability to have multiple data sources open at the same time and control which open data source is active at any point in the session. Using defined dataset names, you can then:

- Merge data (for example, `MATCH FILES`, `ADD FILES`, `UPDATE`) from multiple different source types (for example, text data, database, spreadsheet) without saving each one as an external IBM® SPSS® Statistics data file first.
- Create new datasets that are subsets of open data sources (for example, males in one subset, females in another, people under a certain age in another, or original data in one set and transformed/computed values in another subset).
- Copy and paste variables, cases, and/or variable properties between two or more open data sources in the Data Editor.

The `DATASET ACTIVATE` command makes the named dataset the active dataset in the session.

- If the previous active dataset does not have a defined dataset name, it is no longer available in the session.
- If the previous active dataset has a defined dataset name, it remains available for subsequent use in its current state.
- If the named dataset does not exist, an error occurs, and the command is not executed.
- `DATASET ACTIVATE` cannot be used within transformation structures such as `DO IF`, `DO REPEAT`, or `LOOP`.

### **Basic Specification**

The basic specification for `DATASET ACTIVATE` is the command name followed by a name of a previously defined dataset. For more information, see the topic `DATASET NAME` on p. 583.

### **WINDOW keyword**

The `WINDOW` keyword controls the state of the Data Editor window associated with the dataset.

<b>ASIS</b>	The Data Editor window containing the dataset is not affected. This is the default.
<b>FRONT</b>	The Data Editor window containing the dataset is brought to the front and the dataset becomes the active dataset for dialog boxes.

### **Operations**

- Commands operate on the active dataset. The **active** dataset is the data source most recently opened (for example, by commands such as `GET DATA`, `GET SAS`, `GET STATA`, `GET TRANSLATE`) or most recently activated by a `DATASET ACTIVATE` command.

*Note:* The active dataset can also be changed by clicking anywhere in the Data Editor window of an open data source or selecting a dataset from the list of available datasets in a syntax window toolbar.

- Variables from one dataset are not available when another dataset is the active dataset.
- Transformations to the active dataset—before or after defining a dataset name—are preserved with the named dataset during the session, and any pending transformations to the active dataset are automatically executed whenever a different data source becomes the active dataset.
- Dataset names can be used in most commands that can contain references to SPSS Statistics data files.
- Wherever a dataset name, file handle (defined by the `FILE HANDLE` command), or filename can be used to refer to SPSS Statistics data files, defined dataset names take precedence over file handles, which take precedence over filenames. For example, if *file1* exists as both a dataset name and a file handle, `FILE=file1` in the `MATCH FILES` command will be interpreted as referring to the dataset named *file1*, not the file handle.

### **Example**

```
GET FILE='/data/mydata.sav'.
DATASET NAME file1.
COMPUTE AvgIncome=income/famsize.
GET DATA /TYPE=XLS
  /FILE='/data/exceldata.xls'.
COMPUTE TotIncome=SUM(income1, income2, income3).
DATASET NAME file2.
DATASET ACTIVATE file1.
```

- Reading a new data source automatically changes the active dataset; so the `GET DATA` command changes the active dataset to the data read from the Excel worksheet.

- Since the previous active dataset has a defined dataset name associated with it, it is preserved in its current state for subsequent use in the session. The “current state” includes the new variable *AvgIncome* generated by the `COMPUTE` command, since pending transformations are automatically executed before the Excel worksheet become the active dataset.
- When the dataset *file1* is activated again, any pending transformations associated with dataset *file2* are automatically executed; so the new variable *TotIncome* is preserved with the dataset.

# ***DATASET CLOSE***

```
DATASET CLOSE {name}  
              {* }  
              {ALL }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Release History***

Release 14.0

- Command introduced.

## ***Example***

```
DATASET CLOSE file1.
```

## ***Overview***

The `DATASET` commands (`DATASET NAME`, `DATASET ACTIVATE`, `DATASET DECLARE`, `DATASET COPY`, `DATASET CLOSE`) provide the ability to have multiple data sources open at the same time and control which open data source is active at any point in the session. Using defined dataset names, you can then:

- Merge data (for example, `MATCH FILES`, `ADD FILES`, `UPDATE`) from multiple different source types (for example, text data, database, spreadsheet) without saving each one as an external IBM® SPSS® Statistics data file first.
- Create new datasets that are subsets of open data sources (for example, males in one subset, females in another, people under a certain age in another, or original data in one set and transformed/computed values in another subset).
- Copy and paste variables, cases, and/or variable properties between two or more open data sources in the Data Editor.

The `DATASET CLOSE` command closes the named dataset.

- If the dataset name specified is not the active dataset, that dataset is closed and no longer available in the session.
- If the dataset name specified is the active dataset or if an asterisk (\*) is specified and the active dataset has a name, the association with that name is broken. The active dataset remains active but has no name.
- If `ALL` is specified, all associations with datasets are broken. All the datasets except the active dataset and their data windows are closed and no longer available in the session. The active dataset remains active but has no name.

***Basic Specification***

The only specification for `DATASET CLOSE` is the command name followed by a dataset name, an asterisk (\*), or the keyword `ALL`.

# DATASET COPY

```
DATASET COPY name  
    [WINDOW={MINIMIZED}]  
        {HIDDEN }  
        {FRONT  }
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 14.0

- Command introduced.

## Example

```
DATASET NAME original.  
DATASET COPY males.  
DATASET ACTIVATE males.  
SELECT IF gender=0.  
DATASET ACTIVATE original.  
DATASET COPY females.  
DATASET ACTIVATE females.  
SELECT IF gender=1.
```

## Overview

The DATASET commands (DATASET NAME, DATASET ACTIVATE, DATASET DECLARE, DATASET COPY, DATASET CLOSE) provide the ability to have multiple data sources open at the same time and control which open data source is active at any point in the session. Using defined dataset names, you can then:

- Merge data (for example, MATCH FILES, ADD FILES, UPDATE) from multiple different source types (for example, text data, database, spreadsheet) without saving each one as an external IBM® SPSS® Statistics data file first.
- Create new datasets that are subsets of open data sources (for example, males in one subset, females in another, people under a certain age in another, or original data in one set and transformed/computed values in another subset).
- Copy and paste variables, cases, and/or variable properties between two or more open data sources in the Data Editor.

The DATASET COPY command creates a new dataset that captures the current state of the active dataset. This is particularly useful for creating multiple subsets of data from the same original data source.

- If the active dataset has a defined dataset name, its name remains associated with subsequent changes.

- If this command occurs when there are transformations pending, those transformations are executed, as if EXECUTE had been run prior to making the copy; so the transformations appear in both the original and the copy. The command is illegal where EXECUTE would be illegal. If no transformations are pending, the data are not passed.
- If the specified dataset name is already associated with a dataset, a warning is issued, the old dataset is destroyed, and the specified name becomes associated with the current state of the active dataset.
- If the specified name is associated with the active dataset, it becomes associated with the current state and the active dataset becomes unnamed.

### **Basic Specification**

The basic specification for DATASET COPY is the command name followed by a new dataset name that conforms to variable naming rules. [For more information, see the topic Variable Names on p. 49.](#)

### **WINDOW Keyword**

The WINDOW keyword controls the state of the Data Editor window associated with the dataset.

<b>MINIMIZED</b>	The Data Editor window associated with the new dataset is opened in a minimized state. This is the default.
<b>HIDDEN</b>	The Data Editor window associated with the new dataset is not displayed.
<b>FRONT</b>	The Data Editor window containing the dataset is brought to the front and the dataset becomes the active dataset for dialog boxes.

### **Operations**

- Commands operate on the active dataset. The **active** dataset is the data source most recently opened (for example, by commands such as GET DATA, GET SAS, GET STATA, GET TRANSLATE) or most recently activated by a DATASET ACTIVATE command.  
*Note:* The active dataset can also be changed by clicking anywhere in the Data Editor window of an open data source or selecting a dataset from the list of available datasets in a syntax window toolbar.
- Variables from one dataset are not available when another dataset is the active dataset.
- Transformations to the active dataset—before or after defining a dataset name—are preserved with the named dataset during the session, and any pending transformations to the active dataset are automatically executed whenever a different data source becomes the active dataset.
- Dataset names can be used in most commands that can contain references to SPSS Statistics data files.
- Wherever a dataset name, file handle (defined by the FILE HANDLE command), or filename can be used to refer to SPSS Statistics data files, defined dataset names take precedence over file handles, which take precedence over filenames. For example, if *file1* exists as both a dataset name and a file handle, FILE=file1 in the MATCH FILES command will be interpreted as referring to the dataset named *file1*, not the file handle.



### **Limitations**

Because each window requires a minimum amount of memory, there is a limit to the number of windows, SPSS Statistics or otherwise, that can be concurrently open on a given system. The particular number depends on the specifications of your system and may be independent of total memory due to OS constraints.

### **Example**

```
DATASET NAME original.  
DATASET COPY males.  
DATASET ACTIVATE males.  
SELECT IF gender=0.  
DATASET ACTIVATE original.  
DATASET COPY females.  
DATASET ACTIVATE females.  
SELECT IF gender=1.
```

- The first `DATASET COPY` command creates a new dataset, *males*, that represents the state of the active dataset at the time it was copied.
- The *males* dataset is activated and a subset of males is created.
- The original dataset is activated, restoring the cases deleted from the *males* subset.
- The second `DATASET COPY` command creates a second copy of the original dataset with the name *females*, which is then activated and a subset of females is created.
- Three different versions of the initial data file are now available in the session: the original version, a version containing only data for males, and a version containing only data for females.

# ***DATASET DECLARE***

```
DATASET DECLARE name  
    [WINDOW={MINIMIZED}]  
        {HIDDEN  }  
        {FRONT   }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Release History***

Release 14.0

- Command introduced.

## ***Example***

```
DATASET DECLARE corrmatrix.  
REGRESSION  
  /DEPENDENT=var1  
  /METHOD=ENTER= var2 to var10  
  /OUTFILE=CORE(corrmatrix).  
DATASET ACTIVATE corrmatrix.
```

## ***Overview***

The `DATASET` commands (`DATASET NAME`, `DATASET ACTIVATE`, `DATASET DECLARE`, `DATASET COPY`, `DATASET CLOSE`) provide the ability to have multiple data sources open at the same time and control which open data source is active at any point in the session. Using defined dataset names, you can then:

- Merge data (for example, `MATCH FILES`, `ADD FILES`, `UPDATE`) from multiple different source types (for example, text data, database, spreadsheet) without saving each one as an external IBM® SPSS® Statistics data file first.
- Create new datasets that are subsets of open data sources (for example, males in one subset, females in another, people under a certain age in another, or original data in one set and transformed/computed values in another subset).
- Copy and paste variables, cases, and/or variable properties between two or more open data sources in the Data Editor.

The `DATASET DECLARE` command creates a new dataset name that is not associated with any open dataset. It can become associated with a dataset if it is used in a command that writes SPSS Statistics data files. This is particularly useful if you need to create temporary SPSS Statistics data files as an intermediate step in a program.

### **Basic Specification**

The basic specification for `DATASET DECLARE` is the command name followed by a new dataset name that conforms to variable naming rules. [For more information, see the topic Variable Names on p. 49.](#)

### **WINDOW Keyword**

The `WINDOW` keyword controls the state of the Data Editor window associated with the dataset.

<b>MINIMIZED</b>	The Data Editor window associated with the new dataset is opened in a minimized state. This is the default.
<b>HIDDEN</b>	The Data Editor window associated with the new dataset is not displayed.
<b>FRONT</b>	The Data Editor window containing the dataset is brought to the front and the dataset becomes the active dataset for dialog boxes.

### **Example**

```
DATASET DECLARE corrmatrix.  
REGRESSION  
  /DEPENDENT=var1  
  /METHOD=ENTER= var2 to var10  
  /OUTFILE=CORB(corrmatrix).
```

- The `DATASET DECLARE` command creates a new dataset name, *corrmatrix*, that is initially not assigned to any data source.
- The `REGRESSION` command writes a correlation matrix to a data file.
- Instead of specifying an external data file, the `OUTFILE` subcommand specifies the dataset name *corrmatrix*, which is now available for subsequent use in the session. If not explicitly saved (for example, with the `SAVE` command), this dataset will be automatically deleted at the end of the session.

# ***DATASET DISPLAY***

DATASET DISPLAY

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Release History***

Release 14.0

- Command introduced.

## ***Example***

DATASET DISPLAY.

## ***Overview***

The DATASET commands (DATASET NAME, DATASET ACTIVATE, DATASET DECLARE, DATASET COPY, DATASET CLOSE) provide the ability to have multiple data sources open at the same time and control which open data source is active at any point in the session. Using defined dataset names, you can then:

- Merge data (for example, MATCH FILES, ADD FILES, UPDATE) from multiple different source types (for example, text data, database, spreadsheet) without saving each one as an external IBM® SPSS® Statistics data file first.
- Create new datasets that are subsets of open data sources (for example, males in one subset, females in another, people under a certain age in another, or original data in one set and transformed/computed values in another subset).
- Copy and paste variables, cases, and/or variable properties between two or more open data sources in the Data Editor.

The DATASET DISPLAY command displays a list of currently available datasets. The only specification is the command name DATASET DISPLAY.

# ***DATASET NAME***

```
DATASET NAME name  
    [WINDOW={ASIS }]  
    {FRONT}
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Release History***

Release 14.0

- Command introduced.

## ***Example***

```
GET FILE=' /data/mydata.sav' .  
DATASET NAME file1.  
SORT CASES BY ID.  
GET FILE ' /data/moredata.sav'  
SORT CASES BY ID.  
DATASET NAME file2.  
GET DATA /TYPE=XLS  
    /FILE=' /data/exceldata.xls' .  
SORT CASES BY ID.  
MATCH FILES FILE=*  
    /FILE=file1  
    /FILE=file2  
    /BY ID.  
SAVE OUTFILE=' /data/mergedata.sav' .
```

## ***Overview***

The DATASET commands (DATASET NAME, DATASET ACTIVATE, DATASET DECLARE, DATASET COPY, DATASET CLOSE) provide the ability to have multiple data sources open at the same time and control which open data source is active at any point in the session. Using defined dataset names, you can then:

- Merge data (for example, MATCH FILES, ADD FILES, UPDATE) from multiple different source types (for example, text data, database, spreadsheet) without saving each one as an external IBM® SPSS® Statistics data file first.
- Create new datasets that are subsets of open data sources (for example, males in one subset, females in another, people under a certain age in another, or original data in one set and transformed/computed values in another subset).
- Copy and paste variables, cases, and/or variable properties between two or more open data sources in the Data Editor.

The `DATASET NAME` command:

- Assigns a unique name to the active dataset, which can be used in subsequent file access commands and subsequent `DATASET` commands.
- Makes the current data file available even after other data sources have been opened/activated.

The following general rules apply:

- If the active dataset already has a defined dataset name, the existing association is broken, and the new name is associated with the active file.
- If the name is already associated with another dataset, that association is broken, and the new association is created. The dataset previously associated with that name is closed and is no longer available.

### **Basic Specification**

The basic specification for `DATASET NAME` is the command name followed by a name that conforms to variable naming rules. [For more information, see the topic Variable Names on p. 49.](#)

### **WINDOW Keyword**

The `WINDOW` keyword controls the state of the Data Editor window associated with the dataset.

<b>ASIS</b>	The Data Editor window containing the dataset is not affected. This is the default.
<b>FRONT</b>	The Data Editor window containing the dataset is brought to the front and the dataset becomes the active dataset for dialog boxes.

### **Operations**

- Commands operate on the active dataset. The **active** dataset is the data source most recently opened (for example, by commands such as `GET DATA`, `GET SAS`, `GET STATA`, `GET TRANSLATE`) or most recently activated by a `DATASET ACTIVATE` command.
 

*Note:* The active dataset can also be changed by clicking anywhere in the Data Editor window of an open data source or selecting a dataset from the list of available datasets in a syntax window toolbar.
- Variables from one dataset are not available when another dataset is the active dataset.
- Transformations to the active dataset—before or after defining a dataset name—are preserved with the named dataset during the session, and any pending transformations to the active dataset are automatically executed whenever a different data source becomes the active dataset.
- Dataset names can be used in most commands that can contain references to SPSS Statistics data files.
- Wherever a dataset name, file handle (defined by the `FILE HANDLE` command), or filename can be used to refer to SPSS Statistics data files, defined dataset names take precedence over file handles, which take precedence over filenames. For example, if *file1* exists as both a dataset name and a file handle, `FILE=file1` in the `MATCH FILES` command will be interpreted as referring to the dataset named *file1*, not the file handle.

**Example**

```
GET FILE='/examples/data/mydata.sav'.
SORT CASES BY ID.
DATASET NAME mydata.
GET DATA /TYPE=XLS
  /FILE='/examples/data/excelfile.xls'.
SORT CASES BY ID.
DATASET NAME excelfile.
GET DATA /TYPE=ODBC /CONNECT=
  'DSN=MS Access Database;DBQ=/examples/data/dm_demo.mdb;'+
  'DriverId=25;FIL=MS Access;MaxBufferSize=2048;PageTimeout=5;'
  /SQL='SELECT * FROM main'.
SORT CASES BY ID.
MATCH FILES
  /FILE='mydata'
  /FILE='excelfile'
  /FILE=*
  /BY ID.
```

- A data file in SPSS Statistics format and assigned the dataset name *mydata*. Since it has been assigned a dataset name, it remains available for subsequent use even after other data sources have been opened.
- An Excel file is then read and assigned the dataset name *exceldata*. Like the SPSS Statistics data file, since it has been assigned a dataset name, it remains available after other data sources have been opened.
- Then a table from a database is read. Since it is the most recently opened or activated dataset, it is the active dataset.
- The three datasets are then merged together with `MATCH FILES` command, using the dataset names on the `FILE` subcommands instead of file names.
- An asterisk (\*) is used to specify the active dataset, which is the database table in this example.
- The files are merged together based on the value of the key variable *ID*, specified on the `BY` subcommand.
- Since all the files being merged need to be sorted in the same order of the key variable(s), `SORT CASES` is performed on each dataset.

# DATE

```
DATE keyword [starting value [periodicity]]
      [keyword [starting value [periodicity]]]
      [BY increment]
```

*Keywords for long time periods:*

<b>Keyword</b>	<b>Abbreviation</b>	<b>Default starting value</b>	<b>Default periodicity</b>
YEAR	Y	1	none
QUARTER	Q	1	4
MONTH	M	1	12

*Keywords for short time periods:*

<b>Keyword</b>	<b>Abbreviation</b>	<b>Default starting value</b>	<b>Default periodicity</b>
WEEK	W	1	none
DAY	D	1	7
HOUR	H	0	24
MINUTE	MI	0	60
SECOND	S	0	60

*Keywords for any time periods:*

<b>Keyword</b>	<b>Abbreviation</b>	<b>Default starting value</b>	<b>Default periodicity</b>
CYCLE	C	1	none
OBS	O	none	none

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
DATE Y 1960 M.
```

## **Overview**

DATE generates date identification variables. You can use these variables to label plots and other output, establish periodicity, and distinguish between historical, validation, and forecasting periods.



### **Options**

You can specify the starting value and periodicity. You can also specify an increment for the lowest-order keyword specified.

### **Basic Specification**

The basic specification on DATE is a single keyword.

- For each keyword specified, DATE creates a numeric variable whose name is the keyword with an underscore as a suffix. Values for this variable are assigned to observations sequentially, beginning with the specified starting value. DATE also creates a string variable named *DATE\_*, which combines the information from the numeric date variables and is used for labeling.
- If no starting value is specified, either the default is used or the value is inferred from the starting value of another DATE keyword.
- All variables created by DATE are automatically assigned variable labels that describe periodicity and associated formats. DATE produces a list of the names of the variables it creates and their variable labels.

### **Subcommand Order**

- Keywords can be specified in any order.

### **Operations**

- DATE creates a numeric variable for every keyword specified, plus a string variable *DATE\_*, which combines information from all the specified keywords.
- DATE automatically creates variable labels for each keyword specified indicating the variable name and its periodicity. For the *DATE\_* variable, the label indicates the variable name and format.
- If the highest-order DATE variable specified has a periodicity, the *CYCLE\_* variable will automatically be created. *CYCLE\_* cannot have a periodicity. [For more information, see the topic Example 3 on p. 590.](#)
- Default periodicities are not used for the highest-order keyword specified. The exception is QUARTER, which will always have a default periodicity.
- The periodicity of the lowest-order variable is the default periodicity used by the procedures when periodicity is not defined either within the procedure or by the TSET command.
- The keyword name with an underscore is always used as the new variable name, even if keyword abbreviations are used in the specifications.
- Each time the DATE command is used, any DATE variables already in the active dataset are deleted.
- The DATE command invalidates any previous USE and PREDICT commands specified. The USE and PREDICT periods must be respecified after DATE.

**Limitations**

- There is no limit on the number of keywords on the DATE command. However, keywords that describe long time periods (YEAR, QUARTER, MONTH) cannot be used on the same command with keywords that describe short time periods (WEEK, DAY, HOUR, MINUTE, SECOND).
- User-defined variable names must not conflict with DATE variable names.

**Syntax Rules**

- You can specify more than one keyword per command.
- If a keyword is specified more than once, only the last one is executed.
- Keywords that describe long time periods (YEAR, QUARTER, MONTH) cannot be used on the same command with keywords that describe short time periods (WEEK, DAY, HOUR, MINUTE, SECOND).
- Keywords CYCLE and OBS can be used with any other keyword.
- The lowest-order keyword specified should correspond to the level at which observations occur. For example, if observations are daily, the lowest-order keyword should be DAY.
- Keywords (except MINUTE) can be abbreviated down to the first character. MINUTE must have at least two characters (MI) to distinguish it from keyword MONTH.
- Keywords and additional specifications are separated by commas or spaces.

**Starting Value and Periodicity**

- A starting value and periodicity can be entered for any keyword except CYCLE. CYCLE can have only a starting value.
- Starting value and periodicity *must* be specified for keyword OBS.
- The starting value is specified first, followed by the periodicity, if any.
- You cannot specify a periodicity without first specifying a starting value.
- Starting values for HOUR, MINUTE, and SECOND can range from 0 to the periodicity minus 1 (for example, 0 to 59). For all other keywords, the range is 1 to the periodicity.
- If both MONTH and QUARTER are specified, DATE can infer the starting value of one from the other. [For more information, see the topic Example 5 on p. 592.](#)
- Specifying conflicting starting values for MONTH and QUARTER, such as Q 1 M 4, results in an error.
- For keyword YEAR, the starting value can be specified as the last two digits (93) instead of the whole year (1993) when the series and any forecasting are all within the same century. The same format (2 digits or 4 digits) must be used in all other commands that use year values.
- If you specify keywords that describe short time periods and skip over a level of measurement (for example, if you specify HOUR and SECOND but not MINUTE), you must specify the starting value and periodicity of the keyword after the skipped keywords. Otherwise, inappropriate periodicities will be generated. [For more information, see the topic Example 7 on p. 593.](#)

## BY Keyword

- Keyword `BY` and a positive integer can be specified after the lowest-order keyword on the command to indicate an increment value. This value indicates how much to increment values of the lowest-order date variable as they are assigned to observations. [For more information, see the topic Example 4 on p. 591.](#)
- The increment value must divide evenly into the periodicity of the lowest-order `DATE` variable specified.

## Example 1

```
DATE Y 1960 M.
```

- This command generates variables `DATE_`, `YEAR_`, and `MONTH_`.
- `YEAR_` has a starting value of 1960. `MONTH_` starts at the default value of 1.
- By default, `YEAR_` has no periodicity, and `MONTH_` has a periodicity of 12.

DATE reports the following:

Name	Label
<code>YEAR_</code>	YEAR, not periodic
<code>MONTH_</code>	MONTH, period 12
<code>DATE_</code>	DATE. FORMAT: "MMM YYYY"

The following is a partial listing of the new variables:

YEAR_	MONTH_	DATE_
1960	1	JAN 1960
1960	2	FEB 1960
1960	3	MAR 1960
1960	4	APR 1960
...		
1960	10	OCT 1960
1960	11	NOV 1960
1960	12	DEC 1960
1961	1	JAN 1961
1961	2	FEB 1961
...		
1999	4	APR 1999
1999	5	MAY 1999
1999	6	JUN 1999

## Example 2

```
DATE WEEK DAY 1 5 HOUR 1 8.
```

- This command creates four variables (`DATE_`, `WEEK_`, `DAY_`, and `HOUR_`) in a file where observations occur hourly in a 5-day, 40-hour week.
- For `WEEK`, the default starting value is 1 and the default periodicity is none.

## DATE

- For *DAY\_*, the starting value has to be specified, even though it is the same as the default, because a periodicity is specified. The periodicity of 5 means that observations are measured in a 5-day week.
- For *HOURL\_*, a starting value of 1 is specified. The periodicity of 8 means that observations occur in an 8-hour day.

DATE reports the following:

Name	Label
WEEK_	WEEK, not periodic
DAY_	DAY, period 5
HOURL_	HOURL, period 24
DATE_	DATE. FORMAT: "WWW D HH"

The following is a partial listing of the new variables:

WEEK_	DAY_	HOURL_	DATE_
1	1	1	1 1 1
1	1	2	1 1 2
1	1	3	1 1 3
1	1	4	1 1 4
1	1	5	1 1 5
...			
1	1	22	1 1 22
1	1	23	1 1 23
1	2	0	1 2 0
1	2	1	1 2 1
1	2	2	1 2 2
...			
4	5	16	4 5 16
4	5	17	4 5 17
4	5	18	4 5 18

### Example 3

DATE DAY 1 5 HOURL 3 8.

- This command creates four variables (*DATE\_*, *CYCLE\_*, *DAY\_*, and *HOURL\_*) in a file where observations occur hourly.
- For *HOURL\_*, the starting value is 3 and the periodicity is 8.
- For *DAY\_*, the starting value is 1 and the periodicity is 5. Since *DAY\_* is the highest-order variable and it has a periodicity assigned, variable *CYCLE\_* is automatically created.

DATE reports the following:

Name	Label
CYCLE_	CYCLE, not periodic
DAY_	DAY, period 5
HOURL_	HOURL, period 8
DATE_	DATE. FORMAT: "CCCC D H"

The following is a partial listing of the new variables:

```

CYCLE_ DAY_ HOUR_ DATE_
      1   1   3     1 1 3
      1   1   4     1 1 4
      1   1   5     1 1 5
      1   1   6     1 1 6
      1   1   7     1 1 7
      1   2   0     1 2 0
      1   2   1     1 2 1
      ...
     12   4   6     12 4 6
     12   4   7     12 4 7
     12   5   0     12 5 0
     12   5   1     12 5 1
     12   5   2     12 5 2
     12   5   3     12 5 3
     12   5   4     12 5 4

```

### Example 4

```
DATE DAY HOUR 1 24 BY 2.
```

- This command creates three variables (*DATE\_*, *DAY\_*, and *HOUR\_*) in a file where observations occur every two hours in a 24-hour day.
- *DAY\_* uses the default starting value of 1. It has no periodicity, since none is specified, and it is the highest-order keyword on the command.
- *HOUR\_* starts with a value of 1 and has a periodicity of 24.
- Keyword *BY* specifies an increment of 2 to use in assigning hour values.

DATE reports the following:

Name	Label
DAY_	DAY, not periodic
HOUR_	HOUR, period 24 by 2
DATE_	DATE. FORMAT: "DDDD HH"

The following is a partial listing of the new variables:

```

DAY_ HOUR_ DATE_
      1   1     1 1
      1   3     1 3
      1   5     1 5
      ...
     39  17    39 17
     39  19    39 19
     39  21    39 21
     39  23    39 23
     40   1    40  1
     40   3    40  3
     40   5    40  5
     40   7    40  7
     40   9    40  9
     40  11    40 11

```

## Example 5

DATE Y 1950 Q 2 M.

- This example creates four variables (*DATE\_*, *YEAR\_*, *QUARTER\_*, and *MONTH\_*) in a file where observations are quarterly, starting with April 1950.
- The starting value for *MONTH\_* is inferred from *QUARTER\_*.
- This specification is equivalent to DATE Y 1950 Q M 4. Here, the starting value for *QUARTER\_* (2) would be inferred from MONTH.

DATE reports the following:

Name	Label
YEAR_	YEAR, not periodic
QUARTER_	QUARTER, period 4
MONTH_	MONTH, period 12
DATE_	DATE. FORMAT: "MMM YYYY"

The following is a partial listing of the new variables:

YEAR_	QUARTER_	MONTH_	DATE_
1950	2	4	APR 1950
1950	2	5	MAY 1950
1950	2	6	JUN 1950
1950	3	7	JUL 1950
1950	3	8	AUG 1950
...			
1988	4	11	NOV 1988
1988	4	12	DEC 1988
1989	1	1	JAN 1989
1989	1	2	FEB 1989
1989	1	3	MAR 1989
1989	2	4	APR 1989
1989	2	5	MAY 1989
1989	2	6	JUN 1989
1989	3	7	JUL 1989
1989	3	8	AUG 1989
1989	3	9	SEP 1989

## Example 6

DATE OBS 9 17.

- This command creates variables *DATE\_*, *CYCLE\_*, and *OBS\_* and assigns values to observations sequentially, starting with value 9. The periodicity is 17.

DATE reports the following:

Name	Label
CYCLE_	CYCLE, not periodic
OBS_	OBS, period 17
DATE_	DATE. FORMAT: "CCCC OO"

The following is a partial listing of the new variables:

```

CYCLE_  OBS_  DATE_
      1   9    1   9
      1  10    1  10
      1  11    1  11
      1  12    1  12
      1  13    1  13
      1  14    1  14
      1  15    1  15
      1  16    1  16
      1  17    1  17
      2   1    2   1
      2   2    2   2
      ...
     28  15   28  15
     28  16   28  16
     28  17   28  17
     29   1   29   1
     29   2   29   2
     29   3   29   3
     29   4   29   4
     29   5   29   5
     29   6   29   6

```

## Example 7

```
DATE W H 1 168
```

- This example creates three variables (*DATE\_*, *WEEK\_*, and *HOUR\_*) in a file where observations occur hourly.
- Since the `DAY` keyword is not specified, a periodicity must be specified for `HOUR`. The value 168 indicates that there are 168 hours in a week.
- The starting value of `HOUR` is specified as 1.

DATE reports the following:

Name	Label
WEEK_	WEEK, not periodic
HOUR_	HOUR, period 168
DATE_	DATE. FORMAT: "WWWW HHH"

The following is a partial listing of the new variables:

```

WEEK_  HOUR_  DATE_
      1    1    1    1
      1    2    1    2
      1    3    1    3
      1    4    1    4
      1    5    1    5
      1    6    1    6
      ...
     1  161    1  161
     1  162    1  162
     1  163    1  163
     1  164    1  164
     1  165    1  165

```

---

*DATE*

1	166	1	166
1	167	1	167
2	0	2	0
2	1	2	1
2	2	2	2
2	3	2	3
2	4	2	4
2	5	2	5
...			
3	131	3	131
3	132	3	132
3	133	3	133
3	134	3	134
3	135	3	135
3	136	3	136
3	137	3	137
3	138	3	138



# DEFINE-!ENDDDEFINE

```
DEFINE macro name

  ([{argument name=} [!DEFAULT (string)] [!NOEXPAND] {!TOKENS (n)           }]
   {!POSITIONAL= }                                     {!CHAREND ('char')           }
                                                    {!ENCLOSE ('char', 'char')}
                                                    {!CMDEND                    })

  [{argument name=} ...]
  {!POSITIONAL= }

macro body

!ENDDDEFINE
```

## *SET command controls:*

```
PRESERVE
RESTORE
```

## *Assignment:*

```
!LET var=expression
```

## *Conditional processing:*

```
!IF (expression) !THEN statements
  [!ELSE statements]
!IFEND
```

## *Looping constructs:*

```
!DO !varname=start !TO finish [!BY step]
  statements [!BREAK]
!DOEND

!DO !varname !IN (list)
  statements [!BREAK]
!DOEND
```

## *Macro directives:*

```
!OFFEXPAND
!ONEXPAND
```

## *String manipulation functions:*

```
!LENGTH (string)
!CONCAT (string1,string2)
!SUBSTR (string,from,[length])
!INDEX (string1,string2)
!HEAD (string)
!TAIL (string)
!QUOTE (string)
!UNQUOTE (string)
!UPCASE (string)
!BLANKS (n)
!NULL
!EVAL (string)
```

### **Release History**

Release 14.0

- For syntax processed in interactive mode, modifications to the macro facility may affect macro calls occurring at the end of a command. [For more information, see the topic Overview on p. 596.](#)

### **Example**

```
DEFINE sesvars ()  
    age sex educ religion  
!ENDDDEFINE.
```

## **Overview**

DEFINE-!ENDDDEFINE defines a program macro, which can then be used within a command sequence. A macro can be useful in several different contexts. For example, it can be used to:

- Issue a series of the same or similar commands repeatedly, using looping constructs rather than redundant specifications.
- Specify a set of variables.
- Produce output from several program procedures with a single command.
- Create complex input programs, procedure specifications, or whole sessions that can then be executed.

A macro is defined by specifying any part of a valid command and giving it a macro name. This name is then specified in a macro call within a command sequence. When the program encounters the macro name, it expands the macro.

In the examples of macro definition throughout this reference, the macro name, body, and arguments are shown in lowercase for readability. Macro keywords, which are always preceded by an exclamation point (!), are shown in uppercase.

### **Options**

**Macro Arguments.** You can declare and use arguments in the macro definition and then assign specific values to these arguments in the macro call. You can define defaults for the arguments and indicate whether an argument should be expanded when the macro is called. [For more information, see the topic Macro Arguments on p. 600.](#)

**Macro Directives.** You can turn macro expansion on and off. [For more information, see the topic Macro Directives on p. 607.](#)

**String Manipulation Functions.** You can process one or more character strings and produce either a new character string or a character representation of a numeric result. [For more information, see the topic String Manipulation Functions on p. 607.](#)

**Conditional Processing.** You can build conditional and looping constructs. [For more information, see the topic Conditional Processing on p. 610.](#)

**Macro Variables.** You can directly assign values to macro variables [For more information, see the topic Direct Assignment of Macro Variables on p. 613.](#)

### ***Basic Specification***

All macros must start with `DEFINE` and end with `!ENDDFINE`. These commands identify the beginning and end of a macro definition and are used to separate the macro definition from the rest of the command sequence.

- Immediately after `DEFINE`, specify the **macro name**. All macros must have a name. The name is used in the macro call to refer to the macro. Macro names can begin with an exclamation point (!), but other than this, follow the usual naming conventions. Starting a name with an ! ensures that it will not conflict with the other text or variables in the session.
- Immediately after the macro name, specify an optional **argument** definition in parentheses. This specification indicates the arguments that will be read when the macro is called. If you do not want to include arguments, specify just the parentheses; *the parentheses are required, whether or not they enclose an argument.*
- Next specify the body of the macro. The **macro body** can include commands, parts of commands, or macro statements (macro directives, string manipulation statements, and looping and conditional processing statements).
- At the end of the macro body, specify `!ENDDFINE`.

To invoke the macro, issue a **macro call** in the command sequence. To call a macro, specify the macro name and any necessary arguments. If there are no arguments, only the macro name is required.

### ***Operations***

- When the program reads the macro definition, it translates into uppercase all text (except arguments) not enclosed in quotation marks. Arguments are read in upper- and lowercase.
- The macro facility does not build and execute commands; rather, it expands strings in a process called **macro expansion**. A macro call initiates macro expansion. After the strings are expanded, the commands (or parts of commands) that contain the expanded strings are executed as part of the command sequence.
- Any elements on the macro call that are not used in the macro expansion are read and combined with the expanded strings.
- The expanded strings and the remaining elements from the macro call, if any, must conform to the syntax rules for the program. If not, the program generates either a warning or an error message, depending on the nature of the syntax problem.

### ***Syntax Rules***

Just like other commands, expanded macros must adhere to the rules of the processing mode under which they are run. While it is desirable to create macro syntax that will run in both interactive and batch modes, this may sometimes add a layer of complexity that you may want to avoid. So we recommend that you write macro syntax that adheres to interactive syntax rules and structure your jobs to execute macro syntax under interactive syntax rules.

- The macro `!ENDDFINE` statement should end with a period. A period as the last character on a line is interpreted as a command terminator in interactive mode.
- Other macro statements (for example, `!IF`, `!LOOP`, `!LET`) should not end with a period.
- Text within the body of the macro that represent commands that will be generated when the macro is expanded should include the period at the end of each command, and each command should start on a new line.

The macro body cannot contain the characters `\` or `^` unless they are contained within a quoted string, as in "a valid backslash `\` in a quoted string".

### Example

```
DEFINE !macro1(type = !DEFAULT(1) !TOKENS(1)
                /varlist=!CMDEND)
!IF (!type = 1)!THEN
frequencies variables=!varlist.
!ELSE
descriptives variables=!varlist.
!IFEND
!ENDDFINE.
```

- The macro statements `DEFINE`, `!IF`, `!ELSE`, and `!IFEND` do not end with a period.
- `!ENDDFINE` ends with a period.
- The `FREQUENCIES` and `DESCRIPTIVES` commands generated by the macro each start on a new line and end with a period.

To structure your command syntax jobs so that interactive processing rules are always used instead of batch processing rules:

- Use `INSERT` instead of `INCLUDE` to combine command files containing macros with other command files. [For more information, see the topic INSERT on p. 979.](#)
- In Production Facility jobs, select Interactive for the Syntax Input Format.
- In the Batch Facility (available only with IBM® SPSS® Statistics Server), use the `-i` switch to use interactive processing rules.

### Compatibility

Improvements to the macro facility may cause errors in jobs that previously ran without errors. Specifically, for syntax that is processed with interactive rules, if a macro call occurs at the end of a command, and there is no command terminator (either a period or a blank line), the next command after the macro expansion will be interpreted as a continuation line instead of a new command, as in:

```
DEFINE !macro1()
var1 var2 var3
!ENDDFINE.
FREQUENCIES VARIABLES = !macro1
DESCRIPTIVES VARIABLES = !macro1.
```

In interactive mode, the `DESCRIPTIVES` command will be interpreted as a continuation of the `FREQUENCIES` command, and neither command will run.



- The name of the macro is `check`. The empty parentheses indicate that there are no arguments to the macro.
- The macro definition (between `DEFINE` and `!ENDDDEFINE`) contains the command sequence to be repeated: `SPLIT FILE`, `FREQUENCIES`, `DESCRIPTIVES`, `LIST`, `SPLIT FILE`, and `REGRESSION`.
- The macro is called three times. Every time `check` is encountered, it is replaced with the command sequence `SPLIT FILE`, `FREQUENCIES`, `DESCRIPTIVES`, `LIST`, `SPLIT FILE OFF`, and `REGRESSION`. The command sequence using the macro facility is identical to the command sequence in which the specified commands are explicitly stated three separate times.

**Example**

```
* Macro with an argument.

DEFINE myfreq (vars = !CHAREND('/'))
frequencies variables = !vars
  /format = notable
  /statistics = default skewness kurtosis.
!ENDDDEFINE.

myfreq vars = age sex educ religion /.
```

- The macro definition defines `vars` as the macro argument. In the macro call, four variables are specified as the argument to the macro `myfreq`. When the program expands the `myfreq` macro, it substitutes the argument, `age`, `sex`, `educ`, and `religion`, for `!vars` and executes the resulting commands.

## Macro Arguments

The macro definition can include macro arguments, which can be assigned specific values in the macro call. There are two types of arguments: keyword and positional. Keyword arguments are assigned names in the macro definition; in the macro call, they are identified by name. Positional arguments are defined after the keyword `!POSITIONAL` in the macro definition; in the macro call, they are identified by their relative position within the macro definition.

- There is no limit to the number of arguments that can be specified in a macro.
- All arguments are specified in parentheses and must be separated by slashes.
- If both keyword and positional arguments are defined in the same definition, the positional arguments must be defined, used in the macro body, and invoked in the macro call before the keyword arguments.

**Example**

```
* A keyword argument.

DEFINE macname (arg1 = !TOKENS(1))
frequencies variables = !arg1.
!ENDDDEFINE.

macname arg1 = V1.
```

- The macro definition defines `macname` as the macro name and `arg1` as the argument. The argument `arg1` has one token and can be assigned any value in the macro call.
- The macro call expands the `macname` macro. The argument is identified by its name, `arg1`, and is assigned the value `V1`. `V1` is substituted wherever `!arg1` appears in the macro body. The macro body in this example is the `FREQUENCIES` command.

### Example

```
* A positional argument.

DEFINE macname (!POSITIONAL !TOKENS(1)
               /!POSITIONAL !TOKENS(2))
frequencies variables = !1 !2.
!ENDDDEFINE.

macname V1 V2 V3.
```

- The macro definition defines `macname` as the macro name with two positional arguments. The first argument has one token and the second argument has two tokens. The tokens can be assigned any values in the macro call.
- The macro call expands the `macname` macro. The arguments are identified by their positions. `V1` is substituted for `!1` wherever `!1` appears in the macro body. `V2` and `V3` are substituted for `!2` wherever `!2` appears in the macro body. The macro body in this example is the `FREQUENCIES` command.

## Keyword Arguments

Keyword arguments are called with user-defined keywords that can be specified in any order. In the macro body, the argument name is preceded by an exclamation point. On the macro call, the argument is specified without the exclamation point.

- Keyword argument definitions contain the argument name, an equals sign, and the `!TOKENS`, `!ENCLOSE`, `!CHAREND`, or `!CMDEND` keyword. [For more information, see the topic Assigning Tokens to Arguments on p. 603.](#)
- Argument names are limited to 63 bytes and cannot match the character portion of a macro keyword, such as `DEFINE`, `TOKENS`, `CHAREND`, and so forth.
- The keyword `!POSITIONAL` cannot be used in keyword argument definitions.
- Keyword arguments do not have to be called in the order they were defined.

### Example

```
DATA LIST FILE=MAC / V1 1-2 V2 4-5 V3 7-8.

* Macro definition.
DEFINE macdef2 (arg1 = !TOKENS(1)
               /arg2 = !TOKENS(1)
               /arg3 = !TOKENS(1))
frequencies variables = !arg1 !arg2 !arg3.
!ENDDDEFINE.

* Macro call.
macdef2 arg1=V1 arg2=V2 arg3=V3.
macdef2 arg3=V3 arg1=V1 arg2=V2.
```

- Three arguments are defined: `arg1`, `arg2`, and `arg3`, each with one token. In the first macro call, `arg1` is assigned the value `V1`, `arg2` is assigned the value `V2`, and `arg3` is assigned the value `V3`. `V1`, `V2`, and `V3` are then used as the variables in the `FREQUENCIES` command.
- The second macro call yields the same results as the first one. With keyword arguments, you do not need to call the arguments in the order in which they were defined.

## Positional Arguments

Positional arguments must be defined in the order in which they will be specified on the macro call. In the macro body, the first positional argument is referred to by `!1`, the second positional argument defined is referred to by `!2`, and so on. Similarly, the value of the first argument in the macro call is assigned to `!1`, the value of the second argument is assigned to `!2`, and so on.

- Positional arguments can be collectively referred to in the macro body by specifying `!*`. The `!*` specification concatenates arguments, separating individual arguments with a blank.

### Example

```
DATA LIST FILE='/data/mac.txt' / V1 1-2 V2 4-5 V3 7-8.

* Macro definition.
DEFINE macdef (!POS !TOKENS(1)
              /!POS !TOKENS(1)
              /!POS !TOKENS(1))
frequencies variables = !1 !2 !3.
!ENDDDEFINE.

* Macro call.
macdef V1    V2    V3.
macdef V3    V1    V2.
```

- Three positional arguments with one token each are defined. The first positional argument is referred to by `!1` on the `FREQUENCIES` command, the second by `!2`, and the third by `!3`.
- When the first call expands the macro, the first positional argument (`!1`) is assigned the value `V1`, the second positional argument (`!2`) is assigned the value `V2`, and the third positional argument (`!3`) is assigned the value `V3`.
- In the second call, the first positional argument is assigned the value `V3`, the second positional argument is assigned the value `V1`, and the third positional argument is assigned the value `V2`.

### Example

```
DEFINE macdef (!POS !TOKENS(3))
frequencies variables = !1.
!ENDDDEFINE.

macdef V1    V2    V3.
```

- This example is the same as the previous one, except that it assigns three tokens to one argument instead of assigning one token to each of three arguments. The result is the same.

### Example

```
DEFINE macdef (!POS !TOKENS(1))
```



```

                /!POS !TOKENS(1)
                /!POS !TOKENS(1)
frequencies variables = !*.
!ENDDDEFINE.

macdef V1    V2    V3.

```

- This is a third alternative for achieving the macro expansion shown in the previous two examples. It specifies three arguments but then joins them all together on one `FREQUENCIES` command using the symbol `!*`.

## Assigning Tokens to Arguments

A **token** is a character or group of characters that has a predefined function in a specified context. The argument definition must include a keyword that indicates which tokens following the macro name are associated with each argument.

- Any program keyword, variable name, or delimiter (a slash, comma, and so on) is a valid token.
- The arguments for a given macro can use a combination of the token keywords.

**!TOKENS (n)**      *Assign the next n tokens to the argument.* The value *n* can be any positive integer and must be enclosed in parentheses. `!TOKENS` allows you to specify exactly how many tokens are desired.

**!CHAREND ('char')**      *Assign all tokens up to the specified character to the argument.* The character must be a one-character string specified in apostrophes and enclosed in parentheses. `!CHAREND` specifies the character that ends the argument assignment. This is useful when the number of assigned tokens is arbitrary or not known in advance.

**!ENCLOSE ('char','char')**      *Assign all tokens between the indicated characters to the argument.* The starting and ending characters can be any one-character strings, and they do not need to be the same. The characters are each enclosed in apostrophes and separated by a comma. The entire specification is enclosed in parentheses. `!ENCLOSE` allows you to group multiple tokens within a specified pair of symbols. This is useful when the number of tokens to be assigned to an argument is indeterminate, or when the use of an ending character is not sufficient.

**!CMDEND**      *Assign to the argument all of the remaining text on the macro call, up to the start of the next command.* `!CMDEND` is useful for changing the defaults on an existing command. Since `!CMDEND` reads up to the next command, only the last argument on the argument list can be specified with `!CMDEND`. If `!CMDEND` is not the final argument, the arguments following `!CMDEND` are read as text.

### Example

```

* Keyword !TOKENS.

DEFINE macname (!POSITIONAL !TOKENS (3)
frequencies variables = !1.
!ENDDDEFINE.

macname ABC DEFG HI.

```

- The three tokens following `macname` (`ABC`, `DEFG`, and `HI`) are assigned to the positional argument `!1`, and `FREQUENCIES` is then executed.

**Example**

```
* Keyword !TOKENS.

* Macro definition.
DEFINE earnrep (varrep = !TOKENS (1))
sort cases by !varrep.
report variables = earnings
  /break = !varrep
  /summary = mean.
!ENDDFINE.

* Call the macro three times.
earnrep varrep= SALESMAN. /*First macro call
earnrep varrep = REGION. /*Second macro call
earnrep varrep = MONTH. /*Third macro call
```

- This macro runs a REPORT command three times, each time with a different break variable.
- The macro name is earnrep, and there is one keyword argument, varrep, which has one token.
- In the first macro call, the token SALESMAN is substituted for !varrep when the macro is expanded. REGION and MONTH are substituted for !varrep when the macro is expanded in the second and third calls.

**Example**

```
* Keyword !CHAREND'.

DEFINE macname (!POSITIONAL !CHAREND ('/'))
  /!POSITIONAL !TOKENS(2))
frequencies variables = !1.
correlations variables= !2.
!ENDDFINE.

macname A B C D / E F.
```

- When the macro is called, all tokens up to the slash (A, B, C, and D) are assigned to the positional argument !1. E and F are assigned to the positional argument !2.

**Example**

```
* Keyword !CHAREND.

DEFINE macname (!POSITIONAL !CHAREND ('/'))
frequencies variables = !1.
!ENDDFINE.

macname A B C D / E F.
```

- Although E and F are not part of the positional argument and are not used in the macro expansion, the program still reads them as text and interprets them in relation to where the macro definition ends. In this example, macro definition ends after the expanded variable list (D). E and F are names of variables. Thus, E and F are added to the variable list and FREQUENCIES is executed with six variables: A, B, C, D, E, and F.

**Example**

```
* Keyword !ENCLOSE.

DEFINE macname (!POSITIONAL !ENCLOSE('(', ')'))
frequencies variables = !1
/statistics = default skewness.
!ENDDDEFINE.

macname (A B C) D E.
```

- When the macro is called, the three tokens enclosed in parentheses—A, B, and C—are assigned to the positional argument !1 in the macro body.
- After macro expansion is complete, the program reads the remaining characters on the macro call as text. In this instance, the macro definition ends with keyword SKEWNESS on the STATISTICS subcommand. Adding variable names to the STATISTICS subcommand is not valid syntax. The program generates a warning message but is still able to execute the frequencies command. Frequency tables and the specified statistics are generated for the variables A, B, and C.

**Example**

```
* Keyword !CMDEND'.

DEFINE macname (!POSITIONAL !TOKENS(2)
                /!POSITIONAL !CMDEND)
frequencies variables = !1.
correlations variables= !2.
!ENDDDEFINE.

macname A B C D E.
```

- When the macro is called, the first two tokens following macname (A and B) are assigned to the positional argument !1. C, D, and E are assigned to the positional argument !2. Thus, the variables used for FREQUENCIES are A and B, and the variables used for CORRELATION are C, D, and E.

**Example**

```
* Incorrect order for !CMDEND.

DEFINE macname (!POSITIONAL !CMDEND
                /!POSITIONAL !tokens(2))
frequencies variables = !1.
correlations variables= !2.
!ENDDDEFINE.

macname A B C D E.
```

- When the macro is called, all five tokens, A, B, C, D, and E, are assigned to the first positional argument. No variables are included on the variable list for CORRELATIONS, causing the program to generate an error message. The previous example declares the arguments in the correct order.

**Example**

```
* Using !CMDEND.
```

---

*DEFINE-!ENDDFINE*

```

SUBTITLE 'CHANGING DEFAULTS ON A COMMAND'.

DEFINE myfreq (!POSITIONAL !CMDEND)
frequencies !1
  /statistics=default skewness /* Modify default statistics.
!ENDDFINE.

myfreq VARIABLES = A B /HIST.

```

- The macro `myfreq` contains options for the `FREQUENCIES` command. When the macro is called, `myfreq` is expanded to perform a `FREQUENCIES` analysis on the variables *A* and *B*. The analysis produces default statistics and the skewness statistic, plus a histogram, as requested on the macro call.

**Example**

```

* Keyword arguments: Using a combination of token keywords.

DATA LIST FREE / A B C D E.
DEFINE macdef3 (arg1 = !TOKENS(1)
  /arg2 = !ENCLOSE ('(',')')
  /arg3 = !CHAREND('%'))
frequencies variables = !arg1 !arg2 !arg3.
!ENDDFINE.
macdef arg1 = A arg2=(B C) arg3=D E %.

```

- Because `arg1` is defined with the `!TOKENS` keyword, the value for `arg1` is simply specified as *A*. The value for `arg2` is specified in parentheses, as indicated by `!ENCLOSE`. The value for `arg3` is followed by a percentage sign, as indicated by `!CHAREND`.

**Defining Defaults**

The optional `!DEFAULT` keyword in the macro definition establishes default settings for arguments.

**!DEFAULT**     *Default argument.* After `!DEFAULT`, specify the value you want to use as a default for that argument. A default can be specified for each argument.

**Example**

```

DEFINE macdef (arg1 = !DEFAULT (V1) !TOKENS(1)
  /arg2 = !TOKENS(1)
  /arg3 = !TOKENS(1))
frequencies variables = !arg1 !arg2 !arg3.
!ENDDFINE.

macdef arg2=V2 arg3=V3.

```

- `V1` is defined as the default value for argument `arg1`. Since `arg1` is not specified on the macro call, it is set to `V1`.
- If `!DEFAULT (V1)` were not specified, the value of `arg1` would be set to a null string.

## Controlling Expansion

!NOEXPAND indicates that an argument should not be expanded when the macro is called.

**!NOEXPAND**            *Do not expand the specified argument.* !NOEXPAND applies to a single argument and is useful only when a macro calls another macro (embedded macros).

## Macro Directives

!ONEXPAND and !OFFEXPAND determine whether macro expansion is on or off. !ONEXPAND activates macro expansion and !OFFEXPAND stops macro expansion. All symbols between !OFFEXPAND and !ONEXPAND in the macro definition will not be expanded when the macro is called.

**!ONEXPAND**            *Turn macro expansion on.*

**!OFFEXPAND**          *Turn macro expansion off.* !OFFEXPAND is effective only when SETMEXPAND is ON (the default).

## Macro Expansion in Comments

When macro expansion is on, a macro is expanded when its name is specified in a comment line beginning with \*. To use a macro name in a comment, specify the comment within slashes and asterisks (/...\*/) to avoid unwanted macro expansion. (See COMMENT.)

## String Manipulation Functions

String manipulation functions process one or more character strings and produce either a new character string or a character representation of a numeric result.

- The result of any string manipulation function is treated as a character string.
- The arguments to string manipulation functions can be strings, variables, or even other macros. A macro argument or another function can be used in place of a string.
- The strings within string manipulation functions must be either single tokens, such as ABC, or delimited by apostrophes or quotation marks, as in 'A B C'.

Table 61-1  
Expressions and results

Expression	Result
!UPCASE(abc)	ABC
!UPCASE('abc')	ABC
!UPCASE(a b c)	error
!UPCASE('a b c')	A B C
!UPCASE(a/b/c)	error
!UPCASE('a/b/c')	A/B/C
!UPCASE(!CONCAT(a,b,c))	ABC
!UPCASE(!CONCAT('a','b','c'))	ABC
!UPCASE(!CONCAT(a, b, c))	ABC

<b>Expression</b>	<b>Result</b>
<code>!UPCASE(!CONCAT('a','b','c'))</code>	A B C
<code>!UPCASE(!CONCAT('a,b,c'))</code>	A,B,C
<code>!QUOTE(abc)</code>	'ABC'
<code>!QUOTE('abc')</code>	abc
<code>!QUOTE('Bill's')</code>	'Bill's'
<code>!QUOTE("Bill's")</code>	"Bill's"
<code>!QUOTE(Bill's)</code>	error
<code>!QUOTE(!UNQUOTE('Bill's'))</code>	'Bill's'
<b>!LENGTH (str)</b>	<i>Return the length of the specified string.</i> The result is a character representation of the string length. <code>!LENGTH(abcdef)</code> returns 6. If the string is specified with apostrophes around it, each apostrophe adds 1 to the length. <code>!LENGTH ('abcdef')</code> returns 8. If an argument is used in place of a string and it is set to null, this function will return 0.
<b>!CONCAT (str1,str2 . . .)</b>	<i>Return a string that is the concatenation of the strings.</i> For example, <code>!CONCAT (abc,def)</code> returns abcdef.
<b>!SUBSTR (str,from,[length])</b>	<i>Return a substring of the specified string.</i> The substring starts at the <i>from</i> position and continues for the specified <i>length</i> . If the length is not specified, the substring ends at the end of the input string. For example, <code>!SUBSTR (abcdef, 3, 2)</code> returns cd.
<b>!INDEX (haystack,needle)</b>	<i>Return the position of the first occurrence of the needle in the haystack.</i> If the needle is not found in the haystack, the function returns 0. <code>!INDEX (abcdef,def)</code> returns 4.
<b>!HEAD (str)</b>	<i>Return the first token within a string.</i> The input string is not changed. <code>!HEAD ('a b c')</code> returns a.
<b>!TAIL (str)</b>	<i>Return all tokens except the head token.</i> The input string is not changed. <code>!TAIL ('a b c')</code> returns b c.
<b>!QUOTE (str)</b>	<i>Put apostrophes around the argument.</i> <code>!QUOTE(abc)</code> replicates any embedded apostrophe. <code>!QUOTE(abc)</code> returns 'abc'. If !1 equals Bill's, <code>!QUOTE(!1)</code> returns 'Bill's'.
<b>!UNQUOTE (str)</b>	<i>Remove quotation marks and apostrophes from the enclosed string.</i> If !1 equals 'abc', <code>!UNQUOTE(!1)</code> is abc. Internal paired quotation marks are unpaired; if !1 equals 'Bill's', <code>!UNQUOTE(!1)</code> is Bill's. The specification <code>!UNQUOTE(!QUOTE(Bill))</code> returns Bill.
<b>!UPCASE (str)</b>	<i>Convert all lowercase characters in the argument to uppercase.</i> <code>!UPCASE('abc def')</code> returns ABC DEF.
<b>!BLANKS (n)</b>	<i>Generate a string containing the specified number of blanks.</i> The <i>n</i> specification must be a positive integer. <code>!BLANKS(5)</code> returns a string of five blank spaces. Unless the blanks are quoted, they cannot be processed, since the macro facility compresses blanks.
<b>!NULL</b>	<i>Generate a string of length 0.</i> This can help determine whether an argument was ever assigned a value, as in <code>!IF (!1 !EQ !NULL) !THEN . . .</code>
<b>!EVAL (str)</b>	<i>Scan the argument for macro calls.</i> During macro definition, an argument to a function or an operand in an expression is not scanned for possible macro calls unless the <code>!EVAL</code> function is used. It returns a string that is the expansion of its argument. For example, if <code>mac1</code> is a macro, then <code>!EVAL(mac1)</code> returns the expansion of <code>mac1</code> . If <code>mac1</code> is not a macro, <code>!EVAL(mac1)</code> returns <code>mac1</code> .

## SET Subcommands for Use with Macro

Four subcommands on the SET command were designed for use with the macro facility.

<b>MPRINT</b>	<i>Display a list of commands after macro expansion.</i> The specification on MPRINT is YES or NO (alias ON or OFF). By default, the output does not include a list of commands after macro expansion (MPRINT NO). The MPRINT subcommand on SET is independent of the PRINTBACK command.
<b>MEXPAND</b>	<i>Macro expansion.</i> The specification on MEXPAND is YES or NO (alias ON or OFF). By default, MEXPAND is on. SET MEXPAND OFF prevents macro expansion. Specifying SET MEXPAND ON reestablishes macro expansion.
<b>MNEST</b>	<i>Maximum nesting level for macros.</i> The default number of levels that can be nested is 50. The maximum number of levels depends on storage capacity.
<b>MITERATE</b>	<i>Maximum loop iterations permitted in macro expansions.</i> The default number of iterations is 1000.

## Restoring SET Specifications

The PRESERVE and RESTORE commands bring more flexibility and control over SET. PRESERVE and RESTORE are available generally within the program but are especially useful with macros.

- The settings of all SET subcommands—those set explicitly and those set by default (except MEXPAND)—are saved with PRESERVE. PRESERVE has no further specifications.
- With RESTORE, all SET subcommands are changed to what they were when the PRESERVE command was executed. RESTORE has no further specifications.
- PRESERVE . . . RESTORE sequences can be nested up to five levels.

<b>PRESERVE</b>	<i>Store the SET specifications that are in effect at this point in the session.</i>
<b>RESTORE</b>	<i>Restore the SET specifications to what they were when PRESERVE was specified.</i>

### Example

\* Two nested levels of preserve and restore'.

```
DEFINE macdef ()
preserve.
set format F5.3.
descriptives v1 v2.
+ preserve.
set format F3.0 blanks=999.
descriptives v3 v4.
+ restore.
descriptives v5 v6.
restore.
!ENDDDEFINE.
```

- The first PRESERVE command saves all of the current SET conditions. If none have been specified, the default settings are saved.
- Next, the format is set to F5.3 and descriptive statistics for V1 and V2 are obtained.
- The second PRESERVE command saves the F5.3 format setting and all other settings in effect.
- The second SET command changes the format to F3.0 and sets BLANKS to 999 (the default is SYSMIS). Descriptive statistics are then obtained for V3 and V4.

- The first RESTORE command restores the format to F5.3 and BLANKS to the default, the setting in effect at the second PRESERVE. Descriptive statistics are then obtained for V5 and V6.
- The last RESTORE restores the settings in effect when the first PRESERVE was specified.

## Conditional Processing

The !IF construct specifies conditions for processing. The syntax is as follows:

```
!IF (expression) !THEN statements
                        [!ELSE statements]
!IFEND
```

- !IF, !THEN, and !IFEND are all required. !ELSE is optional.
- If the result of the expression is true, the statements following !THEN are executed. If the result of the expression is false and !ELSE is specified, the statements following !ELSE are executed. Otherwise, the program continues.
- Valid operators for the expressions include !EQ, !NE, !GT, !LT, !GE, !LE, !OR, !NOT, and !AND, or =, ~= (≠), >, <, >=, <=, |, ~ (¬), and &.
- When a macro is expanded, conditional processing constructs are interpreted after arguments are substituted and functions are executed.
- !IF statements can be nested whenever necessary. Parentheses can be used to specify the order of evaluation. The default order is the same as for transformations: !NOT has precedence over !AND, which has precedence over !OR.

### Example

```
DEFINE mymacro(type = !DEFAULT(1) !TOKENS(1))
!IF (!type = 1)!then
frequencies variables=varone.
!ELSE
descriptives variables=vartwo.
!IFEND
!ENDDFINE.
```

## Unquoted String Constants in Conditional !IF Statements

Prior to version 12.0, under certain circumstances unquoted string constants in conditional !IF statements were not case sensitive. Starting with version 12.0, unquoted string constants are case sensitive. For backward compatibility, always use quoted string constants.

### Example

```
DEFINE noquote(type = !DEFAULT(a) !TOKENS(1))
!IF (!type = A)!THEN
frequencies variables=varone.
!ELSE
descriptives variables=vartwo.
!IFEND
!ENDDFINE.
DEFINE yesquote(type = !DEFAULT('a') !TOKENS(1)).
!IF (!type = 'A')!THEN
FREQUENCIES varone.
```



```
!ELSE
DESCRIPTIVES vartwo.
!IFEND.
!ENDDDEFINE.
```

- In the first macro, `!IF(!type = A)` is evaluated as false if the value of the unquoted string constant is lowercase ‘a’—and is therefore evaluated as false in this example.
- Prior to version 12.0, `!IF (!type = A)` was evaluated as true if the value of the unquoted string constant was lowercase ‘a’ or uppercase ‘A’—and was therefore evaluated as true in this example.
- In the second macro, `!IF (!type = 'A')` is always evaluated as false if the value of the string constant is lowercase ‘a.’

## Looping Constructs

Looping constructs accomplish repetitive tasks. Loops can be nested to whatever depth is required, but loops cannot be crossed. The macro facility has two looping constructs: the index loop (DO loop) and the list-processing loop (DO IN loop).

- When a macro is expanded, looping constructs are interpreted after arguments are substituted and functions are executed.

### Index Loop

The syntax of an index loop is as follows:

```
!DO !var = start !TO finish [ !BY step ]
    statements
!BREAK
!DOEND
```

- The indexing variable is `!var` and must begin with an exclamation point.
- The start, finish, and step values must be numbers or expressions that evaluate to numbers.
- The loop begins at the start value and continues until it reaches the finish value (unless a `!BREAK` statement is encountered). The step value is optional and can be used to specify a subset of iterations. If start is set to 1, finish to 10, and step to 3, the loop will be executed four times with the index variable assigned values 1, 4, 7, and 10.
- The statements can be any valid commands or macro keywords. `!DOEND` specifies the end of the loop.
- `!BREAK` is an optional specification. It can be used in conjunction with conditional processing to exit the loop.

#### Example

```
DEFINE macdef (arg1 = !TOKENS(1)
              /arg2 = !TOKENS(1))
!DO !i = !arg1 !TO !arg2.
frequencies variables = !CONCAT(var,!i).
!DOEND
!ENDDDEFINE.
macdef arg1 = 1 arg2 = 3.
```

- The variable *!i* is initially assigned the value 1 (*arg1*) and is incremented until it equals 3 (*arg2*), at which point the loop ends.
- The first loop concatenates *var* and the value for *!I*, which is 1 in the first loop. The second loop concatenates *var* and 2, and the third concatenates *var* and 3. The result is that `FREQUENCIES` is executed three times, with variables *VAR1*, *VAR2*, and *VAR3*, respectively.

## List-Processing Loop

The syntax of a list-processing loop is as follows:

```
!DO !var !IN (list)
    statements
!BREAK
!DOEND
```

- The `!DO` and `!DOEND` statements begin and end the loop. `!BREAK` is used to exit the loop.
- The `!IN` function requires one argument, which must be a list of items. The number of items on the list determines the number of iterations. At each iteration, the index variable *!var* is set to each item on the list.
- The list can be any expression, although it is usually a string. Only one list can be specified in each list-processing loop.

### Example

```
DEFINE macdef (!POS !CHAREND('/'))
!DO !i !IN (!1)
frequencies variables = !i.
!DOEND
!ENDDDEFINE.
macdef VAR1 VAR2 VAR3 /.
```

- The macro call assigns three variables, *VAR1*, *VAR2*, and *VAR3*, to the positional argument `!1`. Thus, the loop completes three iterations.
- In the first iteration, *!i* is set to value *VAR1*. In the second and third iterations, *!I* is set to *VAR2* and *VAR3*, respectively. Thus, `FREQUENCIES` is executed three times, respectively with *VAR1*, *VAR2*, and *VAR3*.

### Example

```
DEFINE macdef (!POS !CHAREND('/'))
!DO !i !IN (!1)
sort cases by !i.
report var = earnings
    /break = !i
    /summary = mean.
!DOEND
!ENDDDEFINE.

macdef SALESMAN REGION MONTH /.
```

- The positional argument `!1` is assigned the three variables *SALESMAN*, *REGION*, and *MONTH*. The loop is executed three times and the index variable *!i* is set to each of the variables in succession. The macro creates three reports.

## ***Direct Assignment of Macro Variables***

The macro command `!LET` assigns values to macro variables. The syntax is as follows:

```
!LET !var = expression
```

- The expression must be either a single token or enclosed in parentheses.
- The macro variable *!var* cannot be a macro keyword, and it cannot be the name of one of the arguments within the macro definition. Thus, `!LET` cannot be used to change the value of an argument.
- The macro variable *!var* can be a new variable or one previously assigned by a `!DO` command or another `!LET` command.

### ***Example***

```
!LET !a = 1
!LET !b = !CONCAT(ABC, !SUBSTR(!1, 3, 1), DEF)
!LET !c = (!2 ~= !NULL)
```

- The first `!LET` sets *!a* equal to 1.
- The second `!LET` sets *!b* equal to ABC followed by 1 character taken from the third position of *!1* followed by DEF.
- The last `!LET` sets *!c* equal to 0 (false) if *!2* is a null string or to 1 (true) if *!2* is not a null string.

# DELETE VARIABLES

```
DELETE VARIABLES varlist.
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
DELETE VARIABLES varX varY thisVar TO thatVar.
```

## **Overview**

DELETE VARIABLES deletes the specified variables from the active dataset.

### **Basic Specification**

- The basic specification is one or more variable names.

### **Syntax Rules**

- The variables must exist in the active dataset.
- The keyword TO can be used to specify consecutive variable in the active dataset.
- This command cannot be executed when there are pending transformations. For example, DELETE VARIABLES cannot be immediately preceded by transformation commands such as COMPUTE or RECODE.
- DELETE VARIABLES cannot be used with TEMPORARY.
- You cannot use this command to delete all variables in the active dataset. If the variable list includes all variables in the active dataset, an error results and the command is not executed. Use NEW FILE to delete all variables.

# DESCRIPTIVES

DESCRIPTIVES is available in the Statistics Base option.

```
DESCRIPTIVES VARIABLES= varname[ (zname)] [varname...]  
  
  [/MISSING={VARIABLE**}  [INCLUDE]]  
            {LISTWISE  }  
  
  [/SAVE]  
  
  [/STATISTICS={DEFAULT**}  [MEAN**]  [MIN**]  [SKEWNESS]]  
                [STDDEV**]  [SEMEAN]  [MAX**]  [KURTOSIS]  
                [VARIANCE ]  [SUM      ]  [RANGE]  [ALL]  
  
  [/SORT=[{MEAN      }]  [{(A)}]]  
          {SMEAN      }  {(D)}  
          {STDDEV     }  
          {VARIANCE   }  
          {KURTOSIS   }  
          {SKEWNESS   }  
          {RANGE      }  
          {MIN        }  
          {MAX        }  
          {SUM        }  
          {NAME       }
```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
DESCRIPTIVES VARIABLES=FOOD RENT, APPL TO COOK
```

## Overview

DESCRIPTIVES computes univariate statistics—including the mean, standard deviation, minimum, and maximum—for numeric variables. Because it does not sort values into a frequency table, DESCRIPTIVES is an efficient means of computing descriptive statistics for continuous variables. Other procedures that display descriptive statistics include FREQUENCIES, MEANS, and EXAMINE.

## Options

**Z Scores.** You can create new variables that contain  $z$  scores (standardized deviation scores from the mean) and add them to the active dataset by specifying  $z$ -score names on the VARIABLES subcommand or by using the SAVE subcommand.

**Statistical Display.** Optional statistics available with the STATISTICS subcommand include the standard error of the mean, variance, kurtosis, skewness, range, and sum. DESCRIPTIVES does not compute the median or mode (see FREQUENCIES or EXAMINE).

**Display Order.** You can list variables in ascending or descending alphabetical order or by the numerical value of any of the available statistics using the `SORT` subcommand.

### **Basic Specification**

The basic specification is the `VARIABLES` subcommand with a list of variables. All cases with valid values for a variable are included in the calculation of statistics for that variable. Statistics include the mean, standard deviation, minimum, maximum, and number of cases with valid values.

### **Subcommand Order**

- Subcommands can be used in any order.

### **Operations**

- If a string variable is specified on the variable list, no statistics are displayed for that variable.
- If there is insufficient memory available to calculate statistics for all variables requested, `DESCRIPTIVES` truncates the variable list.

## **Example**

```
DESCRIPTIVES
  VARIABLES=longmon tollmon equipmon cardmon wiremon
  /STATISTICS=MEAN STDDEV MIN MAX .
```

- The procedure analyzes the variables *longmon*, *tollmon*, *equipmon*, *cardmon*, and *wiremon*.
- The `STATISTICS` subcommand requests the mean, standard deviation, minimum, and maximum.

## **VARIABLES Subcommand**

`VARIABLES` names the variables for which you want to compute statistics.

- The keyword `ALL` can be used to refer to all user-defined variables in the active dataset.
- Only one variable list can be specified.

## **Z Scores**

The *z*-score transformation standardizes variables to the same scale, producing new variables with a mean of 0 and a standard deviation of 1. These variables are added to the active dataset.

- To obtain *z* scores for all specified variables, use the `SAVE` subcommand.
- To obtain *z* scores for a subset of variables, name the new variable in parentheses following the source variable on the `VARIABLES` subcommand and do not use the `SAVE` subcommand.
- Specify new names individually; a list in parentheses is not recognized.
- The new variable name can be any acceptable variable name that is not already part of the active dataset. For information on variable naming rules, see “Variable Names” on p. 36.

**Example**

```
DESCRIPTIVES VARIABLES=NTCSAL NTCPUR (PURCHZ) NTCPRI (PRICEZ) .
```

- `DESCRIPTIVES` creates  $z$ -score variables named *PURCHZ* and *PRICEZ* for *NTCPUR* and *NTCPRI*, respectively. No  $z$ -score variable is created for *NTCSAL*.

**SAVE Subcommand**

`SAVE` creates a  $z$ -score variable for each variable specified on the `VARIABLES` subcommand. The new variables are added to the active dataset.

- When `DESCRIPTIVES` creates new  $z$ -score variables, it displays the source variable names, the new variable names, and their labels in the Notes table.
- `DESCRIPTIVES` automatically supplies variable names for the new variables. The new variable name is created by prefixing the letter *Z* to the source variable name. For example, *ZNTCPRI* is the  $z$ -score variable for *NTCPRI*.
- If the default naming convention duplicates variable names in the active dataset, `DESCRIPTIVES` uses an alternative naming convention: first *ZSC001* through *ZSC099*, then *STDZ01* through *STDZ09*, then *ZZZZ01* through *ZZZZ09*, and then *ZQZQ01* through *ZQZQ09*.
- Variable labels are created by prefixing *ZSCORE* to the source variable label. If the alternative naming convention is used, `DESCRIPTIVES` prefixes *ZSCORE(varname)* to the label. If the source variable does not have a label, `DESCRIPTIVES` uses *ZSCORE(varname)* for the label.
- If you specify new names on the `VARIABLES` subcommand *and* use the `SAVE` subcommand, `DESCRIPTIVES` creates one new variable for each variable on the `VARIABLES` subcommand, using default names for variables not assigned names on `VARIABLES`.
- If at any time you want to change any of the variable names, whether those `DESCRIPTIVES` created or those you previously assigned, you can do so with the `RENAME VARIABLES` command.

**Example**

```
DESCRIPTIVES VARIABLES=ALL
  /SAVE .
```

- `SAVE` creates a  $z$ -score variable for all variables in the active dataset. All  $z$ -score variables receive the default name.

**Example**

```
DESCRIPTIVES VARIABLES=NTCSAL NTCPUR (PURCHZ) NTCPRI (PRICEZ)
  /SAVE .
```

- `DESCRIPTIVES` creates three  $z$ -score variables named *ZNTCSAL* (the default name), *PURCHZ*, and *PRICEZ*.

## STATISTICS Subcommand

By default, `DESCRIPTIVES` displays the mean, standard deviation, minimum, and maximum. Use the `STATISTICS` subcommand to request other statistics.

- When you use `STATISTICS`, `DESCRIPTIVES` displays *only* those statistics you request.
- The keyword `ALL` obtains all statistics.
- You can specify the keyword `DEFAULT` to obtain the default statistics without having to name `MEAN`, `STDDEV`, `MIN`, and `MAX`.
- The median and mode, which are available in `FREQUENCIES` and `EXAMINE`, are not available in `DESCRIPTIVES`. These statistics require that values be sorted, and `DESCRIPTIVES` does not sort values (the `SORT` subcommand does not sort values, it simply lists variables in the order you request).
- If you request a statistic that is not available, `DESCRIPTIVES` issues an error message and the command is not executed.

<b>MEAN</b>	<i>Mean.</i>
<b>SEMEAN</b>	<i>Standard error of the mean.</i>
<b>STDDEV</b>	<i>Standard deviation.</i>
<b>VARIANCE</b>	<i>Variance.</i>
<b>KURTOSIS</b>	<i>Kurtosis and standard error of kurtosis.</i>
<b>SKEWNESS</b>	<i>Skewness and standard error of skewness.</i>
<b>RANGE</b>	<i>Range.</i>
<b>MIN</b>	<i>Minimum observed value.</i>
<b>MAX</b>	<i>Maximum observed value.</i>
<b>SUM</b>	<i>Sum.</i>
<b>DEFAULT</b>	<i>Mean, standard deviation, minimum, and maximum.</i> These are the default statistics.
<b>ALL</b>	<i>All statistics available in <code>DESCRIPTIVES</code>.</i>

## SORT Subcommand

By default, `DESCRIPTIVES` lists variables in the order in which they are specified on `VARIABLES`. Use `SORT` to list variables in ascending or descending alphabetical order of variable name or in ascending or descending order of numeric value of any of the statistics.

- If you specify `SORT` without any keywords, variables are sorted in ascending order of the mean.
- `SORT` can sort variables by the value of any of the statistics available with `DESCRIPTIVES`, but only those statistics specified on `STATISTICS` (or the default statistics) are displayed.

Only one of the following keywords can be specified on `SORT`:

<b>MEAN</b>	<i>Sort by mean.</i> This is the default when <code>SORT</code> is specified without a keyword.
<b>SEMEAN</b>	<i>Sort by standard error of the mean.</i>
<b>STDDEV</b>	<i>Sort by standard deviation.</i>
<b>VARIANCE</b>	<i>Sort by variance.</i>
<b>KURTOSIS</b>	<i>Sort by kurtosis.</i>



<b>SKEWNESS</b>	<i>Sort by skewness.</i>
<b>RANGE</b>	<i>Sort by range.</i>
<b>MIN</b>	<i>Sort by minimum observed value.</i>
<b>MAX</b>	<i>Sort by maximum observed value.</i>
<b>SUM</b>	<i>Sort by sum.</i>
<b>NAME</b>	<i>Sort by variable name.</i>

Sort order can be specified in parentheses following the specified keyword:

<b>A</b>	<i>Sort in ascending order.</i> This is the default when SORT is specified without keywords.
<b>D</b>	<i>Sort in descending order.</i>

### **Example**

```
DESCRIPTIVES VARIABLES=A B C
  /STATISTICS=DEFAULT RANGE
  /SORT=RANGE (D) .
```

- DESCRIPTIVES sorts variables *A*, *B*, and *C* in descending order of range and displays the mean, standard deviation, minimum and maximum values, range, and the number of valid cases.

## **MISSING Subcommand**

MISSING controls missing values.

- By default, DESCRIPTIVES deletes cases with missing values on a variable-by-variable basis. A case with a missing value for a variable will not be included in the summary statistics for that variable, but the case *will* be included for variables where it is not missing.
- The VARIABLE and LISTWISE keywords are alternatives; however, each can be specified with INCLUDE.
- When either the keyword VARIABLE or the default missing-value treatment is used, DESCRIPTIVES reports the number of valid cases for each variable. It always displays the number of cases that would be available if listwise deletion of missing values had been selected.

<b>VARIABLE</b>	<i>Exclude cases with missing values on a variable-by-variable basis.</i> This is the default.
<b>LISTWISE</b>	<i>Exclude cases with missing values listwise.</i> Cases with missing values for any variable named are excluded from the computation of statistics for all variables.
<b>INCLUDE</b>	<i>Include user-missing values.</i>

# DETECTANOMALY

DETECTANOMALY is available in the Data Preparation option.

```
DETECTANOMALY

[/VARIABLES [CATEGORICAL=varlist]
            [SCALE=varlist]
            [ID=variable]
            [EXCEPT=varlist]]

[/HANDLEMISSING [APPLY={NO**}
                {YES }
                [CREATEMISPROPVAR={NO**}]]
                {YES }

[/CRITERIA [MINNUMPEERS={1** }
           {integer}
           [MAXNUMPEERS={15** }
           {integer}
           [MLWEIGHT={6** }
           {number}
           [NUMREASONS={1** }
           {integer}
           [PCTANOMALOUSCASES={5** }
           {number}
           [NUMANOMALOUSCASES={integer}]
           [ANOMALYCUTPOINT={2** }
           {number}
           {NONE }]]]

[/SAVE [ANOMALY[(varname)]]
       [PEERID[(varname)]]
       [PEERSIZE[(varname)]]
       [PEERPCTSIZE[(varname)]]
       [REASONVAR[(rootname)]]
       [REASONMEASURE[(rootname)]]
       [REASONVALUE[(rootname)]]
       [REASONNORM[(rootname)]]]

[/OUTFILE [MODEL=filespec]]

[/PRINT [CPS] [ANOMALYLIST**] [NORMS]
        [ANOMALYSUMMARY] [REASONSUMMARY]
        [NONE]]
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 14.0

- Command introduced.

## Example

```
DETECTANOMALY .
```

## Overview

The Anomaly Detection procedure searches for unusual cases based on deviations from the norms of their cluster groups. The procedure is designed to quickly detect unusual cases for data-auditing purposes in the exploratory data analysis step, prior to any inferential data analysis. This algorithm is designed for generic anomaly detection; that is, the definition of an anomalous case is not specific to any particular application, such as detection of unusual payment patterns in the healthcare industry or detection of money laundering in the finance industry, in which the definition of an anomaly can be well-defined.

### Options

**Methods.** The DETECTANOMALY procedure clusters cases into peer groups based on the similarities of a set of input variables. An anomaly index is assigned to each case to reflect the unusualness of a case with respect to its peer group. All cases are sorted by the values of the anomaly index, and the top portion of the cases is identified as the set of anomalies. For each variable, an impact measure is assigned to each case that reflects the contribution of the variable to the deviation of the case from its peer group. For each case, the variables are sorted by the values of the variable impact measure, and the top portion of variables is identified as the set of reasons why the case is anomalous.

**Data File.** The DETECTANOMALY procedure assumes that the input data is a flat file in which each row represents a distinct case and each column represents a distinct variable. Moreover, it is assumed that all input variables are non-constant and that no case has missing values for all of the input variables.

**Missing Values.** The DETECTANOMALY procedure allows missing values. By default, missing values of continuous variables are substituted by their corresponding grand means, and missing categories of categorical variables are grouped and treated as a valid category. Moreover, an additional variable called the *Missing Proportion Variable*, which represents the proportion of missing variables in each case, is created. The processed variables are used to detect the anomalies in the data. You can turn off either of the options. If the first is turned off, cases with missing values are excluded from the analysis. In this situation, the second option is turned off automatically.

**ID Variable.** A variable that is the unique identifier of the cases in the data can optionally be specified in the ID keyword. If this keyword is not specified, the case sequence number of the active dataset is assumed to be the ID.

**Weights.** The DETECTANOMALY procedure ignores specification on the WEIGHT command.

**Output.** The DETECTANOMALY procedure displays an anomaly list in pivot table output, or offers an option for suppressing it. The procedure can also save the anomaly information to the active dataset as additional variables. Anomaly information can be grouped into three sets of variables: anomaly, peer, and reason. The anomaly set consists of the anomaly index of each case. The peer set consists of the peer group ID of each case, the size, and the percentage size of the peer group. The reason set consists of a number of reasons. Each reason consists of information such as the variable impact, the variable name for this reason, the value of the variable, and the corresponding norm value of the peer group.

### **Basic Specification**

The basic specification is the `DETECTANOMALY` command. By default, all variables in the active dataset are used in the procedure, with the dictionary setting of each variable in the dataset determining its measurement level.

### **Syntax Rules**

- All subcommands are optional.
- Only a single instance of each subcommand is allowed.
- An error occurs if an attribute or keyword is specified more than once within a subcommand.
- Parentheses, slashes, and equals signs shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.
- Empty subcommands are not honored.

### **Operations**

The `DETECTANOMALY` procedure begins by applying the missing value handling option and the create missing proportion variable option to the data.

Then the procedure groups cases into their peer groups based on the similarities of the processed variables. An anomaly index is assigned to each case to measure the overall deviation of the case from its peer group. All cases are sorted by the values of the anomaly index, and the top portion of the cases is identified as the anomaly list.

For each anomalous case, the variables are sorted by their corresponding variable impact values. The top variables, their values, and the corresponding norm values are presented as the reasons why a case is identified as an anomaly.

By default, the anomaly list is presented in a pivot table. Optionally, the anomaly information can be added to the active dataset as additional variable. The anomaly detection model may be written to an XML model file.

### **Limitations**

`WEIGHT` and `SPLIT FILE` settings are ignored with a warning by the `DETECTANOMALY` procedure.

## **Examples**

```
DETECTANOMALY  
  /VARIABLES CATEGORICAL=A B C SCALE=D E  
  /SAVE ANOMALY REASON.
```

- `DETECTANOMALY` treats variables *A*, *B*, and *C* as categorical variables and *D* and *E* as continuous variables.
- All of the processed variables are then used in the analysis to generate the anomaly list. Since no `ID` option is specified, the case number is used as the case identity variable. The size of the list is the number of cases with an anomaly index of at least 2.0 and is not more than 5% of the size of the working file. The anomaly list consists of the `ID` variable, the anomaly index, the peer group ID and size, and the reason. By default, there is one reason for each anomaly.

Each reason consists of information such as the variable impact measure, the variable name for this reason, the value of the variable, and the value of the corresponding peer group. The anomaly list is presented in the pivot table output.

- Since the keywords ANOMALY and REASON are specified in the SAVE subcommand, the additional variables for the anomaly index and the anomaly reason are added to the active dataset.

### **Specifying a Case ID Variable and Excepted Variables**

```
DETECTANOMALY
  /VARIABLES ID=CaseID EXCEPT=GeoID DemoID AddressID.
```

- DETECTANOMALY uses all variables in the active dataset in the analysis, except the variables *GeoID*, *DemoID*, and *AddressID*, which are excluded. Moreover, it treats the variable *CaseID* as the ID variable in the procedure. The processed variables are used in the analysis to generate the anomaly list.

## **VARIABLES Subcommand**

The VARIABLES subcommand specifies the variables to be used in the procedure. If the CATEGORICAL option or the SCALE option are specified, then variables that are listed in these options are used in the analysis. If neither the CATEGORICAL option nor the SCALE option is specified, then all variables in the active dataset are used, except the variable specified in the ID option and those in the EXCEPT option, if any. In the latter situation, the dictionary setting of each variable determines its measurement level. The procedure treats ordinal and nominal variables equivalently as categorical.

### **CATEGORICAL=varlist**

*List of categorical variables.* If this option is specified, at least one variable must be listed. Variables in the list can be numeric or string. They are treated as categorical variables and are used in the analysis. If duplicate variables are specified in the list, the duplicates are ignored. After the EXCEPT option (if specified) is applied, if there are variables also specified in either the continuous list or as an ID variable, an error is issued. TO and ALL keywords may be used.

### **SCALE=varlist**

*List of continuous variables.* If this option is specified, at least one variable must be listed. Variables in the list must be numeric and are used in the analysis. If duplicate variables are specified in the list, the duplicates are ignored. After the EXCEPT option (if specified) is applied, if there are variables also specified in either the categorical list or as an ID variable, an error is issued. TO and ALL keywords may be used.

### **ID=variable**

*Case ID variable.* If this option is specified, one variable must be listed. The variable can be numeric or string. It is used as a unique identifier for the cases in the data file and is not used in the analysis. If this option is not specified, the case number of the active dataset is used as the identifier variable. If the identifier variable is specified in either the categorical list or continuous list, an error is issued.

**EXCEPT=varlist**

*List of variables that are excluded from the analysis.* If this option is specified, at least one variable must be listed. Variables in the list are not used in the analysis, even if they are specified in the continuous or categorical lists. This option ignores duplicate variables and variables that are not specified on the continuous or categorical list. Specifying the ALL keyword causes an error. The TO keyword may be used.

This option can be useful if the categorical list or continuous list contains a large number of variables but there are a few variables that should be excluded.

## **HANDLEMISSING Subcommand**

The HANDLEMISSING subcommand specifies the methods of handling missing values in this procedure.

**APPLY=optionvalue**

*Apply missing value handling.* Valid option values are YES or NO. If YES, the missing values of continuous variables are substituted by their corresponding grand means, and missing categories of categorical variables are grouped and treated as a valid category. The processed variables are used in the analysis. If NO, cases with missing values are excluded from the analysis. The default value is NO.

**CREATEMISPROPVAR=optionvalue**

*Create an additional Missing Proportion Variable and use it in the analysis.* Valid option values are YES or NO. If YES, an additional variable called the *Missing Proportion Variable* that represents the proportion of missing variables in each case is created, and this variable is used in the analysis. If NO, the *Missing Proportion Variable* is not created. The default value is NO.

## **CRITERIA Subcommand**

The CRITERIA subcommand specifies settings for the DETECTANOMALY procedure.

**MINNUMPEERS=integer**

*Minimum number of peer groups.* The procedure will search for the best number of peer groups between the specified value and the value in the MAXNUMPEERS keyword. The specified value must be a positive integer less than or equal to the value in the MAXNUMPEERS keyword. When the specified value is equal to the value in the MAXNUMPEERS keyword, the procedure assumes a fixed number of peer groups. The default value is 1.

*Note:* Depending on the amount of variation in your data, there may be situations in which the number of peer groups that the data can support is less than the number specified in the MINNUMPEERS option. In such a situation, the procedure may produce a smaller number of peer groups.

<b>MAXNUMPEERS=integer</b>	<i>Maxim number of peer groups.</i> The procedure will search for the best number of peer groups between the value in the MINNUMPEERS keyword and the specified value. The specified value must be a positive integer greater than or equal to the value in the MINNUMPEERS keyword. When the specified value is equal to the value in the MINNUMPEERS keyword, the procedure assumes a fixed number of peer groups. The default value is 15.
<b>MLWEIGHT=number</b>	<i>An adjustment weight on the measurement level.</i> This parameter is used to balance the influences between continuous and categorical variables during the calculation of the indices. A large value increases the influence of a continuous variable. Specify a positive number. The default value is 6.
<b>NUMREASONS=integer</b>	<i>Number of reasons in the anomaly list.</i> A reason consists of information such as the variable impact measure, the variable name for this reason, the value of the variable, and the value of the corresponding peer group. Specify a non-negative integer less than or equal to the number of processed variables used in the analysis. The specified option value will be adjusted downward to the maximum number of variables used in the analysis if it is set larger than the number of variables. The default value is 1.
<b>PCTANOMALOUSCASES=number</b>	<i>Percentage of cases considered as anomalies and included in the anomaly list.</i> Specify a non-negative number less than or equal to 100. The default value is 5.
<b>NUMANOMALOUSCASES=integer</b>	<i>Number of cases considered as anomalies and included in the anomaly list.</i> Specify a non-negative integer less than or equal to the total number of cases in the active dataset and used in the analysis. If this option is specified, an option value must be listed. The specified option value will be adjusted downward to the maximum available if it is set larger than the number of cases used in the analysis. This option, if specified, overrides the PCTANOMALOUSCASES option.
<b>ANOMALYCUTPOINT=number</b>	<i>Cut point of the anomaly index to determine whether a case is considered as an anomaly.</i> Specify a non-negative number. A case is considered anomalous if its anomaly index value is larger than or equal to the specified cut point. This option can be used together with the PCTANOMALOUSCASES and NUMANOMALOUSCASES options. For example, if NUMANOMALOUSCASES=50 and ANOMALYCUTPOINT=2 are specified, the anomaly list will consist of at most 50 cases each with an anomaly index value larger than or equal to 2. The default value is 2. If NONE is specified, the option is suppressed and no cut point is set.

## **SAVE Subcommand**

The **SAVE** subcommand specifies the additional variables to save to the active dataset.

- One or more keywords should be specified, each followed by an optional variable name or rootname in parentheses.
- The variable name or the rootname, if specified, must be a valid variable name.
- If no variable name or rootname is specified, a default varname or rootname is used. If the default varname or rootname is used and it conflicts with that of an existing variable, a suffix is added to make the name unique.

- The values of the additional variables are assigned to all cases included in the analysis, even if the cases are not in the anomaly list.
- This subcommand is not affected by the specifications on the PCTANOMALOUSCASES, NUMANOMALOUSCASES, or ANOMALYCUTPOINT keywords in the CRITERIA subcommand.

**ANOMALY(varname)** *The anomaly index.* If an optional varname is not specified, the default varname is *AnomalyIndex*, which is the anomaly index. If an optional varname is specified, the specified varname is used to replace the default varname. For example, if **ANOMALY(MyAnomaly)** is specified, the variable name will be *MyAnomaly*.

**PEERID(varname)** *Peer group ID.* If an optional varname is not specified, the default varname *PeerId* is used. If an optional varname is specified, the specified name is used.

**PEERSIZE(varname)** *Peer group size.* If an optional varname is not specified, the default variable name *PeerSize* is used. If an optional varname is specified, the specified name is used.

**PEERPCTSIZE(varname)** *Peer group size in percentage.* If an optional varname is not specified, the default varname *PeerPctSize* is used. If an optional varname is specified, the specified name is used.

**REASONVAR(rootname)** *The variable associated with a reason.* The number of REASONVAR variables created depends on the number of reasons specified on the CRITERIA subcommand NUMREASONS option. If an optional rootname is not specified, the default rootname *ReasonVar* is used to automatically generate one or more varnames, *ReasonVar<sub>k</sub>*, where *k* is the *k*th reason. If an optional rootname is specified, the specified name is used. If NUMREASONS=0 is specified, this option is ignored and a warning is issued.

**REASONMEASURE(root-name)**

*The variable impact measure associated with a reason.* The number of REASONMEASURE variables created depends on the number of reasons specified on the CRITERIA subcommand NUMREASONS option. If an optional rootname is not specified, the default rootname *ReasonMeasure* is used to automatically generate one or more varnames, *ReasonMeasure<sub>k</sub>*, where *k* is the *k*th reason. If an optional rootname is specified, the specified name is used. If NUMREASONS=0 is specified, this option is ignored and a warning is issued.

**REASONVALUE(root-name)**

*The variable value associated with a reason.* The number of REASONVALUE variables created depends on the number of reasons specified on the CRITERIA subcommand NUMREASONS option. If an optional rootname is not specified, the default rootname *ReasonValue* is used to automatically generate one or more varnames, *ReasonValue<sub>k</sub>*, where *k* is the *k*th reason. If an optional rootname is specified, the specified name is used. If NUMREASONS=0 is specified, this option is ignored and a warning is issued.

**REASONNORM(root-name)**

*The norm value associated with a reason.* The number of REASONNORM variables created depends on the number of reasons specified on the CRITERIA subcommand NUMREASONS option. If an optional rootname is not specified, the default rootname *ReasonNorm* is used to automatically generate one or more varnames, *ReasonNorm<sub>k</sub>*, where *k* is the *k*th reason. If an optional rootname is specified, the



specified name is used. If NUMREASONS=0 is specified, this option is ignored and a warning is issued.

## **OUTFILE Subcommand**

The OUTFILE subcommand directs the DETECTANOMALY procedure to write its model to the specified filename as XML.

**MODEL=filespec** *File specification to which the model is written.*

## **PRINT Subcommand**

The PRINT subcommand controls the display of the output results.

- If the PRINT subcommand is not specified, the default output is the anomaly list. If the PRINT subcommand is specified, DETECTANOMALY displays output only for the keywords that are specified.

<b>CPS</b>	<i>Display a case processing summary.</i> The case processing summary displays the counts and count percentages for all cases in the active dataset, the cases included and excluded in the analysis, and the cases in each peer.
<b>ANOMALYLIST</b>	<i>Display the anomaly index list, the anomaly peer ID list, and the anomaly reason list.</i> The anomaly index list displays the case number and its corresponding anomaly index value. The anomaly peer ID list displays the case number, its corresponding peer group ID, peer size, and size in percent. The anomaly reason list displays the case number, the reason variable, the variable impact value, the value of the variable, and the norm of the variable for each reason. All tables are sorted by anomaly index in descending order. Moreover, the IDs of the cases are displayed if the case identifier variable is specified in the ID option of the VARIABLES subcommand. This is the default output.
<b>NORMS</b>	<i>Display the continuous variable norms table if any continuous variable is used in the analysis, and display the categorical variable norms table if any categorical variable is used in the analysis.</i> In the continuous variable norms table, the mean and standard deviation of each continuous variable for each peer group is displayed. In the categorical variable norms table, the mode (most popular category), its frequency, and frequency percent of each categorical variable for each peer group is displayed. The mean of a continuous variable and the mode of a categorical variable are used as the norm values in the analysis.
<b>ANOMALYSUMMARY</b>	<i>Display the anomaly index summary.</i> The anomaly index summary displays descriptive statistics for the anomaly index of the cases identified as the most unusual.

---

*DETECTANOMALY***REASONSUMMARY**

*Display the reason summary table for each reason.* For each reason, the table displays the frequency and frequency percent of each variable's occurrence as a reason. It also reports the descriptive statistics of the impact of each variable. If `NUMREASONS=0` is specified on the `CRITERIA` subcommand, this option is ignored and a warning is issued.

**NONE**

*Suppress all displayed output except the notes table and any warnings.* If `NONE` is specified with one or more other keywords, the other keywords override `NONE`.

# DISCRIMINANT

DISCRIMINANT is available in the Statistics Base option.

```
DISCRIMINANT GROUPS=varname(min,max) /VARIABLES=varlist

[/SELECT=varname(value)]

[/ANALYSIS=varlist[(level)] [varlist...]]

[/OUTFILE MODEL('file')]

[/METHOD={DIRECT**}] [/TOLERANCE={0.001}]
    {WILKS      }           { n      }
    {MAHAL      }
    {MAXMINF    }
    {MINRESID   }
    {RAO        }

[/MAXSTEPS={n}]

[/FIN={3.84**}] [/FOUT={2.71**}] [/PIN={n}]
    { n      }           { n      }

[/POUT={n}] [/VIN={0**}]
    { n      }

[/FUNCTIONS={g-1,100.0,1.0**}] [/PRIORS={EQUAL** }]}
    {n1 , n2 , n3 }           {SIZE  }
                                {value list}

[/SAVE=[CLASS[=varname]] [PROBS[=rootname]]
    [SCORES[=rootname]]]

[/ANALYSIS=...]

[/MISSING={EXCLUDE**}]
    {INCLUDE  }

[/MATRIX=[OUT({*
    {'savfile'|'dataset'}})] [IN({*
    {'savfile'|'dataset'}})]

[/HISTORY={STEP**} ]
    {NONE  }

[/ROTATE={NONE** }]}
    {COEFF }
    {STRUCTURE}

[/CLASSIFY={NONMISSING } {POOLED } [MEANSUB]]
    {UNSELECTED } {SEPARATE}
    {UNCLASSIFIED}

[/STATISTICS=[MEAN] [COV ] [FPAIR] [RAW ] [STDDEV]
    [GCOV] [UNIVF] [COEFF] [CORR] [TCOV ]
    [BOXM] [TABLE] [CROSSVALID]
    [ALL]]

[/PLOT=[MAP] [SEPARATE] [COMBINED] [CASES[(n)]] [ALL]]
```

**\*\*Default if subcommand or keyword is omitted.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 17.0

- Added support for SET THREADS and SET MCACHE.

### **Example**

```
DISCRIMINANT GROUPS=OUTCOME (1,4)  
/VARIABLES=V1 TO V7.
```

## **Overview**

DISCRIMINANT performs linear discriminant analysis for two or more groups. The goal of discriminant analysis is to classify cases into one of several mutually exclusive groups based on their values for a set of predictor variables. In the analysis phase, a classification rule is developed using cases for which group membership is known. In the classification phase, the rule is used to classify cases for which group membership is not known. The grouping variable must be categorical, and the independent (predictor) variables must be interval or dichotomous, since they will be used in a regression-type equation.

### **Options**

**Variable Selection Method.** In addition to the direct-entry method, you can specify any of several stepwise methods for entering variables into the discriminant analysis using the METHOD subcommand. You can set the values for the statistical criteria used to enter variables into the equation using the TOLERANCE, FIN, PIN, FOUT, POUT, and VIN subcommands, and you can specify inclusion levels on the ANALYSIS subcommand. You can also specify the maximum number of steps in a stepwise analysis using the MAXSTEPS subcommand.

**Case Selection.** You can select a subset of cases for the analysis phase using the SELECT subcommand.

**Prior Probabilities.** You can specify prior probabilities for membership in a group using the PRIORS subcommand. Prior probabilities are used in classifying cases.

**New Variables.** You can add new variables to the active dataset containing the predicted group membership, the probability of membership in each group, and discriminant function scores using the SAVE subcommand.

**Classification Options.** With the CLASSIFY subcommand, you can classify only those cases that were not selected for inclusion in the discriminant analysis, or only those cases whose value for the grouping variable was missing or fell outside the range analyzed. In addition, you can classify cases based on the separate-group covariance matrices of the functions instead of the pooled within-groups covariance matrix.

**Statistical Display.** You can request any of a variety of statistics on the STATISTICS subcommand. You can rotate the pattern or structure matrices using the ROTATE subcommand. You can compare actual with predicted group membership using a classification results table requested with the STATISTICS subcommand or compare any of several types of plots or histograms using the PLOT subcommand.

### **Basic Specification**

The basic specification requires two subcommands:

- `GROUPS` specifies the variable used to group cases.
- `VARIABLES` specifies the predictor variables.

By default, `DISCRIMINANT` enters all variables simultaneously into the discriminant equation (the `DIRECT` method), provided that they are not so highly correlated that multicollinearity problems arise. Default output includes analysis case processing summary, valid numbers of cases in group statistics, variables failing tolerance test, a summary of canonical discriminant functions, standardized canonical discriminant function coefficients, a structure matrix showing pooled within-groups correlations between the discriminant functions and the predictor variables, and functions at group centroids.

### **Subcommand Order**

- The `GROUPS`, `VARIABLES`, and `SELECT` subcommands must precede all other subcommands and may be entered in any order.
- The analysis block follows, which may include `ANALYSIS`, `METHOD`, `TOLERANCE`, `MAXSTEPS`, `FIN`, `FOUT`, `PIN`, `POUT`, `VIN`, `FUNCTIONS`, `PRIORS`, `SAVE`, and `OUTFILE`. Each analysis block performs a single analysis. To do multiple analyses, specify multiple analysis blocks.
- The keyword `ANALYSIS` is optional for the first analysis block. Each new analysis block must begin with an `ANALYSIS` subcommand. Remaining subcommands in the block may be used in any order and apply only to the analysis defined within the same block.
- No analysis block subcommands can be specified after any of the global subcommands, which apply to all analysis blocks. The global subcommands are `MISSING`, `MATRIX`, `HISTORY`, `ROTATE`, `CLASSIFY`, `STATISTICS`, and `PLOT`. If an analysis block subcommand appears after a global subcommand, the program displays a warning and ignores it.

### **Syntax Rules**

- Only one `GROUPS`, one `SELECT`, and one `VARIABLES` subcommand can be specified per `DISCRIMINANT` command.

### **Operations**

- `DISCRIMINANT` first estimates one or more discriminant functions that best distinguish among the groups.
- Using these functions, `DISCRIMINANT` then classifies cases into groups (if classification output is requested).
- If more than one analysis block is specified, the above steps are repeated for each block.
- This procedure uses the multithreaded options specified by `SET THREADS` and `SET MCACHE`.

### **Limitations**

- Pairwise deletion of missing data is not available.

## Examples

```
DISCRIMINANT
  /GROUPS=default(0 1)
  /VARIABLES=employ address debtinc creddebt
  /SELECT=validate(1)
  /ANALYSIS ALL
  /SAVE=CLASS PROBS
  /PRIORS EQUAL
  /STATISTICS=MEANS STDDEV UNIVF BOXM COEFF RAW FPAIR TABLE CROSSVALID CORR
  /CLASSIFY=NONMISSING POOLED .
```

- The procedure builds a model for discriminating between the values 0 and 1 on the variable *default* using the variables *employ*, *address*, *debtinc*, and *creddebt*.
- The `SELECT` subcommand specifies that only cases with a value of 1 on the variable *validate* are used to build the model.
- The `SAVE` subcommand requests predicted group membership and predicted probabilities for each group to be saved to the active dataset.
- The `STATISTICS` subcommand requests Box's *M* test, classification function coefficients, unstandardized discriminant functions, pairwise *F* ratios, classification results and crossvalidated results, and a pooled within-groups correlation matrix in addition to the default output.
- All other options are set to their default values.

### Stepwise variable selection

```
DISCRIMINANT
  /GROUPS=custcat(1 4)
  /VARIABLES=age marital address income ed retire employ gender reside
  /ANALYSIS ALL
  /METHOD=WILKS
  /FIN= 3.84
  /FOUT= 2.71
  /PRIORS EQUAL
  /HISTORY
  /STATISTICS= TABLE
  /PLOT=MAP
  /CLASSIFY=NONMISSING POOLED .
```

- The procedure builds a model for discriminating between the values 1, 2, 3, and 4 on the variable *custcat* using the variables *age*, *marital*, *address*, *income*, *ed*, *retire*, *employ*, *gender*, and *reside*.
- The `METHOD`, `FIN`, and `FOUT` subcommands specify stepwise variable selection using Wilks' lambda as the statistic, an *F* value of 3.84 as the criteria for determining entry into the model, and an *F* value of 2.71 for determining removal from the model.
- The `STATISTICS` subcommand requests a classification table.
- The `PLOT` subcommand requests a territorial map.
- All other options are set to their default values.

## ***GROUPS Subcommand***

GROUPS specifies the name of the grouping variable, which defines the categories or groups, and a range of categories.

- GROUPS is required and can be specified only once.
- The specification consists of a variable name followed by a range of values in parentheses.
- Only one grouping variable may be specified; its values must be integers. To use a string variable as the grouping variable, first use AUTORECODE to convert the string values to integers and then specify the recoded variable as the grouping variable.
- Empty groups are ignored and do not affect calculations. For example, if there are no cases in group 2, the value range (1, 5) will define only four groups.
- Cases with values outside the value range or missing are ignored during the analysis phase but are classified during the classification phase.

## ***VARIABLES Subcommand***

VARIABLES identifies the predictor variables, which are used to classify cases into the groups defined on the GROUPS subcommand. The list of variables follows the usual conventions for variable lists.

- VARIABLES is required and can be specified only once. Use the ANALYSIS subcommand to obtain multiple analyses.
- Only numeric variables can be used.
- Variables should be suitable for use in a regression-type equation, either measured at the interval level or dichotomous.

## ***SELECT Subcommand***

SELECT limits cases used in the analysis phase to those with a specified value for any one variable.

- Only one SELECT subcommand is allowed. It can follow the GROUPS and VARIABLES subcommands but must precede all other subcommands.
- The specification is a variable name and a single integer value in parentheses. Multiple variables or values are not permitted.
- The selection variable does not have to be specified on the VARIABLES subcommand.
- Only cases with the specified value for the selection variable are used in the analysis phase.
- All cases, whether selected or not, are classified by default. Use CLASSIFY=UNSELECTED to classify only the unselected cases.
- When SELECT is used, classification statistics are reported separately for selected and unselected cases, unless CLASSIFY=UNSELECTED is used to restrict classification.

### ***Example***

```
DISCRIMINANT GROUPS=APPROVAL(1,5)  
/VARS=Q1 TO Q10
```

*DISCRIMINANT*

```
/SELECT=COMPLETE(1)
/CLASSIFY=UNSELECTED.
```

- Using only cases with the value 1 for the variable *COMPLETE*, *DISCRIMINANT* estimates a function of *Q1* to *Q10* that discriminates between the categories 1 to 5 of the grouping variable *APPROVAL*.
- Because *CLASSIFY=UNSELECTED* is specified, the discriminant function will be used to classify only the unselected cases (cases for which *COMPLETE* does not equal 1).

***ANALYSIS Subcommand***

*ANALYSIS* is used to request several different discriminant analyses using the same grouping variable, or to control the order in which variables are entered into a stepwise analysis.

- *ANALYSIS* is optional for the first analysis block. By default, all variables specified on the *VARIABLES* subcommand are included in the analysis.
- The variables named on *ANALYSIS* must first be specified on the *VARIABLES* subcommand.
- The keyword *ALL* includes all variables on the *VARIABLES* subcommand.
- If the keyword *TO* is used to specify a list of variables on an *ANALYSIS* subcommand, it refers to the order of variables on the *VARIABLES* subcommand, which is not necessarily the order of variables in the active dataset.

***Example***

```
DISCRIMINANT GROUPS=SUCCESS(0,1)
/VARIABLES=V10 TO V15, AGE, V5
/ANALYSIS=V15 TO V5
/ANALYSIS=ALL.
```

- The first analysis will use the variables *V15*, *AGE*, and *V5* to discriminate between cases where *SUCCESS* equals 0 and *SUCCESS* equals 1.
- The second analysis will use all variables named on the *VARIABLES* subcommand.

***Inclusion Levels***

When you specify a stepwise method on the *METHOD* subcommand (any method other than the default direct-entry method), you can control the order in which variables are considered for entry or removal by specifying inclusion levels on the *ANALYSIS* subcommand. By default, all variables in the analysis are entered according to the criterion requested on the *METHOD* subcommand.

- An **inclusion level** is an integer between 0 and 99, specified in parentheses after a variable or list of variables on the *ANALYSIS* subcommand.
- The default inclusion level is 1.
- Variables with higher inclusion levels are considered for entry before variables with lower inclusion levels.
- Variables with even inclusion levels are entered as a group.



- Variables with odd inclusion levels are entered individually, according to the stepwise method specified on the `METHOD` subcommand.
- Only variables with an inclusion level of 1 are considered for removal. To make a variable with a higher inclusion level eligible for removal, name it twice on the `ANALYSIS` subcommand, first specifying the desired inclusion level and then an inclusion level of 1.
- Variables with an inclusion level of 0 are never entered. However, the statistical criterion for entry is computed and displayed.
- Variables that fail the tolerance criterion are not entered regardless of their inclusion level.

The following are some common methods of entering variables and the inclusion levels that could be used to achieve them. These examples assume that one of the stepwise methods is specified on the `METHOD` subcommand (otherwise, inclusion levels have no effect).

**Direct.** `ANALYSIS=ALL(2)` forces all variables into the equation. (This is the default and can be requested with `METHOD=DIRECT` or simply by omitting the `METHOD` subcommand.)

**Stepwise.** `ANALYSIS=ALL(1)` yields a stepwise solution in which variables are entered and removed in stepwise fashion. (This is the default when anything other than `DIRECT` is specified on the `METHOD` subcommand.)

**Forward.** `ANALYSIS=ALL(3)` enters variables into the equation stepwise but does not remove variables.

**Backward.** `ANALYSIS=ALL(2) ALL(1)` forces all variables into the equation and then allows them to be removed stepwise if they satisfy the criterion for removal.

#### ***Inclusion Levels Used With a Stepwise Method***

```
DISCRIMINANT GROUPS=SUCCESS(0,1)
/VARIABLES=A, B, C, D, E
/ANALYSIS=A TO C (2) D, E (1)
/METHOD=WILKS.
```

- *A*, *B*, and *C* are entered into the analysis first, assuming that they pass the tolerance criterion. Since their inclusion level is even, they are entered together.
- *D* and *E* are then entered stepwise. The one that minimizes the overall value of Wilks' lambda is entered first.
- After entering *D* and *E*, the program checks whether the partial *F* for either one justifies removal from the equation (see the `FOUT` and `POUT` subcommands).

#### ***Inclusion Levels Without a Stepwise Method***

```
DISCRIMINANT GROUPS=SUCCESS(0,1)
/VARIABLES=A, B, C, D, E
/ANALYSIS=A TO C (2) D, E (1).
```

- Since no stepwise method is specified, inclusion levels have no effect and all variables are entered into the model at once.

## **METHOD Subcommand**

METHOD is used to select a method for entering variables into an analysis.

- A variable will never be entered into the analysis if it does not pass the tolerance criterion specified on the TOLERANCE subcommand (or the default).
- A METHOD subcommand applies only to the *preceding* ANALYSIS subcommand, or to an analysis using all predictor variables if no ANALYSIS subcommand has been specified before it.
- If more than one METHOD subcommand is specified within one analysis block, the last is used.

Any one of the following methods can be specified on the METHOD subcommand:

<b>DIRECT</b>	<i>All variables passing the tolerance criteria are entered simultaneously. This is the default method.</i>
<b>WILKS</b>	<i>At each step, the variable that minimizes the overall Wilks' lambda is entered.</i>
<b>MAHAL</b>	<i>At each step, the variable that maximizes the Mahalanobis distance between the two closest groups is entered.</i>
<b>MAXMINF</b>	<i>At each step, the variable that maximizes the smallest F ratio between pairs of groups is entered.</i>
<b>MINRESID</b>	<i>At each step, the variable that minimizes the sum of the unexplained variation for all pairs of groups is entered.</i>
<b>RAO</b>	<i>At each step, the variable that produces the largest increase in Rao's V is entered.</i>

## **OUTFILE Subcommand**

Exports model information to the specified file in XML (PMML) format. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#)

- The minimum specification is the keyword MODEL and a file name enclosed in parentheses.
- The OUTFILE subcommand cannot be used if split file processing is on (SPLIT FILE command).

## **TOLERANCE Subcommand**

TOLERANCE specifies the minimum tolerance a variable can have and still be entered into the analysis. The tolerance of a variable that is a candidate for inclusion in the analysis is the proportion of its within-groups variance not accounted for by other variables in the analysis. A variable with very low tolerance is nearly a linear function of the other variables; its inclusion in the analysis would make the calculations unstable.

- The default tolerance is 0.001.
- You can specify any decimal value between 0 and 1 as the minimum tolerance.

## ***PIN and POUT Subcommands***

`PIN` specifies the minimum probability of  $F$  that a variable can have to enter the analysis and `POUT` specifies the maximum probability of  $F$  that a variable can have and not be removed from the model.

- `PIN` and `POUT` take precedence over `FIN` and `FOUT`. That is, if all are specified, `PIN` and `POUT` values are used.
- If `PIN` and `POUT` are omitted, `FIN` and `FOUT` are used by default.
- You can set `PIN` and `POUT` to any decimal value between 0 and 1. However, `POUT` should be greater than `PIN` if `PIN` is also specified.
- `PIN` and `POUT` apply only to the stepwise methods and are ignored if the `METHOD` subcommand is omitted or if `DIRECT` is specified on `METHOD`.

## ***FIN and FOUT Subcommands***

`FIN` specifies the minimum partial  $F$  value that a variable must have to enter the analysis. As additional variables are entered into the analysis, the partial  $F$  for variables already in the equation changes. `FOUT` specifies the smallest partial  $F$  that a variable can have and not be removed from the model.

- `PIN` and `POUT` take precedence over `FIN` and `FOUT`. That is, if all are specified, `PIN` and `POUT` values are used.
- If `PIN` and `POUT` are omitted, `FIN` and `FOUT` are used by default. If `FOUT` is specified but `FIN` is omitted, the default value for `FIN` is 3.84. If `FIN` is specified, the default value for `FOUT` is 2.71.
- You can set `FIN` and `FOUT` to any non-negative number. However, `FOUT` should be less than `FIN` if `FIN` is also specified.
- `FIN` and `FOUT` apply only to the stepwise methods and are ignored if the `METHOD` subcommand is omitted or if `DIRECT` is specified on `METHOD`.

## ***VIN Subcommand***

`VIN` specifies the minimum Rao's  $V$  that a variable must have to enter the analysis. When you use `METHOD=RAO`, variables satisfying one of the other criteria for entering the equation may actually cause a decrease in Rao's  $V$  for the equation. The default `VIN` prevents this but does not prevent the addition of variables that provide no additional separation between groups.

- You can specify any value for `VIN`. The default is 0.
- `VIN` should be used only when you have specified `METHOD=RAO`. Otherwise, it is ignored.

## ***MAXSTEPS Subcommand***

`MAXSTEPS` is used to decrease the maximum number of steps allowed. By default, the maximum number of steps allowed in a stepwise analysis is the number of variables with inclusion levels greater than 1 plus twice the number of variables with inclusion levels equal to 1. This is the

maximum number of steps possible without producing a loop in which a variable is repeatedly cycled in and out.

- `MAXSTEPS` applies only to the stepwise methods (all except `DIRECT`).
- `MAXSTEPS` applies only to the preceding `METHOD` subcommand.
- The format is `MAX=n`, where  $n$  is the maximum number of steps desired.
- If multiple `MAXSTEPS` subcommands are specified, the last is used.

## ***FUNCTIONS Subcommand***

By default, `DISCRIMINANT` computes all possible functions. This is either the number of groups minus 1 or the number of predictor variables, whichever is less. Use `FUNCTIONS` to set more restrictive criteria for the extraction of functions.

`FUNCTIONS` has three parameters:

- *Maximum number of functions.* The default is the number of groups minus 1 or the number of predictor variables, whichever is less.
  - *Cumulative percentage of the sum of the eigenvalues.* The default is 100.
  - *Significance level of function.* The default is 1.0.
- The parameters must always be specified in sequential order ( $n_1, n_2, n_3$ ). To specify  $n_2$ , you must explicitly specify the default for  $n_1$ . Similarly, to specify  $n_3$ , you must specify the defaults for  $n_1$  and  $n_2$ .
  - If more than one restriction is specified, the program stops extracting functions when any one of the restrictions is met.
  - When multiple `FUNCTIONS` subcommands are specified, the program uses the last; however, if  $n_2$  or  $n_3$  are omitted on the last `FUNCTIONS` subcommand, the corresponding specifications on the previous `FUNCTIONS` subcommands will remain in effect.

### ***Example***

```
DISCRIMINANT GROUPS=CLASS(1,5)
/VARIABLES = SCORE1 TO SCORE20
/FUNCTIONS=4,100,.80.
```

- The first two parameters on the `FUNCTIONS` subcommand are defaults: the default for  $n_1$  is 4 (the number of groups minus 1), and the default for  $n_2$  is 100.
- The third parameter tells `DISCRIMINANT` to use fewer than four discriminant functions if the significance level of a function is greater than 0.80.

## ***PRIORS Subcommand***

By default, `DISCRIMINANT` assumes equal prior probabilities for groups when classifying cases. You can provide different prior probabilities with the `PRIORS` subcommand.

- Prior probabilities are used only during classification.

- If you provide unequal prior probabilities, DISCRIMINANT adjusts the classification coefficients to reflect this.
- If adjacent groups have the same prior probability, you can use the notation  $n*c$  on the value list to indicate that  $n$  adjacent groups have the same prior probability  $c$ .
- You can specify a prior probability of 0. No cases are classified into such a group.
- If the sum of the prior probabilities is not 1, the program rescales the probabilities to sum to 1 and issues a warning.

<b>EQUAL</b>	<i>Equal prior probabilities.</i> This is the default.
<b>SIZE</b>	<i>Proportion of the cases analyzed that fall into each group.</i> If 50% of the cases included in the analysis fall into the first group, 25% in the second, and 25% in the third, the prior probabilities are 0.5, 0.25, and 0.25, respectively. Group size is determined after cases with missing values for the predictor variables are deleted.
<b>Value list</b>	<i>User-specified prior probabilities.</i> The list of probabilities must sum to 1.0. The number of prior probabilities named or implied must equal the number of groups.

### Example

```
DISCRIMINANT GROUPS=TYPE(1,5)
/VARIABLES=A TO H
/PRIORS = 4*.15, .4.
```

- The PRIORS subcommand establishes prior probabilities of 0.15 for the first four groups and 0.4 for the fifth group.

## SAVE Subcommand

SAVE allows you to save casewise information as new variables in the active dataset.

- SAVE applies only to the current analysis block. To save casewise results from more than one analysis, specify a SAVE subcommand in each analysis block.
- You can specify a variable name for CLASS and rootnames for SCORES and PROBS to obtain descriptive names for the new variables.
- If you do not specify a variable name for CLASS, the program forms variable names using the formula  $DSC\_m$ , where  $m$  increments to distinguish group membership variables saved on different SAVE subcommands for different analysis blocks.
- If you do not specify a rootname for SCORES or PROBS, the program forms new variable names using the formula  $DSCn\_m$ , where  $m$  increments to create unique rootnames and  $n$  increments to create unique variable names. For example, the first set of default names assigned to discriminant scores or probabilities are  $DSC1\_1$ ,  $DSC2\_1$ ,  $DSC3\_1$ , and so on. The next set of default names assigned will be  $DSC1\_2$ ,  $DSC2\_2$ ,  $DSC3\_2$ , and so on, regardless of whether discriminant scores or probabilities are being saved or whether they are saved by the same SAVE subcommand.
- The keywords CLASS, SCORES, and PROBS can be used in any order, but the new variables are always added to the end of the active dataset in the following order: first the predicted group, then the discriminant scores, and finally probabilities of group membership.

## DISCRIMINANT

- Appropriate variable labels are automatically generated. The labels describe whether the variables contain predictor group membership, discriminant scores, or probabilities, and for which analysis they are generated.
- The CLASS variable will use the value labels (if any) from the grouping variable specified for the analysis.
- When SAVE is specified with any keyword, DISCRIMINANT displays a classification processing summary table and a prior probabilities for groups table.
- You cannot use the SAVE subcommand if you are replacing the active dataset with matrix materials (see Matrix Output on p. 645).

<b>CLASS [(varname)]</b>	<i>Predicted group membership.</i>
<b>SCORES [(rootname)]</b>	<i>Discriminant scores. One score is saved for each discriminant function derived. If a rootname is specified, DISCRIMINANT will append a sequential number to the name to form new variable names for the discriminant scores.</i>
<b>PROBS [(rootname)]</b>	<i>For each case, the probabilities of membership in each group. As many variables are added to each case as there are groups. If a rootname is specified, DISCRIMINANT will append a sequential number to the name to form new variable names.</i>

**Example**

```
DISCRIMINANT GROUPS=WORLD(1,3)
/VARIABLES=FOOD TO FSALES
/SAVE CLASS=PRDCLASS SCORES=SCORE PROBS=PRB
/ANALYSIS=FOOD SERVICE COOK MANAGER FSALES
/SAVE CLASS SCORES PROBS.
```

- Two analyses are specified. The first uses all variables named on the VARIABLES subcommand and the second narrows down to five variables. For each analysis, a SAVE subcommand is specified.
- For each analysis, DISCRIMINANT displays a classification processing summary table and a prior probabilities for groups table.
- On the first SAVE subcommand, a variable name and two rootnames are provided. With three groups, the following variables are added to each case:

<b>Name</b>	<b>Variable label</b>	<b>Description</b>
PRDCLASS	Predicted group for analysis 1	Predicted group membership
SCORE1	Function 1 for analysis 1	Discriminant score for function 1
SCORE2	Function 2 for analysis 1	Discriminant score for function 2
PRB1	Probability 1 for analysis 1	Probability of being in group 1

Name	Variable label	Description
PRB2	Probability 2 for analysis 1	Probability of being in group 2
PRB3	Probability 3 for analysis 1	Probability of being in group 3

- Since no variable name or rootnames are provided on the second `SAVE` subcommand, `DISCRIMINANT` uses default names. Note that *m* serves only to distinguish variables saved as a set and does not correspond to the sequential number of an analysis. To find out what information a new variable holds, read the variable label, as shown in the following table:

Name	Variable label	Description
DSC_1	Predicted group for analysis 2	Predicted group membership
DSC1_1	Function 1 for analysis 2	Discriminant score for function 1
DSC2_1	Function 2 for analysis 2	Discriminant score for function 2
DSC1_2	Probability 1 for analysis 2	Probability of being in group 1
DSC2_2	Probability 2 for analysis 2	Probability of being in group 2
DSC3_2	Probability 3 for analysis 2	Probability of being in group 3

## STATISTICS Subcommand

By default, `DISCRIMINANT` produces the following statistics for each analysis: analysis case processing summary, valid numbers of cases in group statistics, variables failing tolerance test, a summary of canonical discriminant functions, standardized canonical discriminant function coefficients, a structure matrix showing pooled within-groups correlations between the discriminant functions and the predictor variables, and functions at group centroids.

- *Group statistics.* Only valid number of cases is reported.
- *Summary of canonical discriminant functions.* Displayed in two tables: an eigenvalues table with percentage of variance, cumulative percentage of variance, and canonical correlations and a Wilks' lambda table with Wilks' lambda, chi-square, degrees of freedom, and significance.
- *Stepwise statistics.* Wilks' lambda, equivalent *F*, degrees of freedom, significance of *F* and number of variables are reported for each step. Tolerance, *F*-to-remove, and the value of the statistic used for variable selection are reported for each variable in the equation. Tolerance, minimum tolerance, *F*-to-enter, and the value of the statistic used for variable selection are reported for each variable not in the equation. (These statistics can be suppressed with `HISTORY=NONE`.)
- *Final statistics.* Standardized canonical discriminant function coefficients, the structure matrix of discriminant functions and all variables named in the analysis (whether they were entered into the equation or not), and functions evaluated at group means are reported following the last step.

In addition, you can request optional statistics on the `STATISTICS` subcommand. `STATISTICS` can be specified by itself or with one or more keywords.

- `STATISTICS` without keywords displays `MEAN`, `STDDEV`, and `UNIVF`. If you include a keyword or keywords on `STATISTICS`, only the statistics you request are displayed.

<b>MEAN</b>	<i>Means.</i> Total and group means for all variables named on the <code>ANALYSIS</code> subcommand are displayed.
<b>STDDEV</b>	<i>Standard deviations.</i> Total and group standard deviations for all variables named on the <code>ANALYSIS</code> subcommand are displayed.
<b>UNIVF</b>	<i>Univariate F ratios.</i> The analysis-of-variance <i>F</i> statistic for equality of group means for each predictor variable is displayed. This is a one-way analysis-of-variance test for equality of group means on a single discriminating variable.
<b>COV</b>	<i>Pooled within-groups covariance matrix.</i>
<b>CORR</b>	<i>Pooled within-groups correlation matrix.</i>
<b>FPAIR</b>	<i>Matrix of pairwise F ratios.</i> The <i>F</i> ratio for each pair of groups is displayed. This <i>F</i> is the significance test for the Mahalanobis distance between groups. This statistic is available only with stepwise methods.
<b>BOXM</b>	<i>Box's M test.</i> This is a test for equality of group covariance matrices.
<b>GCOV</b>	<i>Group covariance matrices.</i>
<b>TCOV</b>	<i>Total covariance matrix.</i>
<b>RAW</b>	<i>Unstandardized canonical discriminant functions.</i>
<b>COEFF</b>	<i>Classification function coefficients.</i> Although <code>DISCRIMINANT</code> does not directly use these coefficients to classify cases, you can use them to classify other samples (see the <code>CLASSIFY</code> subcommand).
<b>TABLE</b>	<i>Classification results.</i> If both selected and unselected cases are classified, the results are reported separately. To obtain cross-validated results for selected cases, specify <code>CROSSVALID</code> .
<b>CROSSVALID</b>	<i>Cross-validated classification results.</i> The cross-validation is done by treating $n-1$ out of $n$ observations as the training dataset to determine the discrimination rule and using the rule to classify the one observation left out. The results are displayed only for selected cases.
<b>ALL</b>	<i>All optional statistics.</i>

## ***ROTATE Subcommand***

The coefficient and correlation matrices can be rotated to facilitate interpretation of results. To control varimax rotation, use the `ROTATE` subcommand.

- Neither `COEFF` nor `STRUCTURE` affects the classification of cases.

<b>COEFF</b>	<i>Rotate pattern matrix.</i> <code>DISCRIMINANT</code> displays a varimax transformation matrix, a rotated standardized canonical discriminant function coefficients table, and a correlations between variables and rotated functions table.
<b>STRUCTURE</b>	<i>Rotate structure matrix.</i> <code>DISCRIMINANT</code> displays a varimax transformation matrix, a rotated structure matrix, and a rotated standardized canonical discriminant function coefficients table.
<b>NONE</b>	<i>Do not rotate.</i> This is the default.



## **HISTORY Subcommand**

HISTORY controls the display of stepwise and summary output.

- By default, HISTORY displays both the step-by-step output and the summary table (keyword STEP, alias END).

**STEP**     *Display step-by-step and summary output. Alias END. This is the default. See Stepwise statistics in STATISTICS Subcommand on p. 641.*

**NONE**     *Suppress the step-by-step and summary table. Alias NOSTEP, NOEND.*

## **CLASSIFY Subcommand**

CLASSIFY determines how cases are handled during classification.

- By default, all cases with nonmissing values for all predictors are classified, and the pooled within-groups covariance matrix is used to classify cases.
- The default keywords for CLASSIFY are NONMISSING and POOLED.

**NONMISSING**     *Classify all cases that do not have missing values on any predictor variables. Two sets of classification results are produced, one for selected cases (those specified on the SELECT subcommand) and one for unselected cases. This is the default.*

**UNSELECTED**     *Classify only unselected cases. The classification phase is suppressed for cases selected via the SELECT subcommand. If all cases are selected (when the SELECT subcommand is omitted), the classification phase is suppressed for all cases and no classification results are produced.*

**UNCLASSIFIED**     *Classify only unclassified cases. The classification phase is suppressed for cases that fall within the range specified on the GROUPS subcommand.*

**POOLED**     *Use the pooled within-groups covariance matrix to classify cases. This is the default.*

**SEPARATE**     *Use separate-groups covariance matrices of the discriminant functions for classification. DISCRIMINANT displays the group covariances of canonical discriminant functions and Box's test of equality of covariance matrices of canonical discriminant functions. Since classification is based on the discriminant functions and not the original variables, this option is not necessarily equivalent to quadratic discrimination.*

**MEANSUB**     *Substitute means for missing predictor values during classification. During classification, means are substituted for missing values and cases with missing values are classified. Cases with missing values are not used during analysis.*

## **PLOT Subcommand**

PLOT requests additional output to help you examine the effectiveness of the discriminant analysis.

- If PLOT is specified without keywords, the default is COMBINED and CASES.

- If any keywords are requested on PLOT, only the requested plots are displayed.
- If PLOT is specified with any keyword except MAP, DISCRIMINANT displays a classification processing summary table and a prior probabilities for groups table.

<b>COMBINED</b>	<i>All-groups plot.</i> For each case, the first two function values are plotted.
<b>CASES(n)</b>	<i>Casewise statistics.</i> For each case, classification information, squared Mahalanobis distance to centroid for the highest and second highest groups, and discriminant scores of all functions are plotted. Validated statistics are displayed for selected cases if CROSSVALID is specified on STATISTICS. If <i>n</i> is specified, DISCRIMINANT displays the first <i>n</i> cases only.
<b>MAP</b>	<i>Territorial map.</i> A plot of group centroids and boundaries used for classifying groups.
<b>SEPARATE</b>	<i>Separate-groups plots.</i> These are the same types of plots produced by the keyword COMBINED, except that a separate plot is produced for each group. If only one function is used, a histogram is displayed.
<b>ALL</b>	<i>All available plots.</i>

## MISSING Subcommand

MISSING controls the treatment of cases with missing values in the analysis phase. By default, cases with missing values for any variable named on the VARIABLES subcommand are not used in the analysis phase but are used in classification.

- The keyword INCLUDE includes cases with user-missing values in the analysis phase.
- Cases with missing or out-of-range values for the grouping variable are always excluded.

<b>EXCLUDE</b>	<i>Exclude all cases with missing values.</i> Cases with user or system-missing values are excluded from the analysis. This is the default.
<b>INCLUDE</b>	<i>Include cases with user-missing values.</i> User-missing values are treated as valid values. Only the system-missing value is treated as missing.

## MATRIX Subcommand

MATRIX reads and writes IBM® SPSS® Statistics matrix data files.

- Either IN or OUT and the matrix file in parentheses are required. When both IN and OUT are used in the same DISCRIMINANT procedure, they can be specified on separate MATRIX subcommands or on the same subcommand.

<b>OUT ('savfile' 'dataset')</b>	<i>Write a matrix data file.</i> Specify either a quoted file specification, a previously declared dataset name (DATASET DECLARE command) or an asterisk (*), enclosed in parentheses. If you specify an asterisk (*), the matrix data file replaces the active dataset .
<b>IN ('savfile' 'dataset')</b>	<i>Read a matrix data file.</i> Specify either a quoted file specification, a previously declared dataset name (DATASET DECLARE command) or an asterisk (*), enclosed in parentheses. An asterisk indicates the active dataset. A matrix file read from an a file or dataset does not replace the active dataset.

## Matrix Output

- In addition to Pearson correlation coefficients, the matrix materials written by `DISCRIMINANT` include weighted and unweighted numbers of cases, means, and standard deviations. (See Format of the Matrix Data File on p. 645 for a description of the file.) These materials can be used in subsequent `DISCRIMINANT` procedures.
- Any documents contained in the active dataset are not transferred to the matrix file.
- If `BOXM` or `GCOV` is specified on the `STATISTICS` subcommand or `SEPARATE` is specified on the `CLASSIFY` subcommand when a matrix file is written, the `STDDEV` and `CORR` records in the matrix materials represent within-cell data, and separate covariance matrices are written to the file. When the matrix file is used as input for a subsequent `DISCRIMINANT` procedure, at least one of these specifications must be used on that `DISCRIMINANT` command.

## Matrix Input

- `DISCRIMINANT` can read correlation matrices written by a previous `DISCRIMINANT` command or by other procedures. Matrix materials read by `DISCRIMINANT` must contain records with `ROWTYPE_` values `MEAN`, `N` or `COUNT` (or both), `STDDEV`, and `CORR`.
- If the data do not include records with `ROWTYPE_` value `COUNT` (unweighted number of cases), `DISCRIMINANT` uses information from records with `ROWTYPE_` value `N` (weighted number of cases). Conversely, if the data do not have `N` values, `DISCRIMINANT` uses the `COUNT` values. These records can appear in any order in the matrix input file with the following exceptions: the order of split-file groups cannot be violated and all `CORR` vectors must appear consecutively within each split-file group.
- If you want to use a covariance-type matrix as input to `DISCRIMINANT`, you must first use the `MCONVERT` command to change the covariance matrix to a correlation matrix.
- `DISCRIMINANT` can use a matrix from a previous dataset to classify data in the active dataset. The program checks to make sure that the grouping variable (specified on `GROUPS`) and the predictor variables (specified on `VARIABLES`) are the same in the active dataset as in the matrix file. If they are not, the program displays an error message and the classification will not be executed.
- `MATRIX=IN` cannot be used unless a active dataset has already been defined. To read an existing matrix data file at the beginning of a session, first use `GET` to retrieve the matrix file and then specify `IN(*)` on `MATRIX`.

## Format of the Matrix Data File

- The matrix data file has two special variables created by the program: `ROWTYPE_` and `VARNAME_`. Variable `ROWTYPE_` is a short string variable having values `N`, `COUNT`, `MEAN`, `STDDEV`, and `CORR` (for Pearson correlation coefficient). The variable `VARNAME_` is a short string variable whose values are the names of the variables used to form the correlation matrix.
- When `ROWTYPE_` is `CORR`, `VARNAME_` gives the variable associated with that row of the correlation matrix.
- Between `ROWTYPE_` and `VARNAME_` is the grouping variable, which is specified on the `GROUPS` subcommand of `DISCRIMINANT`.
- The remaining variables are the variables used to form the correlation matrix.

### **Split Files**

- When split-file processing is in effect, the first variables in the matrix data file will be split variables, followed by *ROWTYPE\_*, the grouping variable, *VARNAME\_*, and then the variables used to form the correlation matrix.
- A full set of matrix materials is written for each subgroup defined by the split variables.
- A split variable cannot have the same variable name as any other variable written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split file must be in effect when that matrix is read by another procedure.

### **STDDEV and CORR Records**

Records written with *ROWTYPE\_* values *STDDEV* and *CORR* are influenced by specifications on the *STATISTICS* and *CLASSIFY* subcommands.

- If *BOXM* or *GCOV* is specified on *STATISTICS* or *SEPARATE* is specified on *CLASSIFY*, the *STDDEV* and *CORR* records represent within-cell data and receive values for the grouping variable.
- If none of the above specifications is in effect, the *STDDEV* and *CORR* records represent pooled values. The *STDDEV* vector contains the square root of the mean square error for each variable, and *STDDEV* and *CORR* records receive the system-missing value for the grouping variable.

### **Missing Values**

Missing-value treatment affects the values written to a matrix data file. When reading a matrix data file, be sure to specify a missing-value treatment on *DISCRIMINANT* that is compatible with the treatment that was in effect when the matrix materials were generated.

### **Examples**

#### **Writing Output to a Matrix Data File**

```
GET FILE=' /DATA/UNIONBK.SAV' /KEEP WORLD FOOD SERVICE BUS MECHANIC
      CONSTRUC COOK MANAGER FSALES APPL RENT.
DISCRIMINANT GROUPS=WORLD(1,3)
/VARIABLES=FOOD SERVICE BUS MECHANIC CONSTRUC COOK MANAGER FSALES
/METHOD=WILKS
/PRIORS=SIZE
/MATRIX=OUT(' /DATA/DISCMTX.SAV' ) .
```

- *DISCRIMINANT* reads data from the data file *UNIONBK.SAV* and writes one set of matrix materials to the file *DISCMTX.SAV*.
- The active dataset is still *UNIONBK.SAV*. Subsequent commands are executed on this file.

#### **Using Matrix Output to Classify Data in a Different File**

```
GET FILE=UB2 /KEEP WORLD FOOD SERVICE BUS MECHANIC
```

```

                CONSTRUC COOK MANAGER FSALES APPL RENT.
DISCRIMINANT  GROUPS=WORLD(1,3)
/VARIABLES=FOOD SERVICE BUS MECHANIC CONSTRUC COOK MANAGER FSALES
/METHOD=WILKS
/PRIORS=SIZE
/MATRIX=IN(DISCMTX) .

```

- The matrix data file created in the previous example is used to classify data from the file *UB2*.

### **Replacing the Active Dataset with Matrix Data Output**

```

GET FILE=UNIONBK /KEEP WORLD FOOD SERVICE BUS MECHANIC
                CONSTRUC COOK MANAGER FSALES APPL RENT.
DISCRIMINANT  GROUPS=WORLD(1,3)
/VARIABLES=FOOD SERVICE BUS MECHANIC CONSTRUC COOK MANAGER FSALES
/METHOD=WILKS
/PRIORS=SIZE
/MATRIX=OUT(*).
LIST.

```

- DISCRIMINANT writes the same matrix as in the first example. However, the matrix data file replaces the active dataset.
- The LIST command is executed on the matrix file, not on the *UNIONBK* file.

### **Using the Active Dataset as Matrix Input**

```

GET FILE=DISCMTX.
DISCRIMINANT  GROUPS=WORLD(1,3)
/VARIABLES=FOOD SERVICE BUS MECHANIC CONSTRUC COOK MANAGER FSALES
/METHOD=RAO
/MATRIX=IN(*).

```

- This example assumes that you are starting a new session and want to read an existing matrix data file. GET retrieves the matrix data file *DISCMTX*.
- MATRIX=IN specifies an asterisk because the matrix data file is the active dataset. If MATRIX=IN(DISCMTX) is specified, the program issues an error message.
- If the GET command is omitted, the program issues an error message.

### **Using Matrix Output as Matrix Input in the Active Dataset**

```

GET FILE=UNIONBK /KEEP WORLD FOOD SERVICE BUS MECHANIC
                CONSTRUC COOK MANAGER FSALES APPL RENT.
DISCRIMINANT  GROUPS=WORLD(1,3)
/VARIABLES=FOOD SERVICE BUS MECHANIC CONSTRUC COOK MANAGER FSALES
/CLASSIFY=SEPARATE
/MATRIX=OUT(*).
DISCRIMINANT  GROUPS=WORLD(1,3)
/VARIABLES=FOOD SERVICE BUS MECHANIC CONSTRUC COOK MANAGER FSALES
/STATISTICS=BOXM
/MATRIX=IN(*).

```

- The first DISCRIMINANT command creates a matrix with CLASSIFY=SEPARATE in effect. To read this matrix, the second DISCRIMINANT command must specify either BOXM or GCOV on STATISTICS or SEPARATE on CLASSIFY. STATISTICS=BOXM is used.

# DISPLAY

```
DISPLAY [SORTED] [{NAMES**  }] [/VARIABLES=varlist]
                {INDEX      }
                {VARIABLES  }
                {LABELS     }
                {DICTIONARY  }
                {ATTRIBUTES  }
                {@ATTRIBUTES}

                {[SCRATCH]  }
                {[VECTOR]   }
                {[MACROS]   }
                {[DOCUMENTS]}
```

*\*\*Default if the subcommand is omitted.*

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- ATTRIBUTES keyword introduced.

Release 15.0

- @ATTRIBUTES keyword introduced.

## **Example**

```
DISPLAY SORTED DICTIONARY /VARIABLES=DEPT SALARY SEX TO JOBCAT.
```

## **Overview**

DISPLAY exhibits information from the dictionary of the active dataset. The information can be sorted, and it can be limited to selected variables.

### **Basic Specification**

The basic specification is simply the command keyword, which displays an unsorted list of the variables in the active dataset.

**Syntax Rules**

DISPLAY can be specified by itself or with one of the keywords defined below. NAMES is the default. To specify two or more keywords, use multiple DISPLAY commands.

<b>NAMES</b>	<i>Variable names.</i> A list of the variables in the active dataset is displayed.
<b>DOCUMENTS</b>	<i>Documentary text.</i> Documentary text is provided on the DOCUMENT and ADD DOCUMENT commands. No error message is issued if there is no documentary information in the active dataset.
<b>DICTIONARY</b>	<i>Complete dictionary information for variables.</i> Information includes variable names, labels, role, sequential position of each variable in the file, print and write formats, missing values, and value labels.
<b>ATTRIBUTES</b>	<i>Variable and data file attributes, except attributes with names that begin with "@" or "\$@".</i> Custom attributes defined by the VARIABLE ATTRIBUTE and DATAFILE ATTRIBUTE commands.
<b>@ATTRIBUTES</b>	<i>All variable and data file attributes, including those with names that begin with "@" or "\$@".</i>
<b>INDEX</b>	<i>Variable names and positions.</i>
<b>VARIABLES</b>	<i>Variable names, positions, print and write formats, and missing values.</i>
<b>LABELS</b>	<i>Variable names, positions, and variable labels.</i>
<b>SCRATCH</b>	<i>Scratch variable names.</i>
<b>VECTOR</b>	<i>Vector names.</i>
<b>MACROS</b>	<i>Currently defined macros.</i> The macro names are always sorted.

**Operations**

- DISPLAY directs information to the output.
- If SORTED is not specified, information is displayed according to the order of variables in the active dataset.
- DISPLAY is executed as soon as it is encountered in the command sequence, as long as a dictionary has been defined.

**Examples**

```
GET FILE="/data/hub.sav".
DISPLAY DOCUMENTS.
DISPLAY DICTIONARY.
```

- Each DISPLAY command specifies only one keyword. The first requests documentary text and the second requests complete dictionary information for the *hub.sav* file.

**SORTED Keyword**

SORTED alphabetizes the display by variable name. SORTED can precede the keywords NAMES, DICTIONARY, INDEX, VARIABLES, LABELS, SCRATCH, or VECTOR.

**Example**

```
DISPLAY SORTED DICTIONARY.
```

- This command displays complete dictionary information for variables in the active dataset, sorted alphabetically by variable name.

**VARIABLES Subcommand**

VARIABLES (alias NAMES) limits the displayed information to a set of specified variables. VARIABLES must be the last specification on DISPLAY and can follow any specification that requests information about variables (all except VECTOR, SCRATCH, DOCUMENTS, and MACROS).

- The only specification is a slash followed by a list of variables. The slash is optional.
- If the keyword SORTED is not specified, information is displayed in the order in which variables are stored in the active dataset, regardless of the order in which variables are named on VARIABLES.

**Example**

```
DISPLAY SORTED DICTIONARY  
/VARIABLES=DEPT, SALARY, SEX TO JOBCAT.
```

- DISPLAY exhibits dictionary information only for the variables named and implied by the keyword TO on the VARIABLES subcommand, sorted alphabetically by variable name.



# ***DMCLUSTER***

DMCLUSTER is available in the Direct Marketing option.

This is an alias for TWOSTEP CLUSTER. For more information, see the topic TWOSTEP CLUSTER on p. 2065.

# ***DMLOGISTIC***

DMLOGISTIC is available in the Direct Marketing option.

This is an alias for LOGISTIC REGRESSION. For more information, see the topic LOGISTIC REGRESSION on p. 1028.

# ***DMROC***

DMROC is available in the Direct Marketing option.

This is an alias for ROC. [For more information, see the topic ROC on p. 1801.](#)

# ***DMTABLES***

DMTABLES is available in the Direct Marketing option.

This is an alias for CTABLES. [For more information, see the topic CTABLES on p. 510.](#)

# ***DMTREE***

DMTREE is available in the Direct Marketing option.

This is an alias for TREE. [For more information, see the topic TREE on p. 1978.](#)

# DO IF

```
DO IF [(logical expression)]
transformation commands
[ELSE IF [(logical expression)]]
transformation commands
[ELSE IF [(logical expression)]]
.
.
[ELSE]
transformation commands
END IF
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

*The following relational operators can be used in logical expressions:*

Symbol	Definition
EQ or =	Equal to
NE or $\neq$ or $\neq$ or $\neq$	Not equal to
LT or <	Less than
LE or <=	Less than or equal to
GT or >	Greater than
GE or >=	Greater than or equal to

*The following logical operators can be used in logical expressions:*

Symbol	Definition
AND or &	Both relations must be true
OR or	Either relation can be true
NOT	Reverses the outcome of an expression

## Example

```
DO IF (YearHired GT 87).
COMPUTE      Bonus = 0.
ELSE IF (Dept87 EQ 3).
COMPUTE      Bonus = .1*Salary87.
ELSE IF (Dept87 EQ 1).
COMPUTE      Bonus = .12*Salary87.
ELSE IF (Dept87 EQ 4).
COMPUTE      Bonus = .08*Salary87.
ELSE IF (Dept87 EQ 2).
COMPUTE      Bonus = .14*Salary87.
END IF.
```

## Overview

The DO IF-END IF structure conditionally executes one or more transformations on subsets of cases based on one or more logical expressions. The ELSE command can be used within the structure to execute one or more transformations when the logical expression on DO IF is not true. The ELSE IF command within the structure provides further control.

The DO IF-END IF structure is best used for conditionally executing multiple transformation commands, such as COMPUTE, RECODE, and COUNT. DO IF-END IF transforms data for subsets of cases defined by logical expressions. To perform repeated transformations on the same case, use LOOP-END LOOP.

A DO IF-END IF structure can be used within an input program to define complex files that cannot be handled by standard file definition facilities. [For more information, see the topic Complex File Structures on p. 664.](#)

See END FILE for information on using DO IF-END IF to instruct the program to stop reading data before it encounters the end of the file or to signal the end of the file when creating data.

### Basic Specification

The basic specification is DO IF followed by a logical expression, a transformation command, and the END IF command, which has no specifications.

## Examples

### Simple, One-Condition Example

```
DO IF (YearHired LT 87).
RECODE Ethnicity(1=5) (2=4) (4=2) (5=1).
END IF.
```

- The RECODE command recodes *Ethnicity* for those individuals hired before 1987 (*YearHired* is less than 87). The *Ethnicity* variable is not recoded for individuals hired in 1987 or later.
- The RECODE command is skipped for any case with a missing value for *YearHired*.

### Conditional Execution Based on a Logical Expression

```
DATA LIST          FREE / X(F1).
NUMERIC           #QINIT.
DO IF             NOT #QINIT.
+ PRINT EJECT.
+ COMPUTE        #QINIT = 1.
END IF.
PRINT            / X.

BEGIN DATA
1 2 3 4 5
END DATA.
EXECUTE.
```

- This example shows how to execute a command only once.
- The NUMERIC command creates scratch variable *#QINIT*, which is initialized to 0.

- The NOT logical operator on DO IF reverses the outcome of a logical expression. In this example, the logical expression is a numeric variable that takes only 0 (false) or 1 (true) as its values. The PRINT EJECT command is executed only once, when the value of scratch variable #QINIT equals 0. After the COMPUTE command sets #QINIT to 1, the DO IF structure is skipped for all subsequent cases. A scratch variable is used because it is initialized to 0 and is not reinitialized after each case.

## Syntax Rules

- The ELSE IF command is optional and can be repeated as many times as needed.
- The ELSE command is optional. It can be used only once and must follow any ELSE IF commands.
- The END IF command must follow any ELSE IF and ELSE commands.
- A logical expression must be specified on the DO IF and ELSE IF commands. Logical expressions are not used on the ELSE and END IF commands.
- String values used in expressions must be specified in quotation marks and must include any leading or trailing blanks. Lowercase letters are distinguished from uppercase letters.
- To create a new string variable within a DO IF–END IF structure, you must first declare the variable on the STRING command.
- DO IF–END IF structures can be nested to any level permitted by available memory. They can be nested within LOOP–END LOOP structures, and loop structures can be nested within DO IF structures.

### Example

```
DATA LIST FREE /var1.
BEGIN DATA
1 2 3 4 5
END DATA.
DO IF (var1 > 2) & (var1 < 5).
- COMPUTE var2=1.
ELSE IF (var1=2).
- COMPUTE var2=2.
ELSE.
- COMPUTE var2=3.
END IF.
```

var1	var2
1	3
2	2
3	1
4	1
5	3

### Example

```
INPUT PROGRAM.
+ STRING odd (A3).
+ LOOP numvar=1 TO 5.
+ DO IF MOD(numvar, 2)=0.
```



```

+   COMPUTE odd='No' .
+   ELSE.
+   COMPUTE odd='Yes' .
+   END IF.
+   END CASE.
+   END LOOP.
+   END FILE.
END INPUT PROGRAM.

```

numvar	odd
1	Yes
2	No
3	Yes
4	No
5	Yes

## Logical Expressions

- Logical expressions can be simple logical variables or relations, or they can be complex logical tests involving variables, constants, functions, relational operators, and logical operators. Logical expressions can use any of the numeric or string functions allowed in COMPUTE transformations (see COMPUTE).
- Parentheses can be used to enclose the logical expression itself and to specify the order of operations within a logical expression. Extra blanks or parentheses can be used to make the expression easier to read.
- Blanks (*not* commas) are used to separate relational operators from expressions.
- A relation can include variables, constants, or more complicated arithmetic expressions. Relations cannot be abbreviated. For example, the first relation below is valid; the second is not:  
*Valid:* (A EQ 2 OR A EQ 5) *Not valid:* (A EQ 2 OR 5)
- A relation cannot compare a string variable to a numeric value or variable, or vice versa. A relation cannot compare the result of a logical function (SYSMIS, MISSING, ANY, or RANGE) to a number.

## Operations

- DO IF marks the beginning of the control structure and END IF marks the end. Control for a case is passed out of the structure as soon as a logical condition is met on a DO IF, ELSE IF, or ELSE command.
- A logical expression is evaluated as true, false, or missing. A transformation specified for a logical expression is executed only if the expression is true.
- Logical expressions are evaluated in the following order: functions, exponentiation, arithmetic operations, relations, and finally, logical operators. (For strings, the order is functions, relations, and then logical operators.) When more than one logical operator is used, NOT is evaluated first, followed by AND, and then OR. You can change the order of operations using parentheses.

- Numeric variables created within a DO IF structure are initially set to the system-missing value. By default, they are assigned an F8.2 format.
- New string variables created within a DO IF structure are initially set to a blank value and are assigned the format specified on the STRING command that creates them.
- If the transformed value of a string variable exceeds the variable's defined format, the value is truncated. If the value is shorter than the format, the value is right-padded with blanks.
- If WEIGHT is specified within a DO IF structure, it takes effect unconditionally.
- Commands that take effect immediately without reading the data or executing pending commands (e.g., VALUE LABELS, MISSING VALUES, VARIABLE LABEL), commands that don't require an active dataset (e.g., DISPLAY, HOST, INSERT, OMS, SET) and the GET command are processed unconditionally and immediately each time they are encountered in a DO IF structure. The specified condition is ignored for these commands.
- Commands that require a data pass, such as statistical or charting procedures, cannot be included in a DO IF structure and will cause the command to fail.
- The DO IF-END IF structure (like LOOP-END LOOP) can include commands such as DATA LIST, END CASE, END FILE, and REREAD, which define complex file structures.

### Flow of Control

- If the logical expression on DO IF is true, the commands immediately following DO IF are executed up to the next ELSE IF, ELSE, or END IF command. Control then passes to the first statement following END IF.
- If the expression on DO IF is false, control passes to the following ELSE IF command. Multiple ELSE IF commands are evaluated in the order in which they are specified until the logical expression on one of them is true. Commands following that ELSE IF command are executed up to the ELSE or END IF command, and control passes to the first statement following END IF.
- If none of the expressions are true on the DO IF or any of the ELSE IF commands, the commands following ELSE are executed and control passes out of the structure. If there is no ELSE command, a case goes through the entire structure with no change.
- Missing values returned by the logical expression on DO IF or on any ELSE IF cause control to pass to the END IF command at that point.

### Missing Values and Logical Operators

When two or more relations are joined by logical operators AND and OR, the program always returns missing if all of the relations in the expression are missing. However, if any one of the relations can be determined, the program tries to return true or false according to the logical outcomes shown in the following table. The asterisk indicates situations where the program can evaluate the outcome with incomplete information.

Table 72-1  
Logical outcomes

Expression	Outcome	Expression	Outcome
true AND true	= true	true OR true	= true

Expression	Outcome	Expression	Outcome
true AND false	= false	true OR false	= true
false AND false	= false	false OR false	= false
true AND missing	= missing	true OR missing	= true*
missing AND missing	= missing	missing OR missing	= missing
false AND missing	= false*	false OR missing	= missing

## ELSE Command

ELSE executes one or more transformations when none of the logical expressions on DO IF or any ELSE IF commands is true.

- Only one ELSE command is allowed within a DO IF–END IF structure.
- ELSE must follow all ELSE IF commands (if any) in the structure.
- If the logical expression on DO IF or any ELSE IF command is true, the program ignores the commands following ELSE.

### Example

```
DO IF (X EQ 0).
  COMPUTE Y=1.
ELSE.
  COMPUTE Y=2.
END IF.
```

- Y is set to 1 for all cases with value 0 for X, and Y is 2 for all cases with any other valid value for X.
- The value of Y is not changed by this structure if X is missing.

### Example

```
DO IF (YearHired GT 87).
  COMPUTE          Bonus = 0.
ELSE.
  IF (Dept87 EQ 1) Bonus = .12*Salary87.
  IF (Dept87 EQ 2) Bonus = .14*Salary87.
  IF (Dept87 EQ 3) Bonus = .1*Salary87.
  IF (Dept87 EQ 4) Bonus = .08*Salary87.
END IF.
```

- If an individual was hired after 1987 (*YearHired* is greater than 87), *Bonus* is set to 0 and control passes out of the structure. Otherwise, control passes to the IF commands following ELSE.
- Each IF command evaluates every case. The value of *Bonus* is transformed only when the case meets the criteria specified on IF. Compare this structure with the second example for the ELSE IF command, which performs the same task more efficiently.

### Example

```
* Test for listwise deletion of missing values.
DATA LIST / V1 TO V6 1-6.
```

*DO IF*

```

BEGIN DATA
123456
   56
1 3456
123456
123456
END DATA.

DO IF NMISS(V1 TO V6)=0.
+ COMPUTE SELECT='V' .
ELSE
+ COMPUTE SELECT='M' .
END IF.

FREQUENCIES VAR=SELECT.

```

- If there are no missing values for any of the variables *V1* to *V6*, *COMPUTE* sets the value of *SELECT* equal to *V* (for valid). Otherwise, *COMPUTE* sets the value of *SELECT* equal to *M* (for missing).
- *FREQUENCIES* generates a frequency table for *SELECT*. The table gives a count of how many cases have missing values for one or more variables, and how many cases have valid values for all variables. Commands in this example can be used to determine how many cases are dropped from an analysis that uses listwise deletion of missing values.

***ELSE IF Command***

*ELSE IF* executes one or more transformations when the logical expression on *DO IF* is not true.

- Multiple *ELSE IF* commands are allowed within the *DO IF*–*END IF* structure.
- If the logical expression on *DO IF* is true, the program executes the commands immediately following *DO IF* up to the first *ELSE IF*. Then control passes to the command following the *END IF* command.
- If the result of the logical expression on *DO IF* is false, control passes to *ELSE IF*.

***Example***

```

STRING Stock(A9).
DO IF (ITEM EQ 0).
COMPUTE Stock='New'.
ELSE IF (ITEM LE 9).
COMPUTE Stock='Old'.
ELSE.
COMPUTE Stock='Cancelled'.
END IF.

```

- *STRING* declares string variable *Stock* and assigns it a width of nine characters.
- The first *COMPUTE* is executed for cases with value 0 for *ITEM*, and then control passes out of the structure. Such cases are not reevaluated by *ELSE IF*, even though 0 is less than 9.
- When the logical expression on *DO IF* is false, control passes to the *ELSE IF* command, where the second *COMPUTE* is executed only for cases with *ITEM* less than or equal to 9. Then control passes out of the structure.
- If the logical expressions on both the *DO IF* and *ELSE IF* commands are false, control passes to *ELSE*, where the third *COMPUTE* is executed.

- The DO IF-END IF structure sets *Stock* equal to New when *ITEM* equals 0, to Old when *ITEM* is less than or equal to 9 but not equal to 0 (including negative numbers if they are valid), and to Cancelled for all valid values of *ITEM* greater than 9. The value of *Stock* remains blank if *ITEM* is missing.

### Example

```
DO IF (YearHired GT 87).
COMPUTE          Bonus = 0.
ELSE IF (Dept87 EQ 3).
COMPUTE          Bonus = .1*Salary87.
ELSE IF (Dept87 EQ 1).
COMPUTE          Bonus = .12*Salary87.
ELSE IF (Dept87 EQ 4).
COMPUTE          Bonus = .08*Salary87.
ELSE IF (Dept87 EQ 2).
COMPUTE          Bonus = .14*Salary87.
END IF.
```

- For cases hired after 1987, *Bonus* is set to 0 and control passes out of the structure. For a case that was hired before 1987 with value 3 for *Dept87*, *Bonus* equals 10% of salary. Control then passes out of the structure. The other three ELSE IF commands are not evaluated for that case. This differs from the second example for the ELSE command, where the IF command is evaluated for every case. The DO IF-ELSE IF structure shown here is more efficient.
- If Department 3 is the largest, Department 1 the next largest, and so forth, control passes out of the structure quickly for many cases. For a large number of cases or a command file that will be executed frequently, these efficiency considerations can be important.

## Nested DO IF Structures

To perform transformations involving logical tests on two variables, you can use nested DO IF-END IF structures.

- There must be an END IF command for every DO IF command in the structure.

### Example

```
DO IF (Ethnicity EQ 5).          /*Do whites
+ DO IF (Gender EQ 2).          /*White female
+ COMPUTE Gender_Ethnicity=3.
+ ELSE.                          /*White male
+ COMPUTE Gender_Ethnicity=1.
+ END IF.                          /*Whites done
ELSE IF (Gender EQ 2).          /*Nonwhite female
COMPUTE Gender_Ethnicity=4.
ELSE.                          /*Nonwhite male
COMPUTE Gender_Ethnicity=2.
END IF.                          /*Nonwhites done
```

- This structure creates variable *Gender\_Ethnicity*, which indicates both the sex and minority status of an individual.
- An optional plus sign, minus sign, or period in the first column allows you to indent commands so you can easily see the nested structures.

## Complex File Structures

Some complex file structures may require you to embed more than one `DATA LIST` command inside a `DO IF-END IF` structure. For example, consider a data file that has been collected from various sources. The information from each source is basically the same, but it is in different places on the records:

```
111295100FORD      CHAPMAN AUTO SALES
121199005VW      MIDWEST VOLKSWAGEN SALES
11 395025FORD      BETTER USED CARS
11      CHEVY 195005      HUFFMAN SALES & SERVICE
11      VW 595020      MIDWEST VOLKSWAGEN SALES
11      CHEVY 295015      SAM'S AUTO REPAIR
12      CHEVY 210 20      LONGFELLOW CHEVROLET
 9555032 VW      HYDE PARK IMPORTS
```

In the above file, an automobile part number always appears in columns 1 and 2, and the automobile manufacturer always appears in columns 10 through 14. The location of other information, such as price and quantity, depends on both the part number and the type of automobile. The `DO IF-END IF` structure in the following example reads records for part type 11.

### Example

```
INPUT PROGRAM.
DATA LIST FILE="/data/carparts.txt"
  /PARTNO 1-2 KIND 10-14 (A).

DO IF (PARTNO EQ 11 AND KIND EQ 'FORD').
+ REREAD.
+ DATA LIST /PRICE 3-6 (2) QUANTITY 7-9 BUYER 20-43 (A).
+ END CASE.

ELSE IF (PARTNO EQ 11 AND (KIND EQ 'CHEVY' OR KIND EQ 'VW')).
+ REREAD.
+ DATA LIST /PRICE 15-18 (2) QUANTITY 19-21 BUYER 30-53 (A).
+ END CASE.
END IF.
END INPUT PROGRAM.

PRINT FORMATS PRICE (DOLLAR6.2).
PRINT /PARTNO TO BUYER.
WEIGHT BY QUANTITY.
DESCRIPTIVES PRICE.
```

- The first `DATA LIST` extracts the part number and the type of automobile.
- Depending on the information from the first `DATA LIST`, the records are reread, pulling the price, quantity, and buyer from different places.
- The two `END CASE` commands limit the working file to only those cases with a part number of 11 and automobile type of Ford, Chevrolet, or Volkswagen. Without the `END CASE` commands, cases would be created in the working file for other part numbers and automobile types with missing values for price, quantity, and buyer.
- The results of the `PRINT` command are shown below.

Figure 72-1

*Printed information for part 11*

```
11 FORD $12.95 100 CHAPMAN AUTO SALES
```

11	FORD	\$3.95	25	BETTER USED CARS
11	CHEVY	\$1.95	5	HUFFMAN SALES & SERVICE
11	VW	\$5.95	20	MIDWEST VOLKSWAGEN SALES
11	CHEVY	\$2.95	15	SAM'S AUTO REPAIR

# DO REPEAT-END REPEAT

```
DO REPEAT stand-in var={varlist | ALL } [ /stand-in var=... ]  
    {value list}  
  
transformation commands  
  
END REPEAT [PRINT]
```

## Release History

Release 14.0

- ALL keyword introduced.

## Example

```
DO REPEAT var=var1 to var5  
    /value=1 to 5.  
COMPUTE var=value.  
END REPEAT.
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## Overview

The DO REPEAT–END REPEAT structure repeats the same transformations on a specified set of variables, reducing the number of commands you must enter to accomplish a task. This utility does not reduce the number of commands the program executes, just the number of commands you enter. To display the expanded set of commands the program generates, specify PRINT on END REPEAT.

DO REPEAT uses a **stand-in variable** to represent a **replacement list** of variables or values. The stand-in variable is specified as a placeholder on one or more transformation commands within the structure. When the program repeats the transformation commands, the stand-in variable is replaced, in turn, by each variable or value specified on the replacement list.

The following commands can be used within a DO REPEAT–END REPEAT structure:

- Data transformations: COMPUTE, RECODE, IF, COUNT, and SELECT IF
- Data declarations: VECTOR, STRING, NUMERIC, and LEAVE
- Data definition: DATA LIST, MISSING VALUES (but not VARIABLE LABELS or VALUE LABELS)
- Loop structure commands: LOOP, END LOOP, and BREAK
- Do-if structure commands: DO IF, ELSE IF, ELSE, and END IF
- Print and write commands: PRINT, PRINT EJECT, PRINT SPACE, and WRITE
- Format commands: PRINT FORMATS, WRITE FORMATS, and FORMATS



### **Basic Specification**

The basic specification is `DO REPEAT`, a stand-in variable followed by a required equals sign and a replacement list of variables or values, and at least one transformation command. The structure must end with the `END REPEAT` command. On the transformation commands, a single stand-in variable represents every variable or value specified on the replacement list.

### **Syntax Rules**

- Multiple stand-in variables can be specified on a `DO REPEAT` command. Each stand-in variable must have its own equals sign and associated variable or value list and must be separated from other stand-in variables by a slash. All lists must name or generate the same number of items.
- Stand-in variables can be assigned any valid variable names: permanent, temporary, scratch, system, and so forth. A stand-in variable does not exist outside the `DO REPEAT-END REPEAT` structure and has no effect on variables with the same name that exist outside the structure. However, two stand-in variables cannot have the same name within the same `DO REPEAT` structure.
- A replacement variable list can include either new or existing variables. You cannot mix new and existing variables in the same replacement list.
- Keyword `TO` can be used to name consecutive existing variables and to create a set of new variables, and keyword `ALL` can be used to specify all variables.
- New string variables must be declared on the `STRING` command either before `DO REPEAT` or within the `DO REPEAT` structure.
- All replacement variable and value lists must have the same number of items.
- A replacement value list can be a list of strings or numeric values, or it can be of the form  $n_1$  `TO`  $n_2$ , where  $n_1$  is less than  $n_2$  and both are integers. (Note that the keyword is `TO`, not `THRU`.)

### **Operations**

- `DO REPEAT` marks the beginning of the control structure and `END REPEAT` marks the end. Once control passes out of the structure, all stand-in variables defined within the structure cease to exist.
- The program repeats the commands between `DO REPEAT` and `END REPEAT` once for each variable or value on the replacement list.
- Numeric variables created within the structure are initially set to the system-missing value. By default, they are assigned an `F8.2` format.
- New string variables declared within the structure are initially set to a blank value and are assigned the format specified on the `STRING` command that creates them.
- If `DO REPEAT` is used to create new variables, the order in which they are created depends on how the transformation commands are specified. Variables created by specifying the `TO` keyword (for example, `V1 TO V5`) are not necessarily consecutive in the active dataset. [For more information, see the topic PRINT Subcommand on p. 669.](#)
- Multiple replacement lists are stepped through in parallel, not in a nested fashion, and all replacement lists must name or generate the same number of items.

## Examples

### Creating Multiple New Variables with the Same Value

```
DO REPEAT R=REGION1 TO REGION5.
COMPUTE R=0.
END REPEAT.
```

- DO REPEAT defines the stand-in variable *R*, which represents five new numeric variables: *REGION1*, *REGION2*, *REGION3*, *REGION4*, and *REGION5*.
- The five variables are initialized to 0 by a single COMPUTE specification that is repeated for each variable on the replacement list. Thus, the program generates five COMPUTE commands from the one specified.
- Stand-in variable *R* ceases to exist once control passes out of the DO REPEAT structure.

### Multiple Replacement Lists

```
DO REPEAT existVar=firstVar TO var5
      /newVar=new1 TO new5
      /value=1 TO 5.
COMPUTE newVar=existVar*value.
END REPEAT PRINT.
```

```
****generated COMPUTE commands****
57 +COMPUTE      new1=firstVar*1
58 +COMPUTE      new2=secondVar*2
59 +COMPUTE      new3=var3*3
60 +COMPUTE      new4=fourthVar*4
61 +COMPUTE      new5=var5*5.
```

- existVar=firstVar to var5 includes all existing variables from *firstVar* to *var5*, in file order.
- newVar=new1 TO new5 specifies five new variables: *var1*, *var2*, *var3*, *var4*, and *var5*.
- value=1 to 5 specifies a list of five consecutive values: 1, 2, 3, 4, 5.
- All three replacement lists contain five items, and five COMPUTE commands are generated.

### Generating Data with DO REPEAT, LOOP, and INPUT PROGRAM

\* This example shows a typical application of INPUT PROGRAM, LOOP, and DO REPEAT. A data file containing random numbers is generated.

```
INPUT PROGRAM.
+ LOOP #I = 1 TO 1000.
+   DO REPEAT RESPONSE = R1 TO R400.
+     COMPUTE RESPONSE = UNIFORM(1) > 0.5.
+   END REPEAT.
+   COMPUTE AVG = MEAN(R1 TO R400).
+   END CASE.
+ END LOOP.
+ END FILE.
END INPUT PROGRAM.
```

```
FREQUENCIES VARIABLE=AVG
/FORMAT=CONDENSE
/HISTOGRAM
/STATISTICS=MEAN MEDIAN MODE STDDEV MIN MAX.
```

- The INPUT PROGRAM–END INPUT PROGRAM structure encloses an input program that builds cases from transformation commands.
- The indexing variable (*#I*) on LOOP–END LOOP indicates that the loop should be executed 1000 times.
- The DO REPEAT–END REPEAT structure generates 400 variables, each with a 50% chance of being 0 and a 50% chance of being 1. This is accomplished by specifying a logical expression on COMPUTE that compares the values returned by UNIFORM(1) to the value 0.5. (UNIFORM(1) generates random numbers between 0 and 1.) Logical expressions are evaluated as false (0), true (1), or missing. Thus, each random number returned by UNIFORM that is 0.5 or less is evaluated as false and assigned the value 0, and each random number returned by UNIFORM that is greater than 0.5 is evaluated as true and assigned the value 1.
- The second COMPUTE creates variable *AVG*, which is the mean of *R1* to *R400* for each case.
- END CASE builds a case with the variables created within each loop. Thus, the loop structure creates 1000 cases, each with 401 variables (*R1* to *R400*, and *AVG*).
- END FILE signals the end of the data file generated by the input program. If END FILE were not specified in this example, the input program would go into an infinite loop. No dataset would be built, and the program would display an error message for every procedure that follows the input program.
- FREQUENCIES produces a condensed frequency table, histogram, and statistics for *AVG*. The histogram for *AVG* shows a normal distribution.

## **PRINT Subcommand**

The PRINT subcommand on END REPEAT displays the commands generated by the DO REPEAT–END REPEAT structure. PRINT can be used to verify the order in which commands are executed.

### **Example**

```
DO REPEAT Q=Q1 TO Q5/ R=R1 TO R5.
COMPUTE Q=0.
COMPUTE R=1.
END REPEAT PRINT.
```

- The DO REPEAT–END REPEAT structure initializes one set of variables to 0 and another set to 1.
- The output from the PRINT subcommand is shown below. The generated commands are preceded by plus signs.
- The COMPUTE commands are generated in such a way that variables are created in alternating order: *Q1*, *R1*, *Q2*, *R2*, and so forth. If you plan to use the TO keyword to refer to *Q1* to *Q5* later, you should use two separate DO REPEAT utilities; otherwise, *Q1* to *Q5* will include four of the five *R* variables. Alternatively, use the NUMERIC command to predetermine the order in which variables are added to the active dataset, or specify the replacement value lists as shown in the next example.

**Figure 73-1**  
Output from the *PRINT* subcommand

```

2      0      DO REPEAT Q=Q1 TO Q5/ R=R1 TO R5
3      0      COMPUTE Q=0
4      0      COMPUTE R=1
5      0      END REPEAT PRINT

6      0      +COMPUTE Q1=0
7      0      +COMPUTE R1=1
8      0      +COMPUTE Q2=0
9      0      +COMPUTE R2=1
10     0      +COMPUTE Q3=0
11     0      +COMPUTE R3=1
12     0      +COMPUTE Q4=0
13     0      +COMPUTE R4=1
14     0      +COMPUTE Q5=0
15     0      +COMPUTE R5=1

```

**Example**

```

DO REPEAT Q=Q1 TO Q5, R1 TO R5/ N=0,0,0,0,0,1,1,1,1,1.
COMPUTE Q=N.
END REPEAT PRINT.

```

- In this example, a series of constants is specified as a stand-in value list for *N*. All the *Q* variables are initialized first, and then all the *R* variables, as shown below.

**Figure 73-2**  
Output from the *PRINT* subcommand

```

2      0      DO REPEAT Q=Q1 TO Q5, R1 TO R5/ N=0,0,0,0,0,1,1,1,1,1
3      0      COMPUTE Q=N
4      0      END REPEAT PRINT

5      0      +COMPUTE Q1=0
6      0      +COMPUTE Q2=0
7      0      +COMPUTE Q3=0
8      0      +COMPUTE Q4=0
9      0      +COMPUTE Q5=0
10     0      +COMPUTE R1=1
11     0      +COMPUTE R2=1
12     0      +COMPUTE R3=1
13     0      +COMPUTE R4=1
14     0      +COMPUTE R5=1

```

**Example**

```

DO REPEAT R=REGION1 TO REGION5/ X=1 TO 5.
COMPUTE R=REGION EQ X.
END REPEAT PRINT.

```

- In this example, stand-in variable *R* represents the variable list *REGION1* to *REGION5*. Stand-in variable *X* represents the value list 1 to 5.
- The `DO REPEAT-END REPEAT` structure creates dummy variables *REGION1* to *REGION5* that equal 0 or 1 for each of 5 regions, depending on whether variable *REGION* equals the current value of stand-in variable *X*.
- `PRINT` on `END REPEAT` causes the program to display the commands generated by the structure, as shown below.

**Figure 73-3**  
*Commands generated by DO REPEAT*

```
2      0      DO REPEAT R=REGION1 TO REGION5/ X=1 TO 5
3      0      COMPUTE R=REGION EQ X
4      0      END REPEAT PRINT

5      0      +COMPUTE REGION1=REGION EQ 1
6      0      +COMPUTE REGION2=REGION EQ 2
7      0      +COMPUTE REGION3=REGION EQ 3
8      0      +COMPUTE REGION4=REGION EQ 4
9      0      +COMPUTE REGION5=REGION EQ 5
```

# DOCUMENT

DOCUMENT text

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
DOCUMENT      This file contains a subset of variables from the
               General Social Survey data.  For each case it records
               only the age, sex, education level, marital status,
               number of children, and type of medical insurance
               coverage.
```

## **Overview**

DOCUMENT saves a block of text of any length in data files saved in IBM® SPSS® Statistics format. The documentation can be displayed with the DISPLAY command. (See also ADD DOCUMENT.)

When GET retrieves a data file, or when ADD FILES, MATCH FILES, or UPDATE is used to combine data files, all documents from each specified file are copied into the active dataset. DROP DOCUMENTS can be used to drop those documents from the active dataset. Whether or not DROP DOCUMENTS is used, new documents can be added to the active dataset with the DOCUMENT command.

## **Basic Specification**

The basic specification is DOCUMENT followed by any length of text. The text is stored in the file dictionary when the data file is saved in SPSS Statistics format.

## **Syntax Rules**

- The text can be entered on as many lines as needed.
- Blank lines can be used to separate paragraphs.
- A period at the end of a line terminates the command, so you should not place a period at the end of any line but the last.
- Multiple DOCUMENT commands can be used within the command sequence. However, the DISPLAY command cannot be used to exhibit the text from a particular DOCUMENT command. DISPLAY shows all existing documentation.

## **Operations**

- The documentation and the date it was entered are saved in the data file's dictionary. New documentation is saved along with any documentation already in the active dataset.

- If a `DROP DOCUMENTS` command *follows* a `DOCUMENT` command anywhere in the command sequence, the documentation added by that `DOCUMENT` command is dropped from the active dataset along with all other documentation.

## Examples

### **Adding Descriptive Text to a Data File**

```
GET FILE="/data/gensoc.sav" /KEEP=AGE SEX EDUC MARITAL CHILDREN MED_INS.  
FILE LABEL General Social Survey subset.
```

```
DOCUMENT      This file contains a subset of variables from the  
              General Social Survey data.  For each case it records  
              only the age, sex, education level, marital status,  
              number of children, and type of medical insurance  
              coverage.
```

```
SAVE OUTFILE="/data/subsoc.sav".
```

- `GET` keeps only a subset of variables from the file *gensoc.sav*. All documentation from the file *GENSOC* is copied into the active dataset.
- `FILE LABEL` creates a label for the new active dataset.
- `DOCUMENT` specifies the new document text. Both existing documents from the file *GENSOC* and the new document text are saved in the file *subsoc.sav*.

### **Replacing Existing DOCUMENT Text**

```
GET FILE="/data/gensoc.sav" /KEEP=AGE SEX EDUC MARITAL CHILDREN MED_INS.
```

```
DROP DOCUMENTS.
```

```
FILE LABEL General Social Survey subset.
```

```
DOCUMENT      This file contains a subset of variables from the  
              General Social Survey data.  For each case it records  
              only the age, sex, education level, marital status,  
              number of children, and type of medical insurance  
              coverage.
```

```
SAVE OUTFILE="/data/subsoc.sav".
```

- `DROP DOCUMENTS` drops the documentation from the file *gensoc.sav* as data are copied into the active dataset. Only the new documentation specified on `DOCUMENT` is saved in the file *subsoc.sav*.

# DROP DOCUMENTS

DROP DOCUMENTS

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Overview

When GET retrieves IBM® SPSS® Statistics data files, or when ADD FILES, MATCH FILES, or UPDATE are used to combine SPSS Statistics data files, all documents from each specified file are copied into the active dataset. DROP DOCUMENTS is used to drop these or any documents added with the DOCUMENT command from the active dataset. Whether or not DROP DOCUMENTS is used, new documents can be added to the active dataset with the DOCUMENT command.

### Basic Specification

The only specification is DROP DOCUMENTS. There are no additional specifications.

### Operations

- Documents are dropped from the active dataset only. The original data file is unchanged, unless it is resaved.
- DROP DOCUMENTS drops all documentation, including documentation added by any DOCUMENT commands specified prior to the DROP DOCUMENTS command.

## Examples

```
GET FILE="/data/gensoc.sav"  
  /KEEP=AGE SEX EDUC MARITAL CHILDREN MED_INS.
```

```
DROP DOCUMENTS.
```

```
FILE LABEL  General Social Survey Subset.  
DOCUMENT    This file contains a subset of variables from the  
             General Social Survey data. For each case it records  
             only the age, sex, education level, marital status,  
             number of children, and type of medical insurance  
             coverage.
```

```
SAVE OUTFILE="/data/subsoc.sav" .
```

- DROP DOCUMENTS drops the documentation text from data file. Only the new documentation added with the DOCUMENT command is saved in file *subsoc.sav*.
- The original file *gensoc.sav* is unchanged.



# ***ECHO***

```
ECHO "text" .
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Example***

```
ECHO "Hey! Look at this!" .
```

## ***Overview***

ECHO displays the quoted text string as text output.

### ***Basic Specification***

The basic specification is the command name ECHO followed by a quoted text string.

### ***Syntax Rules***

The text string must be enclosed in single or double quotes, following the standard rules for quoted strings.

- The text string can be continued on multiple lines by enclosing each line in quotes and using the plus sign (+) to combine the strings; the string will be displayed on a single line in output.

# END CASE

END CASE

## **Example**

\* Restructure a data file to make each data item into a single case.

```
INPUT PROGRAM.  
DATA LIST /#X1 TO #X3 (3(F1,1X)).  
  
VECTOR V=#X1 TO #X3.  
  
LOOP #I=1 TO 3.  
- COMPUTE X=V(#I).  
- END CASE.  
END LOOP.  
END INPUT PROGRAM.
```

## **Overview**

END CASE is used in an INPUT PROGRAM–END INPUT PROGRAM structure to signal that a case is complete. Control then passes to the commands immediately following the input program. After these commands are executed for the newly created case, the program returns to the input program and continues building cases by processing the commands immediately after the last END CASE command that was executed. For more information about the flow control in an input program, see INPUT PROGRAM–END INPUT PROGRAM.

END CASE is especially useful for restructuring files, either building a single case from several cases or building several cases from a single case. It can also be used to generate data without any data input (see DO REPEAT for an example).

## **Basic Specification**

The basic specification is simply END CASE. There are no additional specifications.

## **Syntax Rules**

- END CASE is available only within an input program and is generally specified within a loop.
- Multiple END CASE commands can be used within an input program. Each builds a case from the transformation and data definition commands executed since the last END CASE command.
- If no END CASE is explicitly specified, an END CASE command is implied immediately before END INPUT PROGRAM and the input program loops until an end-of-file is encountered or specified (see END FILE).

## **Operations**

- When an END CASE command is encountered, the program suspends execution of the rest of the commands before the END INPUT PROGRAM command and passes control to the commands after the input program. After these commands are executed for the new case,

control returns to the input program. The program continues building cases by processing the commands immediately after the most recent `END CASE` command. Use a loop to build cases from the same set of transformation and data definition commands.

- When multiple `END CASE` commands are specified, the program follows the flow of the input program and builds a case whenever it encounters an `END CASE` command, using the set of commands executed since the last `END CASE`.
- Unless `LEAVE` is specified, all variables are reinitialized each time the input program is resumed.
- When transformations such as `COMPUTE`, definitions such as `VARIABLE LABELS`, and utilities such as `PRINT` are specified between the last `END CASE` command and `END INPUT PROGRAM`, they are executed while a case is being initialized, not when it is complete. This may produce undesirable results.

## Examples

### **Restructuring a data file to make each data item a single case**

```
INPUT PROGRAM.
DATA LIST /#X1 TO #X3 (3(F1,1X)).

VECTOR V=#X1 TO #X3.

LOOP #I=1 TO 3.
- COMPUTE X=V(#I).
- END CASE.
END LOOP.
END INPUT PROGRAM.

BEGIN DATA
2 1 1
3 5 1
END DATA.
FORMAT X(F1.0).
PRINT / X.
EXECUTE.
```

- The input program encloses the commands that build cases from the input file. An input program is required because `END CASE` is used to create multiple cases from single input records.
- `DATA LIST` defines three variables. In the format specification, the number 3 is a repetition factor that repeats the format in parentheses three times, once for each variable. The specified format is `F1` and the `1X` specification skips one column.
- `VECTOR` creates the vector  $V$  with the original scratch variables as its three elements. The indexing expression on the `LOOP` command increments the variable `#I` three times to control the number of iterations per input case and to provide the index for the vector  $V$ .
- `COMPUTE` sets  $X$  equal to each of the scratch variables. `END CASE` tells the program to build a case. Thus, the first loop (for the first case) sets  $X$  equal to the first element of vector  $V$ . Since  $V(1)$  references `#X1`, and `#X1` is 2, the value of  $X$  is 2. Variable  $X$  is then formatted and printed before control returns to the command `END LOOP`. The loop continues, since indexing is not complete. Thus, the program then sets  $X$  to `#X2`, which is 1, builds the second case, and passes it to the `FORMAT` and `PRINT` commands. After the third iteration, which sets  $X$  equal to

1, the program formats and prints the case and terminates the loop. Since the end of the file has not been encountered, `END INPUT PROGRAM` passes control to the first command in the input program, `DATA LIST`, to read the next input case. After the second loop, however, the program encounters `END DATA` and completes building the active dataset.

- The six new cases are shown below.

**Figure 77-1**

*Outcome for multiple cases read from a single case*

```
2
1
1
3
5
1
```

### ***Restructuring a data file to create a separate case for each book order***

```
INPUT PROGRAM.
DATA LIST /ORDER 1-4 #X1 TO #X22 (1X,11(F3.0,F2.0,1X)).

LEAVE ORDER.
VECTOR BOOKS=#X1 TO #X22.

LOOP #I=1 TO 21 BY 2 IF NOT SYSMIS(BOOKS(#I)).
- COMPUTE ISBN=BOOKS(#I).
- COMPUTE QUANTITY=BOOKS(#I+1).
- END CASE.
END LOOP.
END INPUT PROGRAM.
BEGIN DATA
1045 182 2 155 1 134 1 153 5
1046 155 3 153 5 163 1
1047 161 5 182 2 163 4 186 6
1048 186 2
1049 155 2 163 2 153 2 074 1 161 1
END DATA.

SORT CASES ISBN.
DO IF $CASENUM EQ 1.
- PRINT EJECT /'Order ISBN Quantity'.
- PRINT SPACE.
END IF.

FORMATS ISBN (F3)/ QUANTITY (F2).
PRINT /' ' ORDER ' ' ISBN ' ' QUANTITY.

EXECUTE.
```

- Data are extracted from a file whose records store values for an invoice number and a series of book codes and quantities ordered. For example, invoice 1045 is for four different titles and a total of nine books: two copies of book 182, one copy each of 155 and 134, and five copies of book 153. The task is to break each individual book order into a record, preserving the order number on each new case.
- The input program encloses the commands that build cases from the input file. They are required because the `END CASE` command is used to create multiple cases from single input records.

- `DATA LIST` specifies `ORDER` as a permanent variable and defines 22 scratch variables to hold the book numbers and quantities (this is the maximum number of numbers and quantities that will fit in 72 columns). In the format specification, the first element skips one space after the value for the variable `ORDER`. The number 11 repeats the formats that follow it 11 times: once for each book number and quantity pair. The specified format is `F3.0` for book numbers and `F2.0` for quantities. The `1X` specification skips one column after each quantity value.
- `LEAVE` preserves the value of the variable `ORDER` across the new cases to be generated.
- `VECTOR` sets up the vector `BOOKS` with the 22 scratch variables as its elements. The first element is `#X1`, the second is `#X2`, and so on.
- If the element for the vector `BOOKS` is not system-missing, `LOOP` initiates the loop structure that moves through the vector `BOOKS`, picking off the book numbers and quantities. The indexing clause initiates the indexing variable `#I` at 1, to be increased by 2 to a maximum of 21.
- The first `COMPUTE` command sets the variable `ISBN` equal to the element in the vector `BOOKS` indexed by `#I`, which is the current book number. The second `COMPUTE` sets the variable `QUANTITY` equal to the next element in the vector `BOOKS`, `#I + 1`, which is the quantity associated with the book number in `BOOKS(#I)`.
- `END CASE` tells the program to write out a case with the current values of the three variables: `ORDER`, `ISBN`, and `QUANTITY`.
- `END LOOP` terminates the loop structure and control is returned to the `LOOP` command, where `#I` is increased by 2 and looping continues until the entire input case is read or until `#I` exceeds the maximum value of 21.
- `SORT CASES` sorts the new cases by book number.
- The `DO IF` structure encloses a `PRINT EJECT` command and a `PRINT SPACE` command to set up titles for the output.
- `FORMATS` establishes dictionary formats for the new variables `ISBN` and `QUANTITY`. `PRINT` displays the new cases.
- `EXECUTE` runs the commands. The output is shown below.

Figure 77-2  
*PRINT* output showing new cases

```
Order ISBN Quantity
1049 74 1
1045 134 1
1045 153 5
1046 153 5
1049 153 2
1045 155 1
1046 155 3
1049 155 2
1047 161 5
1049 161 1
1046 163 1
1047 163 4
1049 163 2
1045 182 2
1047 182 2
1047 186 6
1048 186 2
```

---

 END CASE

### **Create variable that approximates a log-normal distribution**

```

SET FORMAT=F8.0.

INPUT PROGRAM.
LOOP I=1 TO 1000.
+ COMPUTE SCORE=EXP(NORMAL(1)).
+ END CASE.
END LOOP.
END FILE.
END INPUT PROGRAM.

FREQUENCIES VARIABLES=SCORE /FORMAT=NOTABLE /HISTOGRAM
/PERCENTILES=1 10 20 30 40 50 60 70 80 90 99
/STATISTICS=ALL.

```

- The input program creates 1,000 cases with a single variable *SCORE*. Values for *SCORE* approximate a log-normal distribution.

### **Restructure a data file to create a separate case for each individual**

```

INPUT PROGRAM.
DATA LIST /#RECS 1 HEAD1 HEAD2 3-4(A). /*Read header info
LEAVE HEAD1 HEAD2.

LOOP #I=1 TO #RECS.
DATA LIST /INDIV 1-2(1). /*Read individual info
PRINT /#RECS HEAD1 HEAD2 INDIV.
END CASE. /*Create combined case
END LOOP.
END INPUT PROGRAM.
BEGIN DATA
1 AC
91
2 CC
35
43
0 XX
1 BA
34
3 BB
42
96
37
END DATA.
LIST.

```

- Data are in a file with header records that indicate the type of record and the number of individual records that follow. The number of records following each header record varies. For example, the 1 in the first column of the first header record (AC) says that only one individual record (91) follows. The 2 in the first column of the second header record (CC) says that two individual records (35 and 43) follow. The next header record has no individual records, indicated by the 0 in column 1, and so on.
- The first `DATA LIST` reads the expected number of individual records for each header record into temporary variable `#RECS`. `#RECS` is then used as the terminal value in the indexing variable to read the correct number of individual records using the second `DATA LIST`.
- The variables `HEAD1` and `HEAD2` contain the information in columns 3 and 4, respectively, in the header records. The `LEAVE` command retains `HEAD1` and `HEAD2` so that this information can be spread to the individual records.

- The variable *INDIV* is the information from the individual record. *INDIV* is combined with *#RECS*, *HEAD1*, and *HEAD2* to create the new case. Notice in the output from the `PRINT` command below that no case is created for the header record with 0 for *#RECS*.
- `END CASE` passes each case out of the input program to the `LIST` command. Without `END CASE`, the `PRINT` command would still display the cases because it is inside the loop. However, only one (the last) case per header record would pass out of the input program. The outcome for `LIST` will be quite different.

Figure 77-3

*PRINT output*

```

1 A C 9.1
2 C C 3.5
2 C C 4.3
1 B A 3.4
3 B B 4.2
3 B B 9.6
3 B B 3.7

```

Figure 77-4

*LIST output when END CASE is specified*

```

HEAD1 HEAD2 INDIV
A      C      9.1
C      C      3.5
C      C      4.3
B      A      3.4
B      B      4.2
B      B      9.6
B      B      3.7

```

Figure 77-5

*LIST output when END CASE is not specified*

```

HEAD1 HEAD2 INDIV
A      C      9.1
C      C      4.3
X      X      .
B      A      3.4
B      B      3.7

```

# END FILE

```
END FILE
```

## **Example**

```
INPUT PROGRAM.  
DATA LIST FILE=PRICES /YEAR 1-4 QUARTER 6 PRICE 8-12(2).  
DO IF (YEAR GE 1881). /*Stop reading before 1881  
END FILE.  
END IF.  
END INPUT PROGRAM.
```

## **Overview**

END FILE is used in an INPUT PROGRAM–END INPUT PROGRAM structure to tell the program to stop reading data before it actually encounters the end of the file. END FILE can be used with END CASE to concatenate raw data files by causing the program to delay end-of-file processing until it has read multiple data files. END FILE can also be used with LOOP and END CASE to generate data without any data input.

### **Basic Specification**

The basic specification is simply END FILE. There are no additional specifications. The end of file is defined according to the conditions specified for END FILE in the input program.

### **Syntax Rules**

- END FILE is available only within an INPUT PROGRAM structure.
- Only one END FILE command can be executed per input program. However, multiple END FILE commands can be specified within a conditional structure in the input program.

### **Operations**

- When END FILE is encountered, the program stops reading data and puts an end of file in the active dataset it was building. The case that causes the execution of END FILE is not read. To include this case, use the END CASE command before END FILE (see the examples below).
- END FILE has the same effect as the end of the input data file. It terminates the input program (see INPUT PROGRAM–END INPUT PROGRAM).

## **Examples**

### **Stop reading a file based on a data value**

```
*Select cases.  
  
INPUT PROGRAM.  
DATA LIST FILE=PRICES /YEAR 1-4 QUARTER 6 PRICE 8-12(2).
```



```
DO IF (YEAR GE 1881). /*Stop reading before 1881
END FILE.
END IF.

END INPUT PROGRAM.

LIST.
```

- This example assumes that data records are entered chronologically by year. The DO IF–END IF structure specifies an end of file when the first case with a value of 1881 or later for *YEAR* is reached.
- LIST executes the input program and lists cases in the active dataset. The case that causes the end of the file is not included in the active dataset.
- As an alternative to an input program with END FILE, you can use N OF CASES to select cases if you know the exact number of cases. Another alternative is to use SELECT IF to select cases before 1881, but then the program would unnecessarily read the entire input file.

### **END FILE with END CASE**

```
*Select cases but retain the case that causes end-of-file processing.

INPUT PROGRAM.
DATA LIST FILE=PRICES /YEAR 1-4 QUARTER 6 PRICE 8-12(2).

DO IF (YEAR GE 1881). /*Stop reading before 1881 (or at end of file)
END CASE. /*Create case 1881
END FILE.

ELSE.
END CASE. /*Create all other cases
END IF.
END INPUT PROGRAM.

LIST.
```

- The first END CASE command forces the program to retain the case that causes end-of-file processing.
- The second END CASE indicates the end of case for all other cases and passes them out of the input program one at a time. It is required because the first END CASE command causes the program to abandon default end-of-case processing (see END CASE).

# ERASE

```
ERASE FILE='file'
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
ERASE FILE='PRSNL.DAT'.
```

## **Overview**

ERASE removes a file from a disk.

### **Basic Specification**

The basic specification is the keyword `FILE` followed by a file specification enclosed in quotes. The specified file is erased from the disk. The file specification may vary from operating system to operating system.

### **Syntax Rules**

- The keyword `FILE` is required, but the equals sign is optional.
- ERASE allows one file specification only and does not accept wildcard characters. To erase more than one file, specify multiple ERASE commands.
- The file to be erased must be specified in full. ERASE does not recognize any default file extension.

### **Operations**

ERASE deletes the specified file regardless of its type. No message is displayed unless the command cannot be executed. Use ERASE with caution.

## **Examples**

```
ERASE FILE '/data/textdata.txt'.
```

# EXAMINE

EXAMINE is available in the Statistics Base option.

```
EXAMINE VARIABLES=varlist [[BY varlist] [varname BY varname]]

[/COMPARE={GROUPS** }]
           {VARIABLES}

[/{TOTAL**}]
           {NOTOTAL}

[/ID={case number**}]
           {varname      }

[/PERCENTILES [[({5,10,25,50,75,90,95})={(HAVERAGE )} [NONE]]]
              {value list      } {W(AVERAGE )}
                                          {R(OUND )}
                                          {A(EMPIRICAL)}
                                          {E(MPIRICAL )}

[/PLOT={(STEMLEAF**) [BOXPLOT**] [NPLOT] [SPREADLEVEL(n)] [HISTOGRAM]}]

           [{ALL }]
           {NONE}

[/STATISTICS={(DESCRIPTIVES**) [EXTREME({5)}]}]
           [{ALL }]
           {NONE}

[/CINTERVAL {95**}]
           {n }

[/MESTIMATOR={(NONE**)}]
           {ALL }

           [HUBER({1.339)}] [ANDREW({1.34)}]
           {c } {c }

           [HAMPEL({1.7,3.4,8.5)}]
           {a ,b ,c }

           [TUKEY({4.685)}]
           {c }

[/MISSING={(LISTWISE**)}] [(EXCLUDE**)] [(NOREPORT**)]
           {PAIRWISE } {INCLUDE } {REPORT }
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Examples

```
EXAMINE VARIABLES=ENGFSIZE,COST.
```

```
EXAMINE VARIABLES=MIPERVAL BY MODEL, MODEL BY CYLINDERS.
```

## Overview

EXAMINE provides stem-and-leaf plots, histograms, boxplots, normal plots, robust estimates of location, tests of normality, and other descriptive statistics. Separate analyses can be obtained for subgroups of cases.

### Options

**Cells.** You can subdivide cases into cells based on their values for grouping (factor) variables using the `BY` keyword on the `VARIABLES` subcommand.

**Output.** You can control the display of output using the `COMPARE` subcommand. You can specify the computational method and break points for percentiles with the `PERCENTILES` subcommand, and you can assign a variable to be used for labeling outliers on the `ID` subcommand.

**Plots.** You can request stem-and-leaf plots, histograms, vertical boxplots, spread-versus-level plots with Levene tests for homogeneity of variance, and normal and detrended probability plots with tests for normality. These plots are available through the `PLOT` subcommand.

**Statistics.** You can request univariate statistical output with the `STATISTICS` subcommand and maximum-likelihood estimators with the `MESTIMATORS` subcommand.

### Basic Specification

- The basic specification is `VARIABLES` and at least one dependent variable.
- The default output includes a Descriptives table displaying univariate statistics (mean, median, standard deviation, standard error, variance, kurtosis, kurtosis standard error, skewness, skewness standard error, sum, interquartile range (IQR), range, minimum, maximum, and 5% trimmed mean), a vertical boxplot, and a stem-and-leaf plot. Outliers are labeled on the boxplot with the system variable `$CASENUM`.

### Subcommand Order

Subcommands can be named in any order.

### Limitations

- When string variables are used as factors, only the first eight characters are used to form cells. String variables cannot be specified as dependent variables.
- When more than eight crossed factors (for example, A, B, ... in the specification Y by A by B by ...) are specified, the command is not executed.

## Examples

### Univariate statistics

```
EXAMINE VARIABLES=ENGSIZE, COST.
```

- `ENGSIZE` and `COST` are the dependent variables.

- EXAMINE produces univariate statistics for *ENG*SIZE and *COST* in the Descriptives table and a vertical boxplot and a stem-and-leaf plot for each variable.

### **Descriptive statistics across groups**

```
EXAMINE VARIABLES=toxin BY yield
/PLOT BOXPLOT STEMLEAF
/COMPARE GROUP
/STATISTICS DESCRIPTIVES
/CINTERVAL 95
/MISSING LISTWISE
/NOTOTAL.
```

- The procedure examines values of the variable *toxin* by values of the variable *yield*.
- All other options are set to their default values.

### **Exploring distributions**

```
EXAMINE VARIABLES=temp BY batch /ID= labrunid
/PLOT BOXPLOT STEMLEAF NPLOT
/COMPARE GROUP
/MESTIMATORS HUBER(1.339) ANDREW(1.34) HAMPEL(1.7,3.4,8.5) TUKEY(4.685)
/STATISTICS DESCRIPTIVES EXTREME
/CINTERVAL 95
/MISSING LISTWISE
/NOTOTAL.
```

- The procedure examines values of *temp* by values of *batch*.
- The *ID* subcommand specifies that cases in the output should be labeled by values of *labrunid*.
- The *PLOT* subcommand requests boxplots, stem-and-leaf plots, and normality plots.
- The *MESTIMATORS* subcommand requests Huber's, Andrews', Hampel's, and Tukey's robust estimators of location with default weights.
- The *STATISTICS* subcommand requests a descriptive statistics table and an extreme values table with the 5 largest and 5 smallest values.
- All other options are set to their default values.

## **VARIABLES Subcommand**

*VARIABLES* specifies the dependent variables and the cells. The dependent variables are specified first, followed by the keyword *BY* and the variables that define the cells. Repeated models on the same *EXAMINE* are discarded.

- To create cells defined by the combination of values of two or more factors, specify the factor names separated by the keyword *BY*.

**Caution.** Large amounts of output can be produced if many cells are specified. If there are many factors or if the factors have many values, *EXAMINE* will produce a large number of separate analyses.

### **Example**

```
EXAMINE VARIABLES=SALARY,YRSEDUC BY RACE,SEX,DEPT,RACE BY SEX.
```

- *SALARY* and *YRSEDUC* are dependent variables.
- Cells are formed first for the values of *SALARY* and *YRSEDUC* individually, and then each by values for *RACE*, *SEX*, *DEPT*, and the combination of *RACE* and *SEX*.
- By default, EXAMINE produces Descriptives tables, stem-and-leaf plots, and boxplots.

## COMPARE Subcommand

COMPARE controls how boxplots are displayed. This subcommand is most useful if there is more than one dependent variable and at least one factor in the design.

<b>GROUPS</b>	<i>For each dependent variable, boxplots for all cells are displayed together. With this display, comparisons across cells for a single dependent variable are easily made. This is the default.</i>
<b>VARIABLES</b>	<i>For each cell, boxplots for all dependent variables are displayed together. With this display, comparisons of several dependent variables are easily made. This is useful in situations where the dependent variables are repeated measures of the same variable (see the following example) or have similar scales, or when the dependent variable has very different values for different cells, and plotting all cells on the same scale would cause information to be lost.</i>

### Example

```
EXAMINE VARIABLES=GPA1 GPA2 GPA3 GPA4 BY MAJOR /COMPARE=VARIABLES.
```

- The four GPA variables are summarized for each value of *MAJOR*.
- COMPARE=VARIABLES groups the boxplots for the four GPA variables together for each value of *MAJOR*.

### Example

```
EXAMINE VARIABLES=GPA1 GPA2 GPA3 GPA4 BY MAJOR /COMPARE=GROUPS.
```

- COMPARE=GROUPS groups the boxplots for *GPA1* for all majors together, followed by boxplots for *GPA2* for all majors, and so on.

## TOTAL and NOTOTAL Subcommands

TOTAL and NOTOTAL control the amount of output produced by EXAMINE when factor variables are specified.

- TOTAL is the default. By default, or when TOTAL is specified, EXAMINE produces statistics and plots for each dependent variable overall and for each cell specified by the factor variables.
- NOTOTAL suppresses overall statistics and plots.
- TOTAL and NOTOTAL are alternatives.
- NOTOTAL is ignored when the VARIABLES subcommand does not specify factor variables.

## ID Subcommand

ID assigns a variable from the active dataset to identify the cases in the output. By default the case number is used for labeling outliers and extreme cases in boxplots.

- The identification variable can be either string or numeric. If it is numeric, value labels are used to label cases. If no value labels exist, the values are used.
- Only one identification variable can be specified.

### Example

```
EXAMINE VARIABLES=SALARY BY RACE BY SEX /ID=LASTNAME.
```

- ID displays the value of *LASTNAME* for outliers and extreme cases in the boxplots.

## PERCENTILES Subcommand

PERCENTILES displays the Percentiles table. If PERCENTILES is omitted, no percentiles are produced. If PERCENTILES is specified without keywords, HAVERAGE is used with default break points of 5, 10, 25, 50, 75, 90, and 95.

- Values for break points are specified in parentheses following the subcommand. EXAMINE displays up to six decimal places for user-specified percentile values.
- The method keywords follow the specifications for break points.

For each of the following methods of percentile calculation,  $w$  is the sum of the weights for all nonmissing cases,  $p$  is the specified percentile divided by 100, and  $X_i$  is the value of the  $i$ th case (cases are assumed to be ranked in ascending order). For details on the specific formulas used, see the algorithms documentation included on the installation CD.

<b>HAVERAGE</b>	Weighted average at $X_{(w+1)p}$ . The percentile value is the weighted average of $X_i$ and $X_{i+1}$ , where $i$ is the integer part of $(w+1)p$ . This is the default if PERCENTILES is specified without a keyword.
<b>WAVERAGE</b>	Weighted average at $X_{wp}$ . The percentile value is the weighted average of $X_i$ and $X_{(i+1)}$ , where $i$ is the integer portion of $wp$ .
<b>ROUND</b>	<i>Observation closest to</i> $wp$ . The percentile value is $X_i$ or $X_{i+1}$ , depending upon whether $i$ or $i+1$ is "closer" to $wp$ .
<b>EMPIRICAL</b>	<i>Empirical distribution function</i> . The percentile value is $X_i$ , where $i$ is equal to $wp$ rounded up to the next integer.
<b>AEMPIRICAL</b>	<i>Empirical distribution with averaging</i> . This is equivalent to EMPIRICAL, except when $i=wp$ , in which case the percentile value is the average of $X_i$ and $X_{i+1}$ .
<b>NONE</b>	<i>Suppress percentile output</i> . This is the default if PERCENTILES is omitted.

### Example

```
EXAMINE VARIABLE=SALARY /PERCENTILES(10,50,90)=EMPIRICAL.
```

- PERCENTILES produces the 10th, 50th, and 90th percentiles for the dependent variable *SALARY* using the EMPIRICAL distribution function.

## PLOT Subcommand

PLOT controls plot output. The default is a vertical boxplot and a stem-and-leaf plot for each dependent variable for each cell in the model.

- Spread-versus-level plots can be produced only if there is at least one factor variable on the VARIABLES subcommand. If you request a spread-versus-level plot and there are no factor variables, the program issues a warning and no spread-versus-level plot is produced.
- If you specify the PLOT subcommand, only those plots explicitly requested are produced.

<b>BOXPLOT</b>	<i>Vertical boxplot.</i> The boundaries of the box are Tukey's hinges. The median is identified by a line inside the box. The length of the box is the interquartile range (IQR) computed from Tukey's hinges. Values more than three IQR's from the end of a box are labeled as extreme, denoted with an asterisk (*). Values more than 1.5 IQR's but less than 3 IQR's from the end of the box are labeled as outliers (o).
<b>STEMLEAF</b>	<i>Stem-and-leaf plot.</i> In a stem-and-leaf plot, each observed value is divided into two components—leading digits (stem) and trailing digits (leaf).
<b>HISTOGRAM</b>	<i>Histogram.</i>
<b>SPREADLEVEL(n)</b>	<i>Spread-versus-level plot with the Test of Homogeneity of Variance table.</i> If the keyword appears alone, the natural logs of the interquartile ranges are plotted against the natural logs of the medians for all cells. If a power for transforming the data ( <i>n</i> ) is given, the IQR and median of the transformed data are plotted. If 0 is specified for <i>n</i> , a natural log transformation of the data is done. The slope of the regression line and Levene tests for homogeneity of variance are also displayed. The Levene tests are based on the original data if no transformation is specified and on the transformed data if a transformation is requested.
<b>NPLOT</b>	<i>Normal and detrended Q-Q plots with the Tests of Normality table presenting Shapiro-Wilk's statistic and a Kolmogorov-Smirnov statistic with a Lilliefors significance level for testing normality.</i> If non-integer weights are specified, the Shapiro-Wilk's statistic is calculated when the weighted sample size lies between 3 and 50. For no weights or integer weights, the statistic is calculated when the weighted sample size lies between 3 and 5,000.
<b>ALL</b>	<i>All available plots.</i>
<b>NONE</b>	<i>No plots.</i>

### Example

```
EXAMINE VARIABLES=CYCLE BY TREATMNT /PLOT=NPLOT.
```

- PLOT produces normal and detrended Q-Q plots for each value of *TREATMNT* and a Tests of Normality table.

### Example

```
EXAMINE VARIABLES=CYCLE BY TREATMNT /PLOT=SPREADLEVEL(.5).
```

- PLOT produces a spread-versus-level plot of the medians and interquartile ranges of the square root of *CYCLE*. Each point on the plot represents one of the *TREATMNT* groups.
- A Test of Homogeneity of Variance table displays Levene statistics.



**Example**

```
EXAMINE VARIABLES=CYCLE BY TREATMNT /PLOT=SPREADLEVEL(0).
```

- PLOT generates a spread-versus-level plot of the medians and interquartile ranges of the natural logs of *CYCLE* for each *TREATMENT* group.
- A Test of Homogeneity of Variance table displays Levene statistics.

**Example**

```
EXAMINE VARIABLES=CYCLE BY TREATMNT /PLOT=SPREADLEVEL.
```

- PLOT generates a spread-versus-level plot of the natural logs of the medians and interquartile ranges of *CYCLE* for each *TREATMNT* group.
- A Test of Homogeneity of Variance table displays Levene statistics.

**STATISTICS Subcommand**

STATISTICS requests univariate statistics and determines how many extreme values are displayed. DESCRIPTIVES is the default. If you specify keywords on STATISTICS, only the requested statistics are displayed.

<b>DESCRIPTIVES</b>	<i>Display the Descriptives table showing univariate statistics (the mean, median, 5% trimmed mean, standard error, variance, standard deviation, minimum, maximum, range, interquartile range, skewness, skewness standard error, kurtosis, and kurtosis standard error). This is the default.</i>
<b>EXTREME(n)</b>	<i>Display the Extreme Values table presenting cases with the n largest and n smallest values. If n is omitted, the five largest and five smallest values are displayed. Extreme cases are labeled with their values for the identification variable if the ID subcommand is used or with their values for the system variable \$CASENUM if ID is not specified.</i>
<b>ALL</b>	<i>Display the Descriptives and Extreme Values tables.</i>
<b>NONE</b>	<i>Display neither the Descriptives nor the Extreme Values tables.</i>

**Example**

```
EXAMINE VARIABLE=FAILTIME /ID=BRAND
  /STATISTICS=EXTREME(10) /PLOT=NONE.
```

- STATISTICS identifies the cases with the 10 lowest and 10 highest values for *FAILTIME*. These cases are labeled with the first characters of their values for the variable *BRAND*. The Descriptives table is not displayed.

**CINTERVAL Subcommand**

CINTERVAL controls the confidence level when the default DESCRIPTIVES statistics is displayed. CINTERVAL has a default value of 95.

- You can specify a CINTERVAL value (*n*) between 50 and 99.99 inclusive. If the value you specify is out of range, the program issues a warning and uses the default 95% intervals.

- If you specify a keyword on `STATISTICS` subcommand that turns off the default `DESCRIPTIVES`, the `CINTERVAL` subcommand is ignored.
- The confidence interval appears in the output with the label *n% CI for Mean*, followed by the confidence interval in parentheses. For example, **95% CI for Mean (.0001,.00013)**

The *n* in the label shows up to six decimal places. That is, input `/CINTERVAL 95` displays as *95% CI* while input `/CINTERVAL 95.975` displays as *95.975% CI*.

## **MESTIMATORS Subcommand**

M-estimators are robust maximum-likelihood estimators of location. Four M-estimators are available for display in the M-Estimators table. They differ in the weights they apply to the cases. `MESTIMATORS` with no keywords produces Huber's M-estimator with  $c=1.339$ ; Andrews' wave with  $c=1.34\pi$ ; Hampel's M-estimator with  $a=1.7$ ,  $b=3.4$ , and  $c=8.5$ ; and Tukey's biweight with  $c=4.685$ .

<b>HUBER(c)</b>	<i>Huber's M-estimator.</i> The value of weighting constant $c$ can be specified in parentheses following the keyword. The default is $c=1.339$ .
<b>ANDREW(c)</b>	<i>Andrews' wave estimator.</i> The value of weighting constant $c$ can be specified in parentheses following the keyword. Constants are multiplied by $\pi$ . The default is $1.34\pi$ .
<b>HAMPEL(a,b,c)</b>	<i>Hampel's M-estimator.</i> The values of weighting constants $a$ , $b$ , and $c$ can be specified in order in parentheses following the keyword. The default values are $a=1.7$ , $b=3.4$ , and $c=8.5$ .
<b>TUKEY(c)</b>	<i>Tukey's biweight estimator.</i> The value of weighting constant $c$ can be specified in parentheses following the keyword. The default is $c=4.685$ .
<b>ALL</b>	<i>All four above M-estimators.</i> This is the default when <code>MESTIMATORS</code> is specified with no keyword. The default values for weighting constants are used.
<b>NONE</b>	<i>No M-estimators.</i> This is the default if <code>MESTIMATORS</code> is omitted.

### **Example**

```
EXAMINE VARIABLE=CASTTEST /MESTIMATORS.
```

- `MESTIMATORS` generates all four M-estimators computed with the default constants.

### **Example**

```
EXAMINE VARIABLE=CASTTEST /MESTIMATORS=HAMPELS(2,4,8).
```

- `MESTIMATOR` produces Hampel's M-estimator with weighting constants  $a=2$ ,  $b=4$ , and  $c=8$ .

## **MISSING Subcommand**

`MISSING` controls the processing of missing values in the analysis. The default is `LISTWISE`, `EXCLUDE`, and `NOREPORT`.

- `LISTWISE` and `PAIRWISE` are alternatives and apply to all variables. They are modified for dependent variables by `INCLUDE/EXCLUDE` and for factor variables by `REPORT/NOREPORT`.
- `INCLUDE` and `EXCLUDE` are alternatives; they apply only to dependent variables.

- `REPORT` and `NOREPORT` are alternatives; they determine if missing values for factor variables are treated as valid categories.

<b>LISTWISE</b>	<i>Delete cases with missing values listwise.</i> A case with missing values for any dependent variable or any factor in the model specification is excluded from statistics and plots unless modified by <code>INCLUDE</code> or <code>REPORT</code> . This is the default.
<b>PAIRWISE</b>	<i>Delete cases with missing values pairwise.</i> A case is deleted from the analysis only if it has a missing value for the dependent variable or factor being analyzed.
<b>EXCLUDE</b>	<i>Exclude user-missing values.</i> User-missing values and system-missing values for dependent variables are excluded. This is the default.
<b>INCLUDE</b>	<i>Include user-missing values.</i> Only system-missing values for dependent variables are excluded from the analysis.
<b>NOREPORT</b>	<i>Exclude user- and system-missing values for factor variables.</i> This is the default.
<b>REPORT</b>	<i>Include user- and system-missing values for factor variables.</i> User- and system-missing values for factors are treated as valid categories and are labeled as missing.

### Example

```
EXAMINE VARIABLES=RAINFALL MEANTEMP BY REGION.
```

- `MISSING` is not specified and the default is used. Any case with a user- or system-missing value for `RAINFALL`, `MEANTEMP`, or `REGION` is excluded from the analysis and display.

### Example

```
EXAMINE VARIABLES=RAINFALL MEANTEMP BY REGION
/MISSING=PAIRWISE.
```

- Only cases with missing values for `RAINFALL` are excluded from the analysis of `RAINFALL`, and only cases with missing values for `MEANTEMP` are excluded from the analysis of `MEANTEMP`. Missing values for `REGION` are not used.

### Example

```
EXAMINE VARIABLES=RAINFALL MEANTEMP BY REGION
/MISSING=REPORT.
```

- Missing values for `REGION` are considered valid categories and are labeled as missing.

## References

- Hoaglin, D. C., F. Mosteller, and J. W. Tukey. 1983. *Understanding robust and exploratory data analysis*. New York: John Wiley and Sons.
- Hoaglin, D. C., F. Mosteller, and J. W. Tukey. 1985. *Exploring data tables, trends, and shapes*. New York: John Wiley and Sons.
- Tukey, J. W. 1977. *Exploratory data analysis*. Reading, MA: Addison-Wesley.

Velleman, P. F., and D. C. Hoaglin. 1981. *Applications, basics, and computing of exploratory data analysis*. Boston, Mass.: Duxbury Press.

# EXECUTE

EXECUTE.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Overview

EXECUTE forces the data to be read and executes the transformations that precede it in the command sequence.

### Basic Specification

The basic specification is simply the command keyword. EXECUTE has no additional specifications.

### Operations

- EXECUTE causes the data to be read but has no other influence on the session.
- EXECUTE is designed for use with transformation commands and facilities such as ADD FILES, MATCH FILES, UPDATE, PRINT, and WRITE, which do not read data and are not executed unless followed by a data-reading procedure.

## Examples

```
DATA LIST FILE=RAWDATA / 1 LNAME 1-13 (A) FNAME 15-24 (A)
      MMAIDENL 40-55.
VAR LABELS MMAIDENL 'MOTHER'S MAIDEN NAME'.
DO IF (MMAIDENL EQ 'Smith').
WRITE OUTFILE=SMITHS/LNAME FNAME.
END IF.
EXECUTE.
```

- This example writes the last and first names of all people whose mother's maiden name was Smith to the data file *SMITHS*.
- DO IF-END IF and WRITE do not read data and are executed only when data are read for a procedure. Because there is no procedure in this session, EXECUTE is used to read the data and execute all of the preceding transformation commands. Otherwise, the commands would not be executed.

# EXPORT

```
EXPORT OUTFILE='file'  
  
  [/TYPE={COMM**}  
    {TAPE  }]  
  
  [/UNSELECTED=[{RETAIN}]  
    {DELETE}]  
  
  [/KEEP={ALL**  }] [/DROP=varlist]  
    {varlist}  
  
  [/RENAME=(old varnames=new varnames)...]  
  
  [/MAP]  
  
  [/DIGITS=n]
```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
EXPORT OUTFILE="/data/newdata.por"  
  /RENAME=(V1 TO V3=ID, SEX, AGE) /MAP.
```

## Overview

EXPORT produces a portable data file. A portable data file is a data file created used to transport data between different types of computers and operating systems, or other software using the same portable file format. A portable file contains all of the data and dictionary information stored in the active dataset from which it was created.

EXPORT is similar to the SAVE command. It can occur in the same position in the command sequence as the SAVE command and saves the active dataset. The file includes the results of all permanent transformations and any temporary transformations made just prior to the EXPORT command. The active dataset is unchanged after the EXPORT command.

- In most cases, saving data in portable format is no longer necessary, since IBM® SPSS® Statistics data files should be platform/operating system independent.
- To export data in external data formats (e.g., Excel, SAS, Stata, CSV, tab-delimited), use [SAVE TRANSLATE](#).

## Options

**Format.** You can control the format of the portable file using the TYPE subcommand.

**Variables.** You can save a subset of variables from the active dataset and rename the variables using the DROP, KEEP, and RENAME subcommands. You can also produce a record of all variables and their names on the exported file with the MAP subcommand.

**Precision.** You can specify the number of decimal digits of precision for the values of all numeric variables on the `DIGITS` subcommand.

### **Basic Specification**

The basic specification is the `OUTFILE` subcommand with a file specification. All variables from the active dataset are written to the portable file, with variable names, variable and value labels, missing-value flags, and print and write formats.

### **Subcommand Order**

Subcommands can be named in any order.

### **Operations**

- Portable files are written with 80-character record lengths.
- Portable files may contain some unprintable characters.
- The active dataset is still available for transformations and procedures after the portable file is created.
- The system variables `$CASENUM` and `$DATE` are assigned when the file is read by `IMPORT`.
- If the `WEIGHT` command is used before `EXPORT`, the weighting variable is included in the portable file.
- Variable names that exceed eight bytes are converted to unique eight-byte names—for example, `mylongrootname1`, `mylongrootname2`, and `mylongrootname3` would be converted to `mylongro`, `mylong_2`, and `mylong_3`, respectively.

### **Limitations**

The `EXPORT` command is not supported in Unicode mode. For more information, see [SET command](#), [UNICODE subcommand](#).

## **Examples**

```
EXPORT OUTFILE="/newdata.por"  
  /RENAME=(V1 TO V3=ID,SEX,AGE) /MAP.
```

- The portable file is written to `newdata.por`.
- The variables `V1`, `V2`, and `V3` are renamed `ID`, `SEX`, and `AGE` in the portable file. Their names remain `V1`, `V2`, and `V3` in the active dataset. None of the other variables written to the portable file are renamed.
- `MAP` requests a display of the variables in the portable file.

## **Methods of Transporting Portable Files**

Portable files can be transported on magnetic tape or by a communications program.

## ***Magnetic Tape***

Before transporting files on a magnetic tape, make sure the receiving computer can read the tape being sent. The following tape specifications must be known before you write the portable file on the tape:

- **Number of tracks.** Either 7 or 9.
- **Tape density.** 200, 556, 800, 1600, or 6250 bits per inch (BPI).
- **Parity.** Even or odd. This must be known only when writing a 7-track tape.
- **Tape labeling.** Labeled or unlabeled. Check whether the site can use tape labels. Also make sure that the site has the ability to read multivolume tape files if the file being written uses more than one tape.
- **Blocksize.** The maximum blocksize the receiving computer can accept.

A tape written with the following characteristics can be read by most computers: 9 track, 1600 BPI, unlabeled, and a blocksize of 3200 characters. However, there is no guarantee that a tape written with these characteristics can be read successfully. The best policy is to know the requirements of the receiving computer ahead of time.

The following advice may help ensure successful file transfers by magnetic tape:

- Unless you are certain that the receiving computer can read labels, prepare an unlabeled tape.
- Make sure the record length of 80 is not changed.
- Do not use a separate character translation program, especially ASCII/EBCDIC translations. `EXPORT/IMPORT` takes care of this for you.
- Make sure the same blocking factor is used when writing and reading the tape. A blocksize of 3200 is frequently a good choice.
- If possible, write the portable file directly to tape to avoid possible interference from copy programs. Read the file directly from the tape for the same reason.
- Use the `INFO LOCAL` command to find out about using the program on your particular computer and operating system. `INFO LOCAL` generally includes additional information about reading and writing portable files.

## ***Communications Programs***

Transmission of a portable file by a communications program may not be possible if the program misinterprets any characters in the file as control characters (for example, as a line feed, carriage return, or end of transmission). This can be prevented by specifying `TYPE=COMM` on `EXPORT`. This specification replaces each control character with the character 0. The affected control characters are in positions 0–60 of the `IMPORT/EXPORT` character set. [For more information, see the topic `IMPORT/EXPORT` Character Sets on p. 2189.](#)

The line length that the communications program uses must be set to 80 to match the 80-character record length of portable files. A transmitted file must be checked for blank lines or special characters inserted by the communications program. These must be edited out prior to reading the file with the `IMPORT` command.



## Character Translation

Portable files are character files, not binary files, and they have 80-character records so they can be transmitted over data links. A receiving computer may not use the same character set as the computer where the portable file was written. When it imports a portable file, the program translates characters in the file to the character set used by the receiving computer. Depending on the character set in use, some characters in labels and in string data may be lost in the translation. For example, if a file is transported from a computer using a seven-bit ASCII character set to a computer using a six-bit ASCII character set, some characters in the file may have no matching characters in six-bit ASCII. For a character that has no match, the program generates an appropriate nonprintable character (the null character in most cases).

For a table of the character-set translations available with `IMPORT` and `EXPORT`, refer to Appendix B. A blank in a column of the table means that there is no matching character for that character set and an appropriate nonprintable character will be generated when you import a file.

## OUTFILE Subcommand

`OUTFILE` specifies the portable file. `OUTFILE` is the only required subcommand on `EXPORT`.

## TYPE Subcommand

`TYPE` indicates whether the portable file should be formatted for magnetic tape or for a communications program. You can specify either `COMM` or `TAPE`. [For more information, see the topic Methods of Transporting Portable Files on p. 697.](#)

<b>COMM</b>	<i>Transport portable files by a communications program.</i> When <code>COMM</code> is specified on <code>TYPE</code> , the program removes all control characters and replaces them with the character 0. This is the default.
<b>TAPE</b>	<i>Transport portable files on magnetic tape.</i>

### Example

```
EXPORT TYPE=TAPE /OUTFILE=HUBOUT.
```

- File `HUBOUT` is saved as a tape-formatted portable file.

## UNSELECTED Subcommand

`UNSELECTED` determines whether cases excluded on a previous `FILTER` or `USE` command are to be retained or deleted in the data file. The default is `RETAIN`. The `UNSELECTED` subcommand has no effect when the active dataset does not contain unselected cases.

<b>RETAIN</b>	<i>Retain the unselected cases.</i> All cases in the active dataset are saved. This is the default when <code>UNSELECTED</code> is specified by itself.
<b>DELETE</b>	<i>Delete the unselected cases.</i> Only cases that meet the <code>FILTER</code> or <code>USE</code> criteria are saved in the data file.

## ***DROP and KEEP Subcommands***

DROP and KEEP save a subset of variables in the portable file.

- DROP excludes a variable or list of variables from the portable file. All variables not named are included in the portable file.
- KEEP includes a variable or list of variables in the portable file. All variables not named are excluded.
- Variables can be specified on DROP and KEEP in any order. With the DROP subcommand, the order of variables in the portable file is the same as their order in the active dataset. With the KEEP subcommand, the order of variables in the portable file is the order in which they are named on KEEP. Thus, KEEP can be used to reorder variables in the portable file.
- Both DROP and KEEP can be used on the same EXPORT command; the effect is cumulative. If you specify a variable already named on a previous DROP or one not named on a previous KEEP, the variable is considered nonexistent and the program displays an error message. The command is aborted and no portable file is saved.

### ***Example***

```
EXPORT OUTFILE=NEWSUM /DROP=DEPT TO DIVISION.
```

- The portable file is written to file *NEWSUM*. Variables between and including *DEPT* and *DIVISION* in the active dataset are excluded from the portable file.
- All other variables are saved in the portable file.

## ***RENAME Subcommand***

RENAME renames variables being written to the portable file. The renamed variables retain their original variable and value labels, missing-value flags, and print formats. The names of the variables are not changed in the active dataset.

- To rename a variable, specify the name of the variable in the active dataset, an equals sign, and the new name.
- A variable list can be specified on both sides of the equals sign. The number of variables on both sides must be the same, and the entire specification must be enclosed in parentheses.
- The keyword TO can be used for both variable lists.
- If you specify a renamed variable on a subsequent DROP or KEEP subcommand, the new variable name must be used.

### ***Example***

```
EXPORT OUTFILE=NEWSUM /DROP=DEPT TO DIVISION  
/RENAME=(NAME, WAGE=LNAME, SALARY) .
```

- RENAME renames *NAME* and *WAGE* to *LNAME* and *SALARY*.
- *LNAME* and *SALARY* retain the variable and value labels, missing-value flags, and print formats assigned to *NAME* and *WAGE*.

## MAP Subcommand

MAP displays any changes that have been specified by the RENAME, DROP, or KEEP subcommands.

- MAP can be specified as often as desired.
- Each MAP subcommand maps the results of subcommands that precede it; results of subcommands that follow it are not mapped. When MAP is specified last, it also produces a description of the portable file.

### Example

```
EXPORT OUTFILE=NEWSUM /DROP=DEPT TO DIVISION /MAP  
/RENAME NAME=LNAME WAGE=SALARY /MAP.
```

- The first MAP subcommand produces a listing of the variables in the file after DROP has dropped the specified variables.
- RENAME renames *NAME* and *WAGE*.
- The second MAP subcommand shows the variables in the file after renaming. Since this is the last subcommand, the listing will show the variables as they are written in the portable file.

## DIGITS Subcommand

DIGITS specifies the degree of precision for all noninteger numeric values written to the portable file.

- DIGITS has the general form DIGITS=*n*, where *n* is the number of digits of precision.
- DIGITS applies to all numbers for which rounding is required.
- Different degrees of precision *cannot* be specified for different variables. Thus, DIGITS should be set according to the requirements of the variable that needs the most precision.
- Default precision methods used by EXPORT work perfectly for integers that are not too large and for fractions whose denominators are products of 2, 3, and 5 (all decimals, quarters, eighths, sixteenths, thirds, thirtieths, sixtieths, and so forth.) For other fractions and for integers too large to be represented exactly in the active dataset (usually more than 9 digits, often 15 or more), the representation used in the active dataset contains some error already, so no exact way of sending these numbers is possible. The program sends enough digits to get very close. The number of digits sent in these cases depends on the originating computer: on mainframe IBM versions of the program, it is the equivalent of 13 decimal digits (integer and fractional parts combined). If many numbers on a file require this level of precision, the file can grow quite large. If you do not need the full default precision, you can save some space in the portable file by using the DIGITS subcommand.

### Example

```
EXPORT OUTFILE=NEWSUM /DROP=DEPT TO DIVISION /MAP /DIGITS=4.
```

- DIGITS guarantees the accuracy of values to four significant digits. For example, 12.34567890876 will be rounded to 12.35.

# EXTENSION

```
EXTENSION ACTION={ADD** }  
                {REMOVE}  
  
/SPECIFICATION COMMAND=file.
```

\*\* Default if the subcommand or keyword is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- Command introduced.

## **Example**

```
EXTENSION  
/SPECIFICATION COMMAND="c:\mypls.xml"
```

## **Overview**

EXTENSION allows you to add user-created “extension” commands to the command table, enabling the use of “native” command syntax to drive functions written in an external programming language.

## **Basic Specification**

The minimum specification is the SPECIFICATION subcommand with a COMMAND file specified.

## **ACTION Keyword**

The ACTION keyword specifies the action to be taken on the command table.

- |               |  |
|---------------|--|
| <b>ADD</b>    | <i>Add command name to command table.</i> This is the default. Note that ADD will replace a built-in command of the same name as well as an existing extension command of the same name, regardless of the languages in which the extension commands are written. It will not replace an abbreviated form of a built-in command unless the name matches the abbreviation. For example, an extension command named FREQUENCIES will replace the built-in FREQUENCIES command but the abbreviation FREQ will still run the built-in command. |
| <b>REMOVE</b> | <i>Remove command name from command table.</i> Any built-in command replaced by the specified extension command will be restored when the extension command is removed.  |

For information on creating extension commands, see the article “*Writing IBM® SPSS® Statistics Extension Commands*”, available from Developer Central at <http://www.spss.com/devcentral>. You may also want to consult the chapter on Extension Commands in *Programming and Data Management for SPSS Statistics*, available in PDF from [http://www.spss.com/spss/data\\_management\\_book.htm](http://www.spss.com/spss/data_management_book.htm).

## Examples

### **Adding a New Command**

```
EXTENSION  
  /SPECIFICATION COMMAND="<installdir>/extensions/plscommand.xml".
```

The `SPECIFICATION` subcommand specifies that the syntax diagram for the extension command is defined by *plscommand.xml* in the `\extensions` subdirectory of the main installation directory (replace `<installdir>` with the path to the installation directory on your system). The default action is for the command name provided in the XML file to be added to the command table or replace an existing command name.

### **Removing a Command**

```
EXTENSION ACTION=REMOVE  
  /SPECIFICATION COMMAND="<installdir>/extensions/plscommand.xml".
```

The `ACTION` keyword specifies that the command name provided in *plscommand.xml* should be removed from the command table.

## **SPECIFICATION Subcommand**

The `SPECIFICATION` subcommand allows you to specify the location of the XML file that defines the syntax diagram for the extension command.

*Note:* At startup, the system processes all XML files in the `\extensions` subdirectory of the main installation directory as well as in any directories specified in the optional `SPSS_EXTENSIONS_PATH` environment variable. `EXTENSION` is most useful when you have XML files that are not in one of those locations, or want to add a new extension command to an already-running session.

### **COMMAND Keyword**

The `COMMAND` keyword specifies the location of the XML file that defines the syntax diagram for the extension command. This file provides the command name and is used by the universal parser to pass the correct arguments to the extension command.

# FACTOR

FACTOR is available in the Statistics Base option.

```
FACTOR VARIABLES=varlist† [/MISSING={LISTWISE**} [INCLUDE]]
                                {PAIRWISE }
                                {MEANSUB }
                                {DEFAULT** }

[/MATRIX={IN({COR='savfile' | 'dataset'})} [OUT({COR='savfile' | 'dataset'})]]
          {COR=* } {COR=* }
          {COV='savfile' | 'dataset'} {COV='savfile' | 'dataset'}
          {COV=* } {COV=* }
          {FAC='savfile' | 'dataset'} {FAC='savfile' | 'dataset'}
          {FAC=* } {FAC=* }
          {FSC='savfile' | 'dataset'}
          {FSC=* }

[/METHOD = {CORRELATION**}]
           {COVARIANCE }

[/SELECT=varname(value)]

[/ANALYSIS=varlist...]

[/PRINT={DEFAULT**} [INITIAL**] [EXTRACTION**] [ROTATION**]
          [UNIVARIATE] [CORRELATION] [COVARIANCE] [DET] [INV]
          [REPR] [AIC] [KMO] [FSCORE] [SIG] [ALL]]

[/PLOT={EIGEN} [ROTATION [(n1,n2)]]]

[/DIAGONAL={value list}]
           {DEFAULT** }

[/FORMAT={SORT} [BLANK(n)] [DEFAULT**]]

[/CRITERIA={FACTORS(n)} [MINEIGEN({1.0**})] [ITERATE({25**})]
           {n } {n }

          [RCONVERGE({0.0001**})] [{KAISER**}]
          {n } {NOKAISER}

          [ECONVERGE({0.001**})] [DEFAULT**]
          {n }

[/EXTRACTION={PC** }] [/ROTATION={VARIMAX** }]
          {PA1** } {EQUAMAX }
          {PAF } {QUARTIMAX }
          {ALPHA } {OBLIMIN({0})}
          {IMAGE } {n}
          {ULS } {PROMAX({4})}
          {GLS } {n}
          {ML } {NOROTATE }
          {DEFAULT**} {DEFAULT** }

[/SAVE={REG } ({ALL}[rootname])]]
       {BART } {n }
       {AR }
       {DEFAULT}
```

† Omit VARIABLES with matrix input.

\*\*Default if subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 16.0

- Added support for `SET THREADS` and `SET MCACHE`.

### **Example**

```
FACTOR VARIABLES=V1 TO V12.
```

## **Overview**

FACTOR performs factor analysis based either on correlations or covariances and using one of the seven extraction methods. FACTOR also accepts matrix input in the form of correlation matrices, covariance matrices, or factor-loading matrices and can write the matrix materials to a matrix data file.

### **Options**

**Analysis Phase Options.** You can choose to analyze a correlation or covariance matrix using the `METHOD` subcommand. You can select a subset of cases for the analysis phase using the `SELECT` subcommand. You can tailor the statistical display for an analysis using the `PRINT` subcommand. You can sort the output in the factor pattern and structure matrices with the `FORMAT` subcommand. You can also request scree plots and plots of the variables in factor space on the `PLOT` subcommand.

**Extraction Phase Options.** With the `EXTRACTION` subcommand, you can specify one of six extraction methods in addition to the default principal components extraction: principal axis factoring, alpha factoring, image factoring, unweighted least squares, generalized least squares, and maximum likelihood. You can supply initial diagonal values for principal axis factoring on the `DIAGONAL` subcommand. On the `CRITERIA` subcommand, you can alter the default statistical criteria used in the extraction.

**Rotation Phase Options.** You can control the criteria for factor rotation with the `CRITERIA` subcommand. On the `ROTATION` subcommand, you can choose among four rotation methods (equamax, quartimax, promax, and oblimin) in addition to the default varimax rotation, or you can specify no rotation.

**Factor Scores.** You can save factor scores as new variables in the active dataset using any of the three methods available on the `SAVE` subcommand.

**Matrix Input and Output.** With the `MATRIX` subcommand, you can write a correlation matrix, a covariance matrix, or a factor-loading matrix. You can also read matrix materials written either by a previous FACTOR procedure or by a procedure that writes correlation or covariance matrices.

### **Basic Specification**

The basic specification is the `VARIABLES` subcommand with a variable list. FACTOR performs principal components analysis with a varimax rotation on all variables in the analysis using default criteria.

- When matrix materials are used as input, do not specify `VARIABLES`. Use the `ANALYSIS` subcommand to specify a subset of the variables in the matrix.

### **Subcommand Order**

- `METHOD` and `SELECT` can be specified anywhere. `VARIABLES` must be specified before any other subcommands, unless an input matrix is specified. `MISSING` must be specified before `ANALYSIS`.
- The `ANALYSIS`, `EXTRACTION`, `ROTATION`, and `SAVE` subcommands must be specified in the order they are listed here. If you specify these subcommands out of order, you may get unexpected results. For example, if you specify `EXTRACTION` before `ANALYSIS` and `SAVE` before `ROTATION`, `EXTRACTION` and `SAVE` are ignored. If no `EXTRACTION` and `SAVE` subcommands are specified in proper order, the default will be used (that is, `PC` for `EXTRACTION` and no `SAVE`).
- The `FORMAT` subcommand can be specified anywhere after the `VARIABLES` subcommand.
- If an `ANALYSIS` subcommand is present, the statistical display options on `PRINT`, `PLOT`, or `DIAGONAL` must be specified after it. `PRINT`, `PLOT`, and `DIAGONAL` subcommands specified before the `ANALYSIS` subcommand are ignored. If no such commands are specified after the `ANALYSIS` subcommand, the default is used.
- The `CRITERIA` subcommand can be specified anywhere, but applies only to the subcommands that follow. If no `CRITERIA` subcommand is specified before `EXTRACTION` or `ROTATION`, the default criteria for the respective subcommand are used.

### **Example**

```
FACTOR VAR=V1 TO V12  
  /ANALYSIS=V1 TO V8  
  /CRITERIA=FACTORS(3)  
  /EXTRACTION=PAF  
  /ROTATION=QUARTIMAX.
```

- The default `CORRELATION` method is used. `FACTOR` performs a factor analysis of the correlation matrix based on the first eight variables in the active dataset (`V1` to `V8`).
- The procedure extracts three factors using the principal axis method and quartimax rotation.
- `LISTWISE` (the default for `MISSING`) is in effect. Cases with missing values for any one of the variables from `V1` to `V12` are omitted from the analysis. As a result, if you ask for the factor analysis using `VAR=V1 TO V8` and `ANALYSIS=ALL`, the results may be different even though the variables used in the analysis are the same.

### **Syntax Rules**

- Each `FACTOR` procedure performs only one analysis with one extraction and one rotation. Use multiple `FACTOR` commands to perform multiple analyses.
- `VARIABLES` or `MATRIX=IN` can be specified only once. Any other subcommands can be specified multiple times but only the last in proper order takes effect.



**Operations**

- VARIABLES calculates a correlation and a covariance matrix. If SELECT is specified, only the selected cases are used.
- The correlation or covariance matrix (either calculated from the data or read in) is the basis for the factor analysis.
- Factor scores are calculated for all cases (selected and unselected).
- This procedure uses the multithreaded options specified by SET THREADS and SET MCACHE.

**Examples****For data reduction**

```
FACTOR
/VARIABLES type price engine_s horsepow wheelbas width length curb_wgt
fuel_cap mpg /MISSING LISTWISE /ANALYSIS type price engine_s horsepow
wheelbas width length curb_wgt fuel_cap mpg
/PRINT INITIAL EXTRACTION ROTATION FSCORE AIC
/PLOT EIGEN
/CRITERIA MINEIGEN(1) ITERATE(25)
/EXTRACTION PC
/CRITERIA ITERATE(25)
/ROTATION VARIMAX
/METHOD=CORRELATION
/SAVE = REG(3) .
```

- The procedure requests a principal components analysis of the variables *type* through *mpg*.
- The PRINT subcommand requests the initial communalities, extracted communalities, rotated communalities, factor score coefficient matrix, and anti-image covariance and correlation matrices.
- The PLOT subcommand requests a scree plot.
- The ROTATION subcommand requests a Varimax rotation of the solution for easier interpretation of the results.
- The SAVE subcommand requests the factor scores, computed using the regression method, to be saved to the active dataset. Three scores will be added.
- All other options are set to their default values.

**For structure detection**

```
FACTOR
/VARIABLES longmon tollmon equipmon cardmon wiremon multline voice
ebill pager internet callid callwait forward confer /MISSING LISTWISE /ANALYSIS
longmon tollmon equipmon cardmon wiremon multline voice pager
internet callid callwait forward confer ebill
/PRINT INITIAL KMO EXTRACTION ROTATION
/PLOT EIGEN ROTATION
/CRITERIA MINEIGEN(1) ITERATE(25)
/EXTRACTION PAF
/CRITERIA ITERATE(25)
/ROTATION VARIMAX
/METHOD=CORRELATION .
```

- The procedure requests a principal axis factoring of the variables *longmon* through *confer*.

- The `PRINT` subcommand requests the initial communalities, extracted communalities, rotated communalities, and the Kaiser-Meyer-Olkin measure of sampling adequacy and Bartlett's test of sphericity.
- The `PLOT` subcommand requests a scree plot and loadings plot.
- The `ROTATION` subcommand requests a Varimax rotation of the solution for easier interpretation of the results.
- All other options are set to their default values.

## ***VARIABLES Subcommand***

`VARIABLES` names all the variables to be used in the `FACTOR` procedure.

- `VARIABLES` is required except when matrix input is used. When `FACTOR` reads a matrix data file, the `VARIABLES` subcommand cannot be used.
- The specification on `VARIABLES` is a list of numeric variables.
- Keyword `ALL` on `VARIABLES` refers to all variables in the active dataset.
- Only one `VARIABLES` subcommand can be specified, and it must be specified first.

## ***MISSING Subcommand***

`MISSING` controls the treatment of cases with missing values.

- If `MISSING` is omitted or included without specifications, listwise deletion is in effect.
- `MISSING` must precede the `ANALYSIS` subcommand.
- The `LISTWISE`, `PAIRWISE`, and `MEANSUB` keywords are alternatives, but any one of them can be used with `INCLUDE`.

<b>LISTWISE</b>	<i>Delete cases with missing values listwise.</i> Only cases with nonmissing values for all variables named on the <code>VARIABLES</code> subcommand are used. Cases are deleted even if they have missing values only for variables listed on <code>VARIABLES</code> and have valid values for all variables listed on <code>ANALYSIS</code> . Alias <code>DEFAULT</code> .
<b>PAIRWISE</b>	<i>Delete cases with missing values pairwise.</i> All cases with nonmissing values for each pair of variables correlated are used to compute that correlation, regardless of whether the cases have missing values for any other variable.
<b>MEANSUB</b>	<i>Replace missing values with the variable mean.</i> All cases are used after the substitution is made. If <code>INCLUDE</code> is also specified, user-missing values are included in the computation of the means, and means are substituted only for the system-missing value. If <code>SELECT</code> is in effect, only the values of selected cases are used in calculating the means used to replace missing values for selected cases in analysis and for all cases in computing factor scores.
<b>INCLUDE</b>	<i>Include user-missing values.</i> Cases with user-missing values are treated as valid.

## METHOD Subcommand

METHOD specifies whether the factor analysis is performed on a correlation matrix or a covariance matrix.

- Only one METHOD subcommand is allowed. If more than one is specified, the last is in effect.

<b>CORRELATION</b>	<i>Perform a correlation matrix analysis.</i> This is the default.
<b>COVARIANCE</b>	<i>Perform a covariance matrix analysis.</i> Valid only with principal components, principal axis factoring, or image factoring methods of extraction. The program issues an error if this keyword is specified when the input is a factor-loading matrix or a correlation matrix that does not contain standard deviations (STDDEV or SD).

## SELECT Subcommand

SELECT limits cases used in the analysis phase to those with a specified value for any one variable.

- Only one SELECT subcommand is allowed. If more than one is specified, the last is in effect.
- The specification is a variable name and a valid value in parentheses. A string value must be specified within quotes. Multiple variables or values are not permitted.
- The selection variable does not have to be specified on the VARIABLES subcommand.
- Only cases with the specified value for the selection variable are used in computing the correlation or covariance matrix. You can compute and save factor scores for the unselected cases as well as the selected cases.
- SELECT is not valid if MATRIX = IN is specified.

### Example

```
FACTOR VARIABLES = V1 TO V10
  /SELECT=COMPLETE(1)
  /SAVE (4) .
```

- FACTOR analyzes all ten variables named on VARIABLES, using only cases with a value of 1 for the variable *COMPLETE*.
- By default, FACTOR uses the CORRELATION method and performs the principal components analysis of the selected cases. Those with eigenvalues greater than 1 are rotated using varimax rotation.
- Four factor scores, for both selected and unselected cases, are computed using the default regression method and four new variables are saved in the active dataset.

## ANALYSIS Subcommand

The ANALYSIS subcommand specifies a subset of the variables named on VARIABLES for use in an analysis.

- The specification on ANALYSIS is a list of variables, all of which must have been named on the VARIABLES subcommand. For matrix input, ANALYSIS can specify a subset of the variables in a correlation or covariance matrix.

- Only one ANALYSIS subcommand is allowed. When multiple ANALYSIS subcommands are specified, the last is in effect.
- If no ANALYSIS is specified, all variables named on the VARIABLES subcommand (or included in the matrix input file) are used.
- Keyword TO in a variable list on ANALYSIS refers to the order in which variables are named on the VARIABLES subcommand, not to their order in the active dataset.
- Keyword ALL refers to all variables named on the VARIABLES subcommand.

### Example

```
FACTOR VARIABLES=V1 V2 V3 V4 V5 V6
  /ANALYSIS=V4 TO V6.
```

- This example requests a factor analysis of *V4*, *V5*, and *V6*. Keyword TO on ANALYSIS refers to the order of variables on VARIABLES, not the order in the active dataset.
- Cases with missing values for all variables specified on VARIABLES are omitted from the analysis. (The default setting for MISSING.)
- By default, the CORRELATION method is used and a principal components analysis with a varimax rotation is performed.

## FORMAT Subcommand

FORMAT modifies the format of factor pattern and structure matrices.

- FORMAT can be specified anywhere after VARIABLES and MISSING. If more than one FORMAT is specified, the last is in effect.
- If FORMAT is omitted or included without specifications, variables appear in the order in which they are named on ANALYSIS and all matrix entries are displayed.

**SORT**                    *Order the factor loadings in descending order.* Variables are displayed in descending order of the factor 1 loadings until a loading for factor 2 exceeds the loading for factor 1. The remaining variables are then displayed in descending order of the factor 2 loadings until a loading for factor 3 exceeds the loading for factor 2, and so on. The result shows blocks of variables that are similar.

**BLANK(n)**                *Suppress display of coefficients lower than n in absolute value.* The corresponding cells in the table will be blank.

**DEFAULT**                *Turn off keywords SORT and BLANK.*

### Example

```
FACTOR VARIABLES=V1 TO V12
  /MISSING=MEANSUB
  /FORMAT=SORT BLANK(.3)
  /EXTRACTION=ULS
  /ROTATION=NOROTATE.
```

- This example specifies an analysis of all variables between and including *V1* and *V12* in the active dataset.
- The default CORRELATION method is used.

- The `MISSING` subcommand substitutes variable means for missing values.
- The `FORMAT` subcommand orders variables in factor pattern matrices by descending value of loadings. Factor loadings with an absolute value less than 0.3 are omitted.
- Factors are extracted using unweighted least squares and are not rotated.

## **PRINT Subcommand**

`PRINT` controls the statistical display in the output.

- Keywords `INITIAL`, `EXTRACTION`, and `ROTATION` are the defaults if `PRINT` is omitted or specified without keywords.
- If any keywords are specified, only the output specifically requested is produced.
- The requested statistics are displayed only for variables specified on the last `ANALYSIS` subcommand.
- If more than one `PRINT` subcommand is specified, the last is in effect.
- If any `ANALYSIS` subcommand is explicitly specified, all `PRINT` subcommands specified before the last `ANALYSIS` subcommand are ignored. If no `PRINT` subcommand is specified after the last `ANALYSIS` subcommand, the default takes effect.

<b>INITIAL</b>	<i>Initial communalities for each variable, eigenvalues of the unreduced correlation matrix, and percentage of variance for each factor.</i>
<b>EXTRACTION</b>	<i>Factor pattern matrix, revised communalities, the eigenvalue of each factor retained, and the percentage of variance each eigenvalue represents.</i>
<b>ROTATION</b>	<i>Rotated factor pattern matrix, factor transformation matrix, factor correlation matrix, and the post-rotation sums of squared loadings.</i>
<b>UNIVARIATE</b>	<i>Valid number of cases, means, and standard deviations. (Not available with matrix input.) If <code>MISSING=MEANSUB</code> or <code>PAIRWISE</code>, the output also includes the number of missing cases.</i>
<b>CORRELATION</b>	<i>Correlation matrix. Ignored if the input is a factor-loading matrix.</i>
<b>COVARIANCE</b>	<i>Covariance matrix. Ignored if the input is a factor-loading matrix or a correlation matrix that does not contain standard deviations (<code>STDDEV</code> or <code>SD</code>).</i>
<b>SIG</b>	<i>Matrix of significance levels of correlations.</i>
<b>DET</b>	<i>Determinant of the correlation or covariance matrix, depending on the specification on <code>METHOD</code>.</i>
<b>INV</b>	<i>Inverse of the correlation or covariance matrix, depending on the specification on <code>METHOD</code>.</i>
<b>AIC</b>	<i>Anti-image covariance and correlation matrices(Kaiser, 1970). The measure of sampling adequacy for the individual variable is displayed on the diagonal of the anti-image correlation matrix.</i>
<b>KMO</b>	<i>Kaiser-Meyer-Olkin measure of sampling adequacy and Bartlett's test of sphericity. Always based on the correlation matrix. Not computed for an input matrix when it does not contain <math>N</math> values.</i>
<b>REPR</b>	<i>Reproduced correlations and residuals or reproduced covariance and residuals, depending on the specification on <code>METHOD</code>.</i>
<b>FSCORE</b>	<i>Factor score coefficient matrix. Factor score coefficients are calculated using the method requested on the <code>SAVE</code> subcommand. The default is the regression method.</i>
<b>ALL</b>	<i>All available statistics.</i>
<b>DEFAULT</b>	<i>INITIAL, EXTRACTION, and ROTATION.</i>

**Example**

```
FACTOR VARS=V1 TO V12
  /SELECT=COMPLETE ('yes')
  /MISS=MEANSUB
  /PRINT=DEF AIC KMO REPR
  /EXTRACT=ULS
  /ROTATE=VARIMAX.
```

- This example specifies a factor analysis that includes all variables between and including *V1* and *V12* in the active dataset.
- Only cases with the value “yes” on *COMPLETE* are used.
- Variable means are substituted for missing values. Only values for the selected cases are used in computing the mean. This mean is used to substitute missing values in analyzing the selected cases and in computing factor scores for all cases.
- The output includes the anti-image correlation and covariance matrices, the Kaiser-Meyer-Olkin measure of sampling adequacy, the reproduced correlation and residual matrix, as well as the default statistics.
- Factors are extracted using unweighted least squares.
- The factor pattern matrix is rotated using the varimax rotation.

**PLOT Subcommand**

Use *PLOT* to request scree plots or plots of variables in rotated factor space.

- If *PLOT* is omitted, no plots are produced. If *PLOT* is used without specifications, it is ignored.
- If more than one *PLOT* subcommand is specified, only the last one is in effect.
- If any *ANALYSIS* subcommand is explicitly specified, all *PLOT* subcommands specified before the last *ANALYSIS* subcommand are ignored. If no *PLOT* subcommand is specified after the last *ANALYSIS* subcommand, no plot is produced.

**EIGEN**                    *Scree plot*(Cattell, 1966). The eigenvalues from each extraction are plotted in descending order.

**ROTATION**                *Plots of variables in factor space*. When used without any additional specifications, *ROTATION* can produce only high-resolution graphics. If three or more factors are extracted, a 3-D plot is produced with the factor space defined by the first three factors. You can request two-dimensional plots by specifying pairs of factor numbers in parentheses; for example, *PLOT ROTATION(1, 2) (1, 3) (2, 3)* requests three plots, each defined by two factors. The *ROTATION* subcommand must be explicitly specified when you enter the keyword *ROTATION* on the *PLOT* subcommand.

**DIAGONAL Subcommand**

*DIAGONAL* specifies values for the diagonal in conjunction with principal axis factoring.

- If *DIAGONAL* is omitted or included without specifications, *FACTOR* uses the default method for specifying the diagonal.
- *DIAGONAL* is ignored with extraction methods other than *PAF*. The values are automatically adjusted by corresponding variances if *METHOD=COVARIANCE*.

- If more than one `DIAGONAL` subcommand is specified, only the last one is in effect.
- If any `ANALYSIS` subcommand is explicitly specified, `DIAGONAL` subcommands specified before the last `ANALYSIS` subcommand are ignored. If no `DIAGONAL` is specified after the last `ANALYSIS` subcommand, the default is used.
- Default communality estimates for `PAF` are squared multiple correlations. If these cannot be computed, the maximum absolute correlation between the variable and any other variable in the analysis is used.

**valuelist**                    *Diagonal values.* The number of values supplied must equal the number of variables in the analysis block. Use the notation `n*` before a value to indicate that the value is repeated `n` times.

**DEFAULT**                    *Initial communality estimates.*

### Example

```
FACTOR VARIABLES=V1 TO V12
  /DIAGONAL=.56 .55 .74 2*.56 .70 3*.65 .76 .64 .63
  /EXTRACTION=PAF
  /ROTATION=VARIMAX.
```

- The factor analysis includes all variables between and including `V1` and `V12` in the active dataset.
- `DIAGONAL` specifies 12 values to use as initial estimates of communalities in principal axis factoring.
- The factor pattern matrix is rotated using varimax rotation.

## CRITERIA Subcommand

`CRITERIA` controls extraction and rotation criteria.

- `CRITERIA` can be specified anywhere after `VARIABLES` and `MISSING`.
- Only explicitly specified criteria are changed. Unspecified criteria keep their defaults.
- Multiple `CRITERIA` subcommands are allowed. Changes made by a previous `CRITERIA` subcommand are overwritten by a later `CRITERIA` subcommand.
- Any `CRITERIA` subcommands specified after the last `EXTRACTION` subcommand have no effect on extraction.
- Any `CRITERIA` subcommands specified after the last `ROTATION` subcommand have no effect on rotation.

The following keywords on `CRITERIA` apply to extractions:

**FACTORS(n)**                    *Number of factors extracted.* The default is the number of eigenvalues greater than `MINEIGEN`. When specified, `FACTORS` overrides `MINEIGEN`.

**MINEIGEN(n)**                    *Minimum eigenvalue used to control the number of factors extracted.* If `METHOD=CORRELATION`, the default is 1. If `METHOD=COVARIANCE`, the default is computed as  $(\text{Total Variance}/\text{Number of Variables}) * n$ , where Total Variance is the total weighted variance principal components or principal axis factoring extraction and the total image variance for image factoring extraction.

**ECONVERGE(n)**                    *Convergence criterion for extraction.* The default is 0.001.

The following keywords on CRITERIA apply to rotations:

<b>RCONVERGE(n)</b>	<i>Convergence criterion for rotation. The default is 0.0001.</i>
<b>KAISER</b>	<i>Kaiser normalization in the rotation phase. This is the default. The alternative is NOKAISER.</i>
<b>NOKAISER</b>	<i>No Kaiser normalization.</i>

The following keywords on CRITERIA apply to both extractions and rotations:

<b>ITERATE(n)</b>	<i>Maximum number of iterations for solutions in the extraction or rotation phases. The default is 25.</i>
<b>DEFAULT</b>	<i>Reestablish default values for all criteria.</i>

### **Example**

```
FACTOR VARIABLES=V1 TO V12
  /CRITERIA=FACTORS (6)
  /EXTRACTION=PC
  /ROTATION=NOROTATE
  /PLOT=ROTATION.
```

- This example analyzes all variables between and including *V1* and *V12* in the active dataset.
- Six factors are extracted using the default principal components method, and the factor pattern matrix is not rotated.
- PLOT sends all extracted factors to the graphics editor and shows a 3-D plot of the first three factors.

## **EXTRACTION Subcommand**

EXTRACTION specifies the factor extraction technique.

- Only one EXTRACTION subcommand is allowed. If multiple EXTRACTION subcommands are specified, only the last is performed.
- If any ANALYSIS subcommand is explicitly specified, all EXTRACTION subcommands before the last ANALYSIS subcommand are ignored. If no EXTRACTION subcommand is specified after the last ANALYSIS subcommand, the default extraction is performed.
- If EXTRACTION is not specified or is included without specifications, principal components extraction is used.
- If you specify criteria for EXTRACTION, the CRITERIA subcommand must precede the EXTRACTION subcommand.
- When you specify EXTRACTION, you should always explicitly specify the ROTATION subcommand. If ROTATION is not specified, the factors are not rotated.

<b>PC</b>	<i>Principal components analysis(Harman, 1976). This is the default. PC can also be requested with keyword PA1 or DEFAULT.</i>
<b>PAF</b>	<i>Principal axis factoring. PAF can also be requested with keyword PA2.</i>
<b>ALPHA</b>	<i>Alpha factoring(Kaiser and Caffry, 1965). Invalid if METHOD=COVARIANCE.</i>
<b>IMAGE</b>	<i>Image factoring(Kaiser, 1963).</i>



<b>ULS</b>	<i>Unweighted least squares</i> (Jöreskog, 1977). Invalid if METHOD=COVARIANCE.
<b>GLS</b>	<i>Generalized least squares</i> . Invalid if METHOD=COVARIANCE.
<b>ML</b>	<i>Maximum likelihood</i> (Jöreskog and Lawley, 1968). Invalid if METHOD=COVARIANCE.

**Example**

```
FACTOR VARIABLES=V1 TO V12
  /ANALYSIS=V1 TO V6
  /EXTRACTION=ULS
  /ROTATE=NOROTATE.
```

- This example analyzes variables *V1* through *V6* with an unweighted least-squares extraction. No rotation is performed.

**ROTATION Subcommand**

ROTATION specifies the factor rotation method. It can also be used to suppress the rotation phase entirely.

- Only one ROTATION subcommand is allowed. If multiple ROTATION subcommands are specified, only the last is performed.
- If any ANALYSIS subcommand is explicitly specified, all ROTATION subcommands before the last ANALYSIS subcommand are ignored. If any EXTRACTION subcommand is explicitly specified, all ROTATION subcommands before the last EXTRACTION subcommand are ignored.
- If ROTATION is omitted together with EXTRACTION, varimax rotation is used.
- If ROTATION is omitted but EXTRACTION is not, factors are not rotated.
- Keyword NOROTATE on the ROTATION subcommand produces a plot of variables in unrotated factor space if the PLOT subcommand is also included for the analysis.

<b>VARIMAX</b>	<i>Varimax rotation</i> . This is the default if ROTATION is entered without specifications or if EXTRACTION and ROTATION are both omitted. Varimax rotation can also be requested with keyword DEFAULT.
<b>EQUAMAX</b>	<i>Equamax rotation</i> .
<b>QUARTIMAX</b>	<i>Quartimax rotation</i> .
<b>OBLIMIN(n)</b>	<i>Direct oblimin rotation</i> . This is a nonorthogonal rotation; thus, a factor correlation matrix will also be displayed. You can specify a delta ( $n \leq 0.8$ ) in parentheses. The value must be less than or equal to 0.8. The default is 0.
<b>PROMAX(n)</b>	<i>Promax rotation</i> . This is a nonorthogonal rotation; thus, a factor correlation matrix will also be displayed. For this method, you can specify a real-number value greater than 1. The default is 4.
<b>NOROTATE</b>	<i>No rotation</i> .

**Example**

```
FACTOR VARIABLES=V1 TO V12
  /EXTRACTION=ULS
  /ROTATION
  /ROTATION=OBLIMIN.
```

- The first ROTATION subcommand specifies the default varimax rotation.
- The second ROTATION subcommand specifies an oblimin rotation based on the same extraction of factors.

## SAVE Subcommand

SAVE allows you to save factor scores from any rotated or unrotated extraction as new variables in the active dataset. You can use any of the three methods for computing the factor scores.

- Only one SAVE subcommand is executed. If you specify multiple SAVE subcommands, only the last is executed.
- SAVE must follow the last ROTATION subcommand.
- If no ROTATION subcommand is specified after the last EXTRACTION subcommand, SAVE must follow the last EXTRACTION subcommand and no rotation is used.
- If neither ROTATION nor EXTRACTION is specified, SAVE must follow the last ANALYSIS subcommand and the default extraction and rotation are used to compute the factor scores.
- SAVE subcommands before any explicitly specified ANALYSIS, EXTRACTION, or ROTATION subcommands are ignored.
- You cannot use the SAVE subcommand if you are replacing the active dataset with matrix materials. (For more information, see the topic [Matrix Output](#) on p. 718.)
- The new variables are added to the end of the active dataset.

Keywords to specify the method of computing factor scores are:

<b>REG</b>	<i>Regression method.</i> This is the default.
<b>BART</b>	<i>Bartlett method.</i>
<b>AR</b>	<i>Anderson-Rubin method.</i>
<b>DEFAULT</b>	<i>The same as REG.</i>

- After one of the above keywords, specify in parentheses the number of scores to save and a rootname to use in naming the variables.
- You can specify either an integer or the keyword ALL. The maximum number of scores you can specify is the number of factors in the solution.
- FACTOR forms variable names by appending sequential numbers to the rootname you specify. The rootname must begin with a letter and conform to the rules for variable names. For information on variable naming rules, see [Variable Names](#) on p. 49.
- If you do not specify a rootname, FACTOR forms unique variable names using the formula  $FACn\_m$ , where  $m$  increments to create a new rootname and  $n$  increments to create a unique variable name. For example,  $FAC1\_1$ ,  $FAC2\_1$ ,  $FAC3\_1$ , and so on will be generated for the first set of saved scores and  $FAC1\_2$ ,  $FAC2\_2$ ,  $FAC3\_2$ , and so on for the second set.
- FACTOR automatically generates variable labels for the new variables. Each label contains information about the method of computing the factor score, its sequential number, and the sequential number of the analysis.

**Example**

```
FACTOR VARIABLES=V1 TO V12
  /CRITERIA FACTORS(4)
  /ROTATION
  /SAVE REG (4,PCOMP) .
```

- Since there is no `EXTRACTION` subcommand before the `ROTATION` subcommand, the default principal components extraction is performed.
- The `CRITERIA` subcommand specifies that four principal components should be extracted.
- The `ROTATION` subcommand requests the default varimax rotation for the principal components.
- The `SAVE` subcommand calculates scores using the regression method. Four scores will be added to the file: `PCOMP1`, `PCOMP2`, `PCOMP3`, and `PCOMP4`.

**MATRIX Subcommand**

`MATRIX` reads and writes matrix data files.

- `MATRIX` must always be specified first.
- Only one `IN` and one `OUT` keyword can be specified on the `MATRIX` subcommand. If either `IN` or `OUT` is specified more than once, the `FACTOR` procedure is not executed.
- The matrix type must be indicated on `IN` or `OUT`. The types are `COR` for a correlation matrix, `COV` for a covariance matrix, and `FAC` for a factor-loading matrix. Indicate the matrix type within parentheses immediately before you identify the matrix file.
- If you use both `IN` and `OUT` on `MATRIX`, you can specify them in either order. You cannot write a covariance matrix if the input matrix is a factor-loading matrix or a correlation matrix that does not contain standard deviations (`STDDEV` or `SD`).
- If you read in a covariance matrix and write out a factor-loading matrix, the output factor loadings are rescaled.

**OUT (matrix type=  
'savfile'|'dataset')**

*Write a matrix data file.* Specify the matrix type (`COR`, `COV`, `FAC`, or `FSC`) and the matrix file in parentheses. For the matrix data file, specify a filename to store the matrix materials on disk, a previously declared dataset available in the current session, or an asterisk to replace the active dataset. If you specify an asterisk or a dataset name, the matrix data file is not stored on disk unless you use `SAVE` or `XSAVE`.

**IN (matrix type=  
'savfile'|'dataset')**

*Read a matrix data file.* Specify the matrix type (`COR`, `COV`, or `FAC`) and the matrix file in parentheses. For the matrix data file, specify an asterisk if the matrix data file is the active dataset. If the matrix file is another file, specify the filename or dataset name in parentheses. A matrix file read from an external file or another dataset in the current session does not replace the active dataset.

## Matrix Output

FACTOR can write matrix materials in the form of a correlation matrix, a covariance matrix, a factor-loading matrix, or a factor score coefficients matrix.

- The correlation and covariance matrix materials include counts, means, and standard deviations in addition to correlations or covariances.
- The factor-loading matrix materials contain only factor values and no additional statistics.
- The factor score coefficients materials include means and standard deviations, in addition to factor score coefficients.
- See [Format of the Matrix Data File](#) on p. 718 for a description of the file.
- FACTOR generates one matrix per split file.
- Any documents contained in the active dataset are not transferred to the matrix file.

## Matrix Input

- FACTOR can read matrix materials written either by a previous FACTOR procedure or by a procedure that writes correlation or covariance matrices. [For more information, see the topic Universals on p. 36.](#)
- MATRIX=IN cannot be used unless a active dataset has already been defined. To read an existing matrix data file at the beginning of a session, first use GET to retrieve the matrix file and then specify IN(COR=\*), IN(COV=\*) or IN(FAC=\*) on MATRIX.
- The VARIABLES subcommand cannot be used with matrix input.
- For correlation and covariance matrix input, the ANALYSIS subcommand can specify a subset of the variables in the matrix. You cannot specify a subset of variables for factor-loading matrix input. By default, the ANALYSIS subcommand uses all variables in the matrix.

## Format of the Matrix Data File

- For correlation or covariance matrices, the matrix data file has two special variables created by the program: ROWTYPE\_ and VARNAME\_. Variable ROWTYPE\_ is a short string variable with the value CORR (for Pearson correlation coefficient) or COV (for covariance) for each matrix row. Variable VARNAME\_ is a short string variable whose values are the names of the variables used to form the correlation matrix.
- For factor-loading matrices, the program generates two special variables named ROWTYPE\_ and FACTOR\_. The value for ROWTYPE\_ is always FACTOR. The values for FACTOR\_ are the ordinal numbers of the factors.
- For factor score coefficient matrices, the matrix data file has two special variables created: ROWTYPE\_ and VARNAME\_. If split-file processing is in effect, the split variables appear first in the matrix output file, followed by ROWTYPE\_, VARNAME\_, and the variables in the analysis. ROWTYPE\_ is a short string with three possible values: MEAN, STDDEV, and FSCOEf. There is always one occurrence of the value MEAN. If /METHOD = CORRELATION then there is one occurrence of the value STDDEV. Otherwise, this value does not appear. There are as many occurrences of FSCOEf as the number of extracted factors. VARNAME\_ is a short string whose values are FAC $n$  where  $n$  is the sequence of the saved factor when ROWTYPE\_ equals FSCOEf. Otherwise the value is empty.

- The remaining variables are the variables used to form the matrix.

### **Split Files**

- FACTOR can read or write split-file matrices.
- When split-file processing is in effect, the first variables in the matrix data file are the split variables, followed by *ROWTYPE\_*, *VARNAME\_* (or *FACTOR\_*), and then the variables used to form the matrix.
- A full set of matrix materials is written for each split-file group defined by the split variables.
- A split variable cannot have the same variable name as any other variable written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split file must be in effect when that matrix is read by any other procedure.

### **Example: Factor Correlation Matrix Output to External File**

```
GET FILE='/data/GSS80.sav'
/KEEP ABDEFECT TO ABSINGLE.
FACTOR VARIABLES=ABDEFECT TO ABSINGLE
/MATRIX OUT(COR='/data/cormtx.sav').
```

- FACTOR retrieves the *GSS80.sav* file and writes a factor correlation matrix to the file *cormtx.sav*.
- The active dataset is still *GSS80.sav*. Subsequent commands will be executed on this file.

### **Example: Factor Correlation Matrix Output Replacing Active Dataset**

```
GET FILE='/data/GSS80.sav'
/KEEP ABDEFECT TO ABSINGLE.
FACTOR VARIABLES=ABDEFECT TO ABSINGLE
/MATRIX OUT(COR=*) .
LIST.
```

- FACTOR writes the same matrix as in the previous example.
- The active dataset is replaced with the correlation matrix. The LIST command is executed on the matrix file, not on *GSS80*.

### **Example: Factor-Loading Matrix Output Replacing Active Dataset**

```
GET FILE='/dataGSS80.sav'
/KEEP ABDEFECT TO ABSINGLE.
FACTOR VARIABLES=ABDEFECT TO ABSINGLE
/MATRIX OUT(FAC=*) .
```

- FACTOR generates a factor-loading matrix that replaces the active dataset.

### **Example: Matrix Input from active dataset**

```
GET FILE='/data/country.sav'
/KEEP SAVINGS POP15 POP75 INCOME GROWTH.
```

```
REGRESSION MATRIX OUT(*)
/VARS=SAVINGS TO GROWTH
/MISS=PAIRWISE
/DEP=SAVINGS /ENTER.
FACTOR MATRIX IN(COR=*) /MISSING=PAIRWISE.
```

- The GET command retrieves the *country.sav* file and selects the variables needed for the analysis.
- The REGRESSION command computes correlations among five variables with pairwise deletion. MATRIX=OUT writes a matrix data file, which replaces the active dataset.
- MATRIX IN(COR=\*) on FACTOR reads the matrix materials REGRESSION has written to the active dataset. An asterisk is specified because the matrix materials are in the active dataset. FACTOR uses pairwise deletion, since this is what was in effect when the matrix was built.

### Example: Matrix Input from External File

```
GET FILE='/data/country.sav'
/KEEP SAVINGS POP15 POP75 INCOME GROWTH.
REGRESSION
/VARS=SAVINGS TO GROWTH
/MISS=PAIRWISE
/DEP=SAVINGS /ENTER.
FACTOR MATRIX IN(COR=CORMTX) .
```

- This example performs a regression analysis on file *country.sav* and then uses a different file for FACTOR. The file is an existing matrix data file.
- MATRIX=IN specifies the matrix data file *CORMTX*.
- *CORMTX* does not replace *country.sav* as the active dataset.

### Example: Matrix Input from active dataset

```
GET FILE='/data/cormtx.sav'.
FACTOR MATRIX IN(COR=*) .
```

- This example starts a new session and reads an existing matrix data file. GET retrieves the matrix data file *cormtx.sav*.
- MATRIX=IN specifies an asterisk because the matrix data file is the active dataset. If MATRIX=IN(*cormtx.sav*) is specified, the program issues an error message.
- If the GET command is omitted, the program issues an error message.

### Example: Using Saved Coefficients to Score an External File

```
MATRIX.
GET A /FILE="fsc.sav".
GET B /FILE="ext_data_standardized.sav" /VAR=varlist.
COMPUTE SCORES=A*TRANSPOS(B).
SAVE SCORES /OUTFILE="scored.sav".
END MATRIX.
```

- This example scores an external file using the factor score coefficients from a previous analysis, say, upon *original.sav*.

- Factor score coefficients are read from *fsc.sav* into *A*.
- The data are read from *ext\_data\_standardized.sav* into *B*. The variable values in the external file should be standardized. Note that if the values in *ext\_data\_standardized.sav* are standardized based upon the means and standard deviations of the unstandardized values in *ext\_data.sav*, and not upon the means and standard deviations of values in *original.sav*, then this analysis will be different from one in which *original.sav* and *ext\_data.sav* are merged and FACTOR is run using the SELECT subcommand with cases from *original.sav* selected for use in computing the correlation or covariance matrix.

If there are missing values, add /MISSING=OMIT or /MISSING=0 to the second GET statement to remove cases with missing values or impute the mean (0, since the variables are standardized).

- The scores are saved to *scored.sav*.

## References

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# FILE HANDLE

```
FILE HANDLE handle /NAME='path and/or file specifications'  
                [/MODE={CHARACTER }] [/RECFORM={FIXED } [/LRECL=n]  
                {BINARY } {VARIABLE}  
                {MULTIPUNCH} {SPANNED }  
                {IMAGE }  
                {360 }  
                [/ENCODING = 'encoding specification']
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Release History

### Release 13.0

- The NAME subcommand is modified to accept a path and/or file.

### Release 16.0

- ENCODING subcommand added for Unicode support.

## Example

```
FILE HANDLE thisMonthFile /NAME='/sales/data/july.sav'.  
FILE HANDLE dataDirectory /NAME='/sales/data'.  
GET FILE 'thisMonthFile'.  
GET FILE 'dataDirectory/july.sav'.
```

## Overview

FILE HANDLE assigns a unique *file handle* to a path and/or file and supplies operating system specifications for the file. A defined file handle can be specified on any subsequent FILE, OUTFILE, MATRIX, or WRITE subcommands of various procedures.

## Syntax Rules

- FILE HANDLE is required for reading data files with record lengths greater than 8,192. [For more information, see the topic LRECL Subcommand on p. 724.](#)
- FILE HANDLE is required for reading IBM VSAM datasets, EBCDIC data files, binary data files, and character data files that are not delimited by ASCII line feeds.
- If you specify 360 on the MODE subcommand, you must specify RECFORM.
- If you specify IMAGE on the MODE subcommand, you must specify LRECL.

## Operations

A file handle is used only during the current session. The handle is never saved as part of a data file. The normal quoting conventions for file specifications apply, with or without file handles.



## Example

```
FILE HANDLE thisMonthFile /NAME='/sales/data/july.sav'.
FILE HANDLE dataDirectory /NAME='/sales/data'.
GET FILE 'thisMonthFile'.
GET FILE 'dataDirectory/july.sav'.
```

- The first FILE HANDLE command defines a file handle that refers to a specific file.
- The second FILE HANDLE command only specifies a directory path.
- The two subsequent GET FILE commands are functionally equivalent. Note that both file specifications are enclosed in quotes (a good general practice).

## NAME Subcommand

NAME specifies the path and/or file you want to refer to by the file handle. The file specifications must conform to the file naming convention for the type of computer and operating system on which the program is run. See the documentation for your system for specific information about the file naming convention.

If NAME specifies a relative path or does not include a path, the path is set to the current working directory at the time the FILE HANDLE command is executed.

## MODE Subcommand

MODE specifies the type of file you want to refer to by the file handle.

<b>CHARACTER</b>	<i>Character file whose logical records are delimited by ASCII line feeds.</i>
<b>BINARY</b>	<i>Unformatted binary file generated by Microsoft FORTRAN.</i>
<b>MULTIPUNCH</b>	<i>Column binary file.</i>
<b>IMAGE</b>	<i>Binary file consisting of fixed-length records.</i>
<b>360</b>	<i>EBCDIC data file.</i>

### Example

```
FILE HANDLE ELE48 /NAME='/data/ELE48.DAT' /MODE=MULTIPUNCH.
DATA LIST FILE=ELE48.
```

- FILE HANDLE defines *ELE48* as the handle for the file.
- The MODE subcommand indicates that the file contains multipunch data.
- The file specification on NAME conforms to VMS convention: the file *ELE48.DAT* is located in the directory *data*.
- The FILE subcommand on DATA LIST refers to the handle defined on the FILE HANDLE command.

## RECFORM Subcommand

RECFORM specifies the record format and is necessary when you specify 360 on MODE. RECFORM has no effect with other specifications on MODE.

<b>FIXED</b>	<i>Fixed-length record.</i> All records have the same length. Alias F. When FIXED is specified, the record length must be specified on the LRECL subcommand.
<b>VARIABLE</b>	<i>Variable-length record.</i> No logical record is larger than one physical block. Alias V.
<b>SPANNED</b>	<i>Spanned record.</i> Records may be larger than fixed-length physical blocks. Alias VS.

## LRECL Subcommand

LRECL specifies the length of each record in the file. When you specify IMAGE under UNIX, OS/2, or Microsoft Windows, or 360 for IBM360 EBCDIC data files, you must specify LRECL. You can specify a record length greater than the default (8,192) for an image file, a character file, or a binary file. The maximum record length is 2,147,483,647. Do not use LRECL with MULTIPUNCH.

### Example

```
FILE HANDLE TRGT1 /NAME='/data/RGT.DAT'
                /MODE=IMAGE LRECL=16.
DATA LIST FILE=TRGT1.
```

- IMAGE is specified on the MODE subcommand. Subcommand LRECL must be specified.
- The file handle is used on the DATA LIST command.

## ENCODING Subcommand

ENCODING specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is UTF8. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), UTF8, UTF16, UTF16BE (big endian), UTF16LE (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- In Unicode mode, the defined width of string variables is tripled for code page and UTF-16 text data files. Use [ALTER TYPE](#) to automatically adjust the defined width of string variables.

# FILE LABEL

```
FILE LABEL label text
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
FILE LABEL Original survey responses prior to recoding.
```

## **Overview**

FILE LABEL provides a descriptive label for a data file.

### **Basic Specification**

The basic specification is the command name followed by the label text.

### **Syntax Rules**

- The label text cannot exceed 64 bytes.
- Labels do not need to be enclosed in quotes.
- If the label is enclosed in quotes—or starts with a quotation mark (single or double)—standard rules for quoted strings apply. [For more information, see the topic String Values in Command Specifications on p. 40.](#)

### **Operations**

- If the data file is saved in IBM® SPSS® Statistics format, the label is saved in the dictionary.
- The file label is displayed in the Notes tables generated by procedures.
- A data file can only contain one file label. Subsequent FILE LABEL commands replace the label text.

## **Example**

```
FILE LABEL Respondent's original data.  
FILE LABEL "Respondent's original data".  
FILE LABEL 'Respondent's original data.'
```

- The first two commands are functionally equivalent. The enclosing double-quotes in the second command are not included as part of the label text, and the apostrophe (single quote) is preserved.
- In the last command, everything after the apostrophe in *Respondent's* will be omitted from the label because the apostrophe will be interpreted as the closing single quote to match the opening single quote.

# FILE TYPE-END FILE TYPE

*For mixed file types:*

```
FILE TYPE MIXED [FILE='file'] [ENCODING='encoding specification']
  RECORD=[varname] column location [(format)]

  [WILD={NOWARN}]
    {WARN }
```

*For grouped file types:*

```
FILE TYPE GROUPED [FILE='file'] [ENCODING='encoding specification']
  RECORD=[varname] column location [(format)]

CASE=[varname] column location [(format)]

[WILD={WARN  }] [DUPLICATE={WARN  }]
  {NOWARN}          {NOWARN}

[MISSING={WARN  }] [ORDERED={YES}]
  {NOWARN}          {NO }
```

*For nested file types:*

```
FILE TYPE NESTED [FILE='file'] [ENCODING='encoding specification']
  RECORD=[varname] column location [(format)]

  [CASE=[varname] column location [(format)]]

  [WILD={NOWARN}] [DUPLICATE={NOWARN}]
    {WARN  }          {WARN  }
                    {CASE  }

  [MISSING={NOWARN}]
    {WARN  }

END FILE TYPE
```

## **Release History**

Release 16.0

- ENCODING subcommand added for Unicode support.

## **Example**

```
FILE TYPE MIXED RECORD=RECID 1-2.
RECORD TYPE 23.
DATA LIST /SEX 5 AGE 6-7 DOSAGE 8-10 RESULT 12.
END FILE TYPE.
```

```
BEGIN DATA
21 145010 1
22 257200 2
25 235 250 2
35 167          300 3
24 125150 1
23 272075 1
21 149050 2
25 134 035 3
30 138          300 3
```

32 229                    500    3  
 END DATA.

## Overview

The FILE TYPE-END FILE TYPE structure defines data for any one of the three types of complex raw data files: **mixed files**, which contain several types of records that define different types of cases; **hierarchical** or **nested files**, which contain several types of records with a defined relationship among the record types; or **grouped files**, which contain several records for each case with some records missing or duplicated. A fourth type of complex file, files with **repeating groups** of information, can be defined with the REPEATING DATA command.

FILE TYPE must be followed by at least one RECORD TYPE command and one DATA LIST command. Each pair of RECORD TYPE and DATA LIST commands defines one type of record in the data. END FILE TYPE signals the end of file definition.

Within the FILE TYPE structure, the lowest-level record in a nested file can be read with a REPEATING DATA command rather than a DATA LIST command. In addition, any record in a mixed file can be read with REPEATING DATA.

### Basic Specification

The basic specification on FILE TYPE is one of the three file type keywords (MIXED, GROUPED, or NESTED) and the RECORD subcommand. RECORD names the record identification variable and specifies its column location. If keyword GROUPED is specified, the CASE subcommand is also required. CASE names the case identification variable and specifies its column location.

The FILE TYPE-END FILE TYPE structure must enclose at least one RECORD TYPE and one DATA LIST command. END FILE TYPE is required to signal the end of file definition.

- RECORD TYPE specifies the values of the record type identifier (see RECORD TYPE).
- DATA LIST defines variables for the record type specified on the preceding RECORD TYPE command (see DATA LIST).
- Separate pairs of RECORD TYPE and DATA LIST commands must be used to define each different record type.

The resulting active dataset is always a rectangular file, regardless of the structure of the original data file.

### Syntax Rules

- For mixed files, if the record types have different variables or if they have the same variables recorded in different locations, separate RECORD TYPE and DATA LIST commands are required for each record type.
- For mixed files, the same variable name can be used on different DATA LIST commands, since each record type defines a separate case.
- For mixed files, if the same variable is defined for more than one record type, the format type and length of the variable should be the same on all DATA LIST commands. The program refers to the *first* DATA LIST command that defines a variable for the print and write formats to include in the dictionary of the active dataset.

- For grouped and nested files, the variable names on each `DATA LIST` must be unique, since a case is built by combining all record types together into a single record.
- For nested files, the order of the `RECORD TYPE` commands defines the hierarchical structure of the file. The first `RECORD TYPE` defines the highest-level record type, the next `RECORD TYPE` defines the next highest-level record, and so forth. The last `RECORD TYPE` command defines a case in the active dataset. By default, variables from higher-level records are spread to the lowest-level record.
- For nested files, the `SPREAD` subcommand on `RECORD TYPE` can be used to spread the values in a record type only to the *first* case built from each record of that type. All other cases associated with that record are assigned the system-missing value for the variables defined on that type. See `RECORD TYPE` for more information.
- String values specified on the `RECORD TYPE` command must be enclosed in quotes.

### Operations

- For mixed file types, the program skips all records that are not specified on one of the `RECORD TYPE` commands.
- If different variables are defined for different record types in mixed files, the variables are assigned the system-missing value for those record types on which they are not defined.
- For nested files, the first record in the file should be the type specified on the first `RECORD TYPE` command—the highest level of the hierarchy. If the first record in the file is not the highest-level type, the program skips all records until it encounters a record of the highest-level type. If `MISSING` or `DUPLICATE` has been specified, these records may produce warning messages but will not be used to build a case in the active dataset.
- When defining complex files, you are effectively building an input program and can use only commands that are allowed in the input state.

## Examples

### Reading multiple record types from a mixed file

```
FILE TYPE MIXED FILE='/data/treatmnt.txt' RECORD=RECID 1-2.
+ RECORD TYPE 21,22,23,24.
+ DATA LIST /SEX 5 AGE 6-7 DOSAGE 8-10 RESULT 12.
+ RECORD TYPE 25.
+ DATA LIST /SEX 5 AGE 6-7 DOSAGE 10-12 RESULT 15.
END FILE TYPE.
```

- Variable `DOSAGE` is read from columns 8–10 for record types 21, 22, 23, and 24 and from columns 10–12 for record type 25. `RESULT` is read from column 12 for record types 21, 22, 23, and 24, and from column 15 for record type 25.
- The active dataset contains values for all variables defined on the `DATA LIST` commands for record types 21 through 25. All other record types are skipped.

### Reading only one record type from a mixed file

```
FILE TYPE MIXED RECORD=RECID 1-2.
RECORD TYPE 23.
DATA LIST /SEX 5 AGE 6-7 DOSAGE 8-10 RESULT 12.
```

```

END FILE TYPE.

BEGIN DATA
21 145010 1
22 257200 2
25 235 250 2
35 167          300 3
24 125150 1
23 272075 1
21 149050 2
25 134 035 3
30 138          300 3
32 229          500 3
END DATA.

```

- FILE TYPE begins the file definition and END FILE TYPE indicates the end of file definition. FILE TYPE specifies a mixed file type. Since the data are included between BEGIN DATA-END DATA, the FILE subcommand is omitted. The record identification variable *RECID* is located in columns 1 and 2.
- RECORD TYPE indicates that records with value 23 for variable *RECID* will be copied into the active dataset. All other records are skipped. the program does not issue a warning when it skips records in mixed files.
- DATA LIST defines variables on records with the value 23 for variable *RECID*.

### **A grouped file of student test scores**

```

FILE TYPE GROUPED RECORD=#TEST 6 CASE=STUDENT 1-4.
RECORD TYPE 1.
DATA LIST /ENGLISH 8-9 (A).
RECORD TYPE 2.
DATA LIST /READING 8-10.
RECORD TYPE 3.
DATA LIST /MATH 8-10.
END FILE TYPE.

```

```

BEGIN DATA
0001 1 B+
0001 2 74
0001 3 83
0002 1 A
0002 2 100
0002 3 71
0003 1 B-
0003 2 88
0003 3 81
0004 1 C
0004 2 94
0004 3 91
END DATA.

```

- FILE TYPE identifies the file as a grouped file. As required for grouped files, all records for a single case are together in the data. The record identification variable *#TEST* is located in column 6. A scratch variable is specified so it won't be saved in the active dataset. The case identification variable *STUDENT* is located in columns 1-4.
- Because there are three record types, there are three RECORD TYPE commands. For each RECORD TYPE, there is a DATA LIST to define variables on that record type.
- END FILE TYPE signals the end of file definition.

- The program builds four cases—one for each student. Each case includes the case identification variable plus the variables defined for each record type (the test scores). The values for #*TEST* are not saved in the active dataset. Thus, each case in the active dataset has four variables: *STUDENT*, *ENGLISH*, *READING*, and *MATH*.

### **A nested file of accident records**

```
FILE TYPE NESTED RECORD=6 CASE=ACCID 1-4.
RECORD TYPE 1.
DATA LIST /ACC_ID 9-11 WEATHER 12-13 STATE 15-16 (A) DATE 18-24 (A).
RECORD TYPE 2.
DATA LIST /STYLE 11 MAKE 13 OLD 14 LICENSE 15-16(A) INSURNCE 18-21 (A).
RECORD TYPE 3.
DATA LIST /PSNGR_NO 11 AGE 13-14 SEX 16 (A) INJURY 18 SEAT 20-21 (A)
          COST 23-24.
END FILE TYPE.

BEGIN DATA
0001 1 322 1 IL 3/13/88 /* Type 1:  accident record
0001 2 1 44MI 134M /* Type 2:  vehicle record
0001 3 1 34 M 1 FR 3 /* Type 3:  person record
0001 2 2 16IL 322F /*          vehicle record
0001 3 1 22 F 1 FR 11 /*          person record
0001 3 2 35 M 1 FR 5 /*          person record
0001 3 3 59 M 1 BK 7 /*          person record
0001 2 3 21IN 146M /*          vehicle record
0001 3 1 46 M 0 FR 0 /*          person record
END DATA.
```

- `FILE TYPE` specifies a nested file type. The record identifier, located in column 6, is not assigned a variable name, so the default scratch variable name `####RECD` is used. The case identification variable *ACCID* is located in columns 1–4.
- Because there are three record types, there are three `RECORD TYPE` commands. For each `RECORD TYPE`, there is a `DATA LIST` command to define variables on that record type. The order of the `RECORD TYPE` commands defines the hierarchical structure of the file.
- `END FILE TYPE` signals the end of file definition.
- The program builds a case for each lowest-level (type 3) record, representing each person in the file. There can be only one type 1 record for each type 2 record, and one type 2 record for each type 3 record. Each vehicle can be in only one accident, and each person can be in only one vehicle. The variables from the type 1 and type 2 records are spread to their corresponding type 3 records.

## **Specification Order**

- `FILE TYPE` must be the first command in the `FILE TYPE-END FILE TYPE` structure. `FILE TYPE` subcommands can be named in any order.
- Each `RECORD TYPE` command must precede its corresponding `DATA LIST` command.
- `END FILE TYPE` must be the last command in the structure.



## Types of Files

The first specification on `FILE TYPE` is a file type keyword, which defines the structure of the data file. There are three file type keywords: `MIXED`, `GROUPED`, and `NESTED`. Only one of the three types can be specified on `FILE TYPE`.

<b>MIXED</b>	<i>Mixed file type.</i> <code>MIXED</code> specifies a file in which each record type named on a <code>RECORD TYPE</code> command defines a case. You do not need to define all types of records in the file. In fact, <code>FILE TYPE MIXED</code> is useful for reading only one type of record because the program can decide whether to execute the <code>DATA LIST</code> for a record by simply reading the variable that identifies the record type.
<b>GROUPED</b>	<i>Grouped file type.</i> <code>GROUPED</code> defines a file in which cases are defined by grouping together record types with the same identification number. Each case usually has one record of each type. All records for a single case must be together in the file. By default, the program assumes that the records are in the same sequence within each case.
<b>NESTED</b>	<i>Nested file type.</i> <code>NESTED</code> defines a file in which the record types are related to each other hierarchically. The record types are grouped together by a case identification number that identifies the highest level—the first record type—of the hierarchy. Usually, the last record type specified—the lowest level of the hierarchy—defines a case. For example, in a file containing household records and records for each person living in the household, each person record defines a case. Information from higher record types may be <i>spread</i> to each case. For example, the value for a variable on the household record, such as <code>CITY</code> , can be spread to the records for each person in the household.

## Subcommands and Their Defaults for Each File Type

The specifications on the `FILE TYPE` differ for each type of file. The following table shows whether each subcommand is required or optional and, where applicable, what the default specification is for each file type. N/A indicates that the subcommand is not applicable to that type of file.

**Table 87-1**  
Summary of `FILE TYPE` subcommands for different file types

<b>Subcommand</b>	<b>Mixed</b>	<b>Grouped</b>	<b>Nested</b>
<code>FILE</code>	Conditional	Conditional	Conditional
<code>RECORD</code>	Required	Required	Required
<code>CASE</code>	Not Applicable	Required	Optional
<code>WILD</code>	<code>NOWARN</code>	<code>WARN</code>	<code>NOWARN</code>
<code>DUPLICATE</code>	N/A	<code>WARN</code>	<code>NOWARN</code>
<code>MISSING</code>	N/A	<code>WARN</code>	<code>NOWARN</code>
<code>ORDERED</code>	N/A	<code>YES</code>	N/A

- `FILE` is required unless data are inline (included between `BEGIN DATA-END DATA`).
- `RECORD` is always required.
- `CASE` is required for grouped files.
- The subcommands `CASE`, `DUPLICATE`, and `MISSING` can also be specified on the associated `RECORD TYPE` commands for grouped files. However, `DUPLICATE=CASE` is invalid.

- For nested files, `CASE` and `MISSING` can be specified on the associated `RECORD TYPE` commands.
- If the subcommands `CASE`, `DUPLICATE`, or `MISSING` are specified on a `RECORD TYPE` command, the specification on the `FILE TYPE` command (or the default) is overridden only for the record types listed on that `RECORD TYPE` command. The `FILE TYPE` specification or default applies to all other record types.

## **FILE Subcommand**

`FILE` specifies a text file containing the data. `FILE` is not used when the data are inline.

### **Example**

```
FILE TYPE MIXED FILE='/data/treatmnt.txt' RECORD=RECID 1-2.
```

- Data are in the file *treatmnt.txt*. The file type is mixed. The record identification variable *RECID* is located in columns 1 and 2 of each record.

## **ENCODING Subcommand**

`ENCODING` specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is `UTF8`. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), `UTF8`, `UTF16`, `UTF16BE` (big endian), `UTF16LE` (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- In Unicode mode, the defined width of string variables is tripled for code page and UTF-16 text data files. Use [ALTER TYPE](#) to automatically adjust the defined width of string variables.
- If there is no `FILE` subcommand, the `ENCODING` subcommand is ignored.

## **RECORD Subcommand**

`RECORD` specifies the name and column location of the record identification variable.

- The column location of the record identifier is required. The variable name is optional.
- If you do not want to save the record type variable, you can assign a scratch variable name by using the `#` character as the first character of the name. If a variable name is not specified on `RECORD`, the record identifier is defined as the scratch variable `####RECD`.

- The value of the identifier for each record type must be unique and must be in the same location on all records. However, records do not have to be sorted according to type.
- A column-style format can be specified for the record identifier. For example, the following two specifications are valid:

```
RECORD=V1 1-2 (N)
RECORD=V1 1-2 (F, 1)
```

FORTTRAN-like formats cannot be used because the column location must be specified explicitly.

- Specify `A` in parentheses after the column location to define the record type variable as a string variable.

### Example

```
FILE TYPE MIXED FILE='/data/treatmnt.txt' RECORD=RECID 1-2.
```

- The record identifier is variable *RECID*, located in columns 1 and 2 of the hospital treatment data file.

## CASE Subcommand

CASE specifies a name and column location for the case identification variable. CASE is required for grouped files and optional for nested files. It cannot be used with mixed files.

- For grouped files, each unique value for the case identification variable defines a case in the active dataset.
- For nested files, the case identification variable identifies the highest-level record of the hierarchy. The program issues a warning message for each record with a case identification number not equal to the case identification number on the last highest-level record. However, the record with the invalid case number is used to build the case.
- The column location of the case identifier is required. The variable name is optional.
- If you do not want to save the case identification variable, you can assign a scratch variable name by using the `#` character as the first character of the name. If a variable name is not specified on CASE, the case identifier is defined as the scratch variable `####CASE`.
- A column-style format can be specified for the case identifier. For example, the following two specifications are valid:

```
CASE=V1 1-2 (N)
CASE=V1 1-2 (F, 1)
```

FORTTRAN-like formats cannot be used because the column location must be specified explicitly.

- Specify `A` in parentheses after the column location to define the case identification variable as a string variable.
- If the case identification number is not in the same columns on all record types, use the CASE subcommand on the RECORD TYPE commands as well as on the FILE TYPE command (see RECORD TYPE).

**Example**

\* A grouped file of student test scores.

```
FILE TYPE GROUPED RECORD=#TEST 6 CASE=STUDENT 1-4.
RECORD TYPE 1.
DATA LIST /ENGLISH 8-9 (A).
RECORD TYPE 2.
DATA LIST /READING 8-10.
RECORD TYPE 3.
DATA LIST /MATH 8-10.
END FILE TYPE.
```

```
BEGIN DATA
0001 1 B+
0001 2 74
0001 3 83
0002 1 A
0002 2 100
0002 3 71
0003 1 B-
0003 2 88
0003 3 81
0004 1 C
0004 2 94
0004 3 91
END DATA.
```

- CASE is required for grouped files. CASE specifies variable *STUDENT*, located in columns 1–4, as the case identification variable.
- The data contain four different values for *STUDENT*. The active dataset therefore has four cases, one for each value of *STUDENT*. In a grouped file, each unique value for the case identification variable defines a case in the active dataset.
- Each case includes the case identification variable plus the variables defined for each record type. The values for *#TEST* are not saved in the active dataset. Thus, each case in the active dataset has four variables: *STUDENT*, *ENGLISH*, *READING*, and *MATH*.

**Example**

\* A nested file of accident records.

```
FILE TYPE NESTED RECORD=6 CASE=ACCID 1-4.
RECORD TYPE 1.
DATA LIST /ACC_ID 9-11 WEATHER 12-13 STATE 15-16 (A) DATE 18-24 (A).
RECORD TYPE 2.
DATA LIST /STYLE 11 MAKE 13 OLD 14 LICENSE 15-16 (A) INSURNCE 18-21 (A).
RECORD TYPE 3.
DATA LIST /PSNGR_NO 11 AGE 13-14 SEX 16 (A) INJURY 18 SEAT 20-21 (A)
COST 23-24.
END FILE TYPE.
```

```
BEGIN DATA
0001 1 322 1 IL 3/13/88 /* Type 1: accident record
0001 2 1 44MI 134M /* Type 2: vehicle record
0001 3 1 34 M 1 FR 3 /* Type 3: person record
0001 2 2 16IL 322F /* vehicle record
0001 3 1 22 F 1 FR 11 /* person record
0001 3 2 35 M 1 FR 5 /* person record
0001 3 3 59 M 1 BK 7 /* person record
0001 2 3 21IN 146M /* vehicle record
0001 3 1 46 M 0 FR 0 /* person record
```

END DATA.

- CASE specifies variable *ACCID*, located in columns 1–4, as the case identification variable. *ACCID* identifies the highest level of the hierarchy: the level for the accident records.
- As each case is built, the value of the variable *ACCID* is checked against the value of *ACCID* on the last highest-level record (record type 1). If the values do not match, a warning message is issued. However, the record is used to build the case.
- The data in this example contain only one value for *ACCID*, which is spread across all cases. In a nested file, the lowest-level record type determines the number of cases in the active dataset. In this example, the active dataset has five cases because there are five person records.

### Example

```
* Specifying case on the RECORD TYPE command.

FILE TYPE GROUPED FILE=HUBDATA RECORD=#RECID 80 CASE=ID 1-5.
RECORD TYPE 1.
DATA LIST /MOHIRED YRHIRED 12-15 DEPT79 TO DEPT82 SEX 16-20.
RECORD TYPE 2.
DATA LIST /SALARY79 TO SALARY82 6-25 HOURLY81 HOURLY82 40-53 (2)
          PROMO81 72 AGE 54-55 RAISE82 66-70.
RECORD TYPE 3 CASE=75-79.
DATA LIST /JOB CAT 6 NAME 25-48 (A).
END FILE TYPE.
```

- The CASE subcommand on FILE TYPE indicates that the case identification number is located in columns 1–5. However, for type 3 records, the case identification number is located in columns 75–79. The CASE subcommand is therefore specified on the third RECORD TYPE command to override the case setting for type 3 records.
- The format of the case identification variable must be the same on all records. If the case identification variable is defined as a string on the FILE TYPE command, it cannot be defined as a numeric variable on the RECORD TYPE command, and vice versa.

## WILD Subcommand

WILD determines whether the program issues a warning when it encounters undefined record types in the data file. Regardless of whether the warning is issued, undefined records are not included in the active dataset.

- The only specification on WILD is keyword WARN or NOWARN.
- WARN cannot be specified if keyword OTHER is specified on the last RECORD TYPE command to indicate all other record types (see RECORD TYPE).

<b>WARN</b>	<i>Issue warning messages.</i> The program displays a warning message and the first 80 characters of the record for each record type that is not mentioned on a RECORD TYPE command. This is the default for grouped file types.
<b>NOWARN</b>	<i>Suppress warning messages.</i> The program simply skips all record types not mentioned on a RECORD TYPE command and does not display warning messages. This is the default for mixed and nested file types.

**Example**

```
FILE TYPE MIXED FILE='/data/treatmnt.txt' RECORD=RECID 1-2 WILD=WARN.
```

- WARN is specified on the WILD subcommand. The program displays a warning message and the first 80 characters of the record for each record type that is not mentioned on a RECORD TYPE command.

**DUPLICATE Subcommand**

DUPLICATE determines how the program responds when it encounters more than one record of each type for a single case. DUPLICATE is optional for grouped and nested files. DUPLICATE cannot be used with mixed files.

- The only specification on DUPLICATE is keyword WARN, NOWARN, or CASE.

<b>WARN</b>	<i>Issue warning messages.</i> The program displays a warning message and the first 80 characters of the last record of the duplicate set of record types. Only the <i>last</i> record from a set of duplicates is included in the active dataset. This is the default for grouped files.
<b>NOWARN</b>	<i>Suppress warning messages.</i> The program does not display warning messages when it encounters duplicate record types. Only the <i>last</i> record from a set of duplicates is included in the active dataset. This is the default for nested files.
<b>CASE</b>	<i>Build a case in the active dataset for each duplicate record.</i> The program builds one case in the active dataset for each duplicate record, spreading information from any higher-level records and assigning system-missing values to the variables defined on lower-level records. This option is available only for nested files.

**Example**

```
* A nested file of accident records.
* Issue a warning for duplicate record types.

FILE TYPE NESTED RECORD=6 CASE=ACCID 1-4 DUPLICATE=WARN.
RECORD TYPE 1.
DATA LIST /ACC_ID 9-11 WEATHER 12-13 STATE 15-16 (A) DATE 18-24 (A).
RECORD TYPE 2.
DATA LIST /STYLE 11 MAKE 13 OLD 14 LICENSE 15-16 (A) INSURNCE 18-21 (A).
RECORD TYPE 3.
DATA LIST /PSNGR_NO 11 AGE 13-14 SEX 16 (A) INJURY 18 SEAT 20-21 (A)
        COST 23-24.
END FILE TYPE.

BEGIN DATA
0001 1 322 1 IL 3/13/88 /*          accident record
0001 2 1 44MI 134M /*          vehicle record
0001 3 1 34 M 1 FR 3 /*          person record
0001 2 1 31IL 134M /* duplicate vehicle record
0001 2 2 16IL 322F /*          vehicle record
0001 3 1 22 F 1 FR 11 /*          person record
0001 3 2 35 M 1 FR 5 /*          person record
0001 3 3 59 M 1 BK 7 /*          person record
0001 2 3 21IN 146M /*          vehicle record
0001 3 1 46 M 0 FR 0 /*          person record
END DATA.
```

- In the data, there are two vehicle (type 2) records above the second set of person (type 3) records. This implies that an empty (for example, parked) vehicle was involved, or that each of the three persons was in two vehicles, which is impossible.
- `DUPLICATE` specifies keyword `WARN`. The program displays a warning message and the first 80 characters of the second of the duplicate set of type 2 records. The first duplicate record is skipped, and only the second is included in the active dataset. This assumes that no empty vehicles were involved in the accident.
- If the duplicate record represents an empty vehicle, it can be included in the active dataset by specifying keyword `CASE` on `DUPLICATE`. The program builds one case in the active dataset for the first duplicate record, spreading information to that case from the previous type 1 record and assigning system-missing values to the variables defined for type 3 records. The second record from the duplicate set is used to build the three cases for the associated type 3 records.

## ***MISSING Subcommand***

`MISSING` determines whether the program issues a warning when it encounters a missing record type for a case. Regardless of whether the program issues the warning, it builds the case in the active dataset with system-missing values for the variables defined on the missing record. `MISSING` is optional for grouped and nested files.

- `MISSING` cannot be used with mixed files and is optional for grouped and nested files.
- For grouped and nested files, the program verifies that each defined case includes one record of each type.
- The only specification is keyword `WARN` or `NOWARN`.

**WARN**                    *Issue a warning message when a record type is missing for a case. This is the default for grouped files.*

**NOWARN**                *Suppress the warning message when a record type is missing for a case. This is the default for nested files.*

### ***Example***

\* A grouped file with missing records.

```
FILE TYPE GROUPED RECORD=#TEST 6 CASE=STUDENT 1-4 MISSING=NOWARN.
RECORD TYPE 1.
DATA LIST /ENGLISH 8-9 (A).
RECORD TYPE 2.
DATA LIST /READING 8-10.
RECORD TYPE 3.
DATA LIST /MATH 8-10.
END FILE TYPE.
```

```
BEGIN DATA
0001 1 B+
0001 2 74
0002 1 A
0002 2 100
0002 3 71
0003 3 81
0004 1 C
0004 2 94
0004 3 91
```

END DATA.

- The data contain records for three tests administered to four students. However, not all students took all tests. The first student took only the English and reading tests. The third student took only the math test.
- One case in the active dataset is built for each of the four students. If a student did not take a test, the system-missing value is assigned in the active dataset to the variable for the missing test. Thus, the first student has the system-missing value for the math test, and the third student has missing values for the English and reading tests.
- Keyword `NOWARN` is specified on `MISSING`. Therefore, no warning messages are issued for the missing records.

### Example

\* A nested file with missing records.

```
FILE TYPE NESTED RECORD=6 CASE=ACCID 1-4 MISSING=WARN.
RECORD TYPE 1.
DATA LIST /ACC_ID 9-11 WEATHER 12-13 STATE 15-16 (A) DATE 18-24 (A).
RECORD TYPE 2.
DATA LIST /STYLE 11 MAKE 13 OLD 14 LICENSE 15-16 (A) INSURNCE 18-21 (A).
RECORD TYPE 3.
DATA LIST /PSNGR_NO 11 AGE 13-14 SEX 16 (A) INJURY 18 SEAT 20-21 (A)
        COST 23-24.
END FILE TYPE.
```

```
BEGIN DATA
0001 1 322 1 IL 3/13/88 /*          accident record
0001 3 1 34 M 1 FR 3 /*          person record
0001 2 2 16IL 322F /*          vehicle record
0001 3 1 22 F 1 FR 11 /*          person record
0001 3 2 35 M 1 FR 5 /*          person record
0001 3 3 59 M 1 BK 7 /*          person record
0001 2 3 21IN 146M /*          vehicle record
0001 3 1 46 M 0 FR 0 /*          person record
END DATA.
```

- The data contain records for one accident. The first record is a type 1 (accident) record, and the second record is a type 3 (person) record. However, there is no type 2 record, and therefore no vehicle associated with the first person. The person may have been a pedestrian, but it is also possible that the vehicle record is missing.
- One case is built for each person record. The first case has missing values for the variables specified on the vehicle record.
- Keyword `WARN` is specified on `MISSING`. A warning message is issued for the missing record.

## ORDERED Subcommand

`ORDERED` indicates whether the records are in the same order as they are defined on the `RECORD TYPE` commands. Regardless of the order of the records in the data file and the specification on `ORDERED`, the program builds cases in the active dataset with records in the order defined on the `RECORD TYPE` commands.

- `ORDERED` can be used only for grouped files.



- The only specification is keyword YES or NO.
- If YES is in effect but the records are not in the order defined on the RECORD TYPE commands, the program issues a warning for each record that is out of order. The program still uses these records to build cases.

**YES**                      *Records for each case are in the same order as they are defined on the RECORD TYPE commands. This is the default.*

**NO**                        *Records are not in the same order within each case.*

### **Example**

\* A grouped file with records out of order.

```
FILE TYPE GROUPED RECORD=#TEST 6 CASE=STUDENT 1-4 MISSING=NOWARN
  ORDERED=NO.
RECORD TYPE 1.
DATA LIST /ENGLISH 8-9 (A).
RECORD TYPE 2.
DATA LIST /READING 8-10.
RECORD TYPE 3.
DATA LIST /MATH 8-10.
END FILE TYPE.
```

```
BEGIN DATA
0001 2 74
0001 1 B+
0002 3 71
0002 2 100
0002 1 A
0003 2 81
0004 2 94
0004 1 C
0004 3 91
END DATA.
```

- The first RECORD TYPE command specifies record type 1, the second specifies record type 2, and the third specifies record type 3. However, records for each case are not always ordered type 1, type 2, and type 3.
- NO is specified on ORDERED. The program builds cases without issuing a warning that they are out of order in the data.
- Regardless of whether YES or NO is in effect for ORDERED, the program builds cases in the active dataset in the same order specified on the RECORD TYPE commands.

# ***FILTER***

```
FILTER {BY var}  
{OFF }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Example***

```
FILTER BY SEX.  
FREQUENCIES BONUS.
```

## ***Overview***

`FILTER` is used to exclude cases from program procedures without deleting them from the active dataset. When `FILTER` is in effect, cases with a zero or missing value for the specified variable are not used in program procedures. Those cases are not actually deleted and are available again if the filter is turned off. To see the current filter status, use the `SHOW` command.

### ***Basic Specification***

The basic specification is keyword `BY` followed by a variable name. Cases that have a zero or missing value for the filter variable are excluded from subsequent procedures.

### ***Syntax Rules***

- Only one numeric variable can be specified. The variable can be one of the original variables in the data file or a variable computed with transformation commands.
- Keyword `OFF` turns off the filter. All cases in the active dataset become available to subsequent procedures.
- If `FILTER` is specified without a keyword, `FILTER OFF` is assumed but the program displays a warning message.
- `FILTER` can be specified anywhere in the command sequence. Unlike `SELECT IF`, `FILTER` has the same effect within an input program as it does outside an input program. Attention must be paid to the placement of any transformation command used to compute values for the filter variable (see `INPUT PROGRAM`).

### ***Operations***

- `FILTER` performs case selection without changing the active dataset. Cases that have a zero or missing value are excluded from subsequent procedures but are not deleted from the file.
- Both system-missing and user-missing values are treated as missing. The `FILTER` command does not offer options for changing selection criteria. To set up different criteria for exclusion, create a numeric variable and conditionally compute its values before specifying it on `FILTER`.

- If `FILTER` is specified after `TEMPORARY`, `FILTER` affects the next procedure only. After that procedure, the filter status reverts to whatever it was before the `TEMPORARY` command.
- The filter status does not change until another `FILTER` command is specified, a `USE` command is specified, or the active dataset is replaced.
- `FILTER` and `USE` are mutually exclusive. `USE` automatically turns off any previous `FILTER` command, and `FILTER` automatically turns off any previous `USE` command.
- If the specified filter variable is renamed, it is still in effect. The `SHOW` command will display the new name of the filter variable. However, the filter is turned off if the filter variable is recoded into a string variable or is deleted from the file.
- If the active dataset is replaced after a `MATCH FILES`, `ADD FILES`, or `UPDATE` command and the active dataset is one of the input files, the filter remains in effect if the new active dataset has a numeric variable with the name of the filter variable. If the active dataset does not have a numeric variable with that name (for example, if the filter variable was dropped or renamed), the filter is turned off.
- If the active dataset is replaced by an entirely new data file (for example, by a `DATA LIST`, `GET`, or `IMPORT` command), the filter is turned off.
- The `FILTER` command changes the filter status and takes effect when a procedure is executed or an `EXECUTE` command is encountered.

## Examples

### **Filter by a variable with values of 0 and 1**

```
FILTER BY SEX.  
FREQUENCIES BONUS.
```

- This example assumes that `SEX` is a numeric variable, with male and female coded as 0 and 1, respectively. The `FILTER` command excludes males and cases with missing values for `SEX` from the subsequent procedures. The `FREQUENCIES` command generates a frequency table of `BONUS` for females only.

### **Recoding the filter variable to change the filter criterion**

```
RECODE SEX (1=0) (0=1) .  
FILTER BY SEX.  
FREQUENCIES BONUS.
```

- This example assumes the same coding scheme for `SEX` as the previous example. Before `FILTER` is specified, variable `SEX` is recoded. The `FILTER` command then excludes females and cases with missing values for `SEX`. The `FREQUENCIES` command generates a frequency table of `BONUS` for males only.

# FINISH

FINISH

## Overview

FINISH causes the program to stop reading commands.

### Operations

- FINISH immediately causes the program to stop reading commands.
- The appearance of FINISH on the printback of commands in the display file indicates that the session has been completed.

## Example

```
DATA LIST FILE=RAWDATA /NAME 1-15(A) V1 TO V15 16-30.  
LIST.  
FINISH.  
REPORT FORMAT=AUTO LIST /VARS=NAME V1 TO V10.
```

- FINISH causes the program to stop reading commands after LIST is executed. The REPORT command is not executed.

## Basic Specification

The basic specification is keyword FINISH. There are no additional specifications.

## Command Files

- FINISH is optional in a command file and is used to mark the end of a session.
- FINISH causes the program to stop reading commands. Anything following FINISH in the command file is ignored. Any commands following FINISH in an INCLUDE file are ignored.
- FINISH cannot be used within a DO IF structure to end a session conditionally. FINISH within a DO IF structure will end the session unconditionally.

## Prompted Sessions

- FINISH is required in a prompted session to terminate the session.
- Because FINISH is a program command, it can be used only after the command line prompt for the program, which expects a procedure name. FINISH cannot be used to end a prompted session from a DATA>, CONTINUE>, HELP>, or DEFINE> prompt.

# FIT

```
FIT [[ERRORS=] residual series names]
    [/OBS=observed series names]
    [{DFE=error degrees of freedom    }]
    {DFH=hypothesis degrees of freedom}
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
FIT.
```

## Overview

FIT displays a variety of descriptive statistics computed from the residual series as an aid in evaluating the goodness of fit of one or more models.

### Options

**Statistical Output.** You can produce statistics for a particular residual series by specifying the names of the series after FIT. You can also obtain percent error statistics for specified residual series by specifying observed series on the OBS subcommand.

**Degrees of Freedom.** You can specify the degrees of freedom for the residual series using the DFE or DFH subcommands.

### Basic Specification

The basic specification is simply the command keyword FIT. All other specifications are optional.

- By default, FIT calculates the mean error, mean percent error, mean absolute error, mean absolute percent error, sum of squared errors, mean square error, root mean square error, and the Durbin-Watson statistic for the last *ERR\_n* (residual) series generated and the corresponding observed series in the active dataset.
- If neither residual nor observed series are specified, percent error statistics for the default residual and observed series are included.

### Syntax Rules

- If OBS is specified, the ERRORS subcommand naming the residual series is required.

### Operations

- Observed series and degrees of freedom are matched with residual series according to the order in which they are specified.

- If residual series are explicitly specified but observed series are not, percent error statistics are not included in the output. If neither residual nor observed series are specified, percent error statistics for the default residual and observed series are included.
- If subcommand `DFH` is specified, `FIT` calculates the `DFE` (error degrees of freedom) by subtracting the `DFH` (hypothesis degrees of freedom) from the number of valid cases in the series.
- If a `PREDICT` period (validation period) starts before the end of the observed series, statistics are reported separately for the `USE` period (historical period) and the `PREDICT` period.

### **Limitations**

- There is no limit on the number of residual series specified. However, the number of observed series must equal the number of residual series.

## **Example**

```
FIT ERR_4 ERR_5 ERR_6 .
```

- This command requests goodness-of-fit statistics for the residual series `ERR_4`, `ERR_5`, and `ERR_6`, which were generated by previous procedures. Percent error statistics are not included in the output, since only residual series are named.

## **ERRORS Subcommand**

`ERRORS` specifies the residual (error) series.

- The actual keyword `ERRORS` can be omitted. `VARIABLES` is an alias for `ERRORS`.
- The minimum specification on `ERRORS` is a residual series name.
- The `ERRORS` subcommand is required if the `OBS` subcommand is specified.

## **OBS Subcommand**

`OBS` specifies the observed series to use for calculating the mean percentage error and mean absolute percentage error.

- `OBS` can be used only when the residual series are explicitly specified.
- The number and order of observed series must be the same as that of the residual series.
- If more than one residual series was calculated from a single observed series, the observed series is specified once for each residual series that is based on it.

### **Example**

```
FIT ERRORS=ERR#1 ERR#2
/OBS=VAR1 VAR1 .
```

- This command requests `FIT` statistics for two residual series, `ERR#1` and `ERR#2`, which were computed from the same observed series, `VAR1`.

## ***DFE and DFH Subcommands***

DFE and DFH specify the degrees of freedom for each residual series. With DFE, error degrees of freedom are entered directly. DFH specifies hypothesis degrees of freedom so FIT can compute the DFE.

- Only one DFE or DFH subcommand should be specified. If both are specified, only the last one is in effect.
- The specification on DFE or DFH is a list of numeric values. The order of these values should correspond to the order of the residual series list.
- The error degrees of freedom specified on DFE are used to compute the mean square error (MSE) and root mean square (RMS).
- The value specified for DFH should equal the number of parameters in the model (including the constant if it is present). Differencing is not considered in calculating DFH, since any observations lost due to differencing are system-missing.
- If neither DFE or DFH are specified, FIT sets DFE equal to the number of observations.

### ***Example***

```
FIT ERR#1 ERR#2
  /OBS=VAR1 VAR2
  /DFE=47 46.
```

- In this example, the error degrees of freedom for the first residual series, *ERR#1*, is 47. The error degrees of freedom for the second residual series, *ERR#2*, is 46.

## ***Output Considerations for SSE***

The sum of squared errors (SSE) reported by FIT may not be the same as the SSE reported by the estimation procedure. The SSE from the procedure is an estimate of sigma squared for that model. The SSE from FIT is simply the sum of the squared residuals.

## ***References***

Makridakis, S., S. C. Wheelwright, and V. E. McGee. 1983. *Forecasting: Methods and applications*. New York: John Wiley and Sons.

McLaughlin, R. L. 1984. *Forecasting techniques for decision making*. Rockville, Md.: Control Data Management Institute.

# FLIP

```
FLIP [VARIABLES= {ALL }]  
      {varlist}  
  
      [/NEWNAMES=variable]
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
FLIP.
```

## Overview

The program requires a file structure in which the variables are the columns and observations (cases) are the rows. If a file is organized such that variables are in rows and observations are in columns, you need to use `FLIP` to reorganize it. `FLIP` transposes the rows and columns of the data in the active dataset so that, for example, row 1, column 2 becomes row 2, column 1, and so forth.

## Options

**Variable Subsets.** You can transpose specific variables (columns) from the original file using the `VARIABLES` subcommand.

**Variable Names.** You can use the values of one of the variables from the original file as the variable names in the new file, using the `NEWNAMES` subcommand.

## Basic Specification

The basic specification is the command keyword `FLIP`, which transposes all rows and columns.

- By default, `FLIP` assigns variable names `VAR001` to `VARn` to the variables in the new file. It also creates the new variable `CASE_LBL`, whose values are the variable names that existed before transposition.

## Subcommand Order

`VARIABLES` must precede `NEWNAMES`.

## Operations

- `FLIP` replaces the active dataset with the transposed file and displays a list of variable names in the transposed file.
- `FLIP` discards any previous `VARIABLE LABELS`, `VALUE LABELS`, and `WEIGHT` settings. Values defined as user-missing in the original file are translated to system-missing in the transposed file.



- FLIP obeys any SELECT IF, N, and SAMPLE commands in effect.
- FLIP does not obey the TEMPORARY command. Any transformations become permanent when followed by FLIP.
- String variables in the original file are assigned system-missing values after transposition.
- Numeric variables are assigned a default format of F8.2 after transposition (with the exceptions of CASE\_LBL and the variable specified on NEWNAMES).
- The variable CASE\_LBL is created and added to the active dataset each time FLIP is executed.
- If CASE\_LBL already exists as the result of a previous FLIP, its current values are used as the names of variables in the new file (if NEWNAMES is not specified).

## Example

The following is the LIST output for a data file arranged in a typical spreadsheet format, with variables in rows and observations in columns:

A	B	C	D
Income	22.00	31.00	43.00
Price	34.00	29.00	50.00
Year	1970.00	1971.00	1972.00

The command

```
FLIP.
```

transposes all variables in the file. The LIST output for the transposed file is as follows:

CASE_LBL	VAR001	VAR002	VAR003
A	.	.	.
B	22.00	34.00	1970.00
C	31.00	29.00	1971.00
D	43.00	50.00	1972.00

- The values for the new variable CASE\_LBL are the variable names from the original file.
- Case A has system-missing values, since variable A had the string values *Income*, *Price*, and *Year*.
- The names of the variables in the new file are CASE\_LBL, VAR001, VAR002, and VAR003.

## VARIABLES Subcommand

VARIABLES names one or more variables (columns) to be transposed. The specified variables become observations (rows) in the new active dataset.

- The VARIABLES subcommand is optional. If it is not used, all variables are transposed.
- If the VARIABLES subcommand is specified, variables that are not named are discarded.

### Example

Using the untransposed file from the previous example, the command

```
FLIP VARIABLES=A TO C.
```

transposes only variables *A* through *C*. Variable *D* is not transposed and is discarded from the active dataset. The LIST output for the transposed file is as follows:

CASE_LBL	VAR001	VAR002	VAR003
A	.	.	.
B	22.00	34.00	1970.00
C	31.00	29.00	1971.00

## NEWNAMES Subcommand

NEWNAMES specifies a variable whose values are used as the new variable names.

- The NEWNAMES subcommand is optional. If it is not used, the new variable names are either *VAR001* to *VARn*, or the values of *CASE\_LBL* if it exists.
- Only one variable can be specified on NEWNAMES.
- The variable specified on NEWNAMES does not become an observation (case) in the new active dataset, regardless of whether it is specified on the VARIABLES subcommand.
- If the variable specified is numeric, its values become a character string beginning with the prefix *K\_*.
- Characters not allowed in variables names, such as blank spaces, are replaced with underscore (`_`) characters.
- If the variable's values are not unique, unique variable names are created by appending a sequential suffix of the general form *\_A*, *\_B*, *\_C*,...*\_AA*, *\_AB*, *\_AC*,...*\_AAA*, *\_AAB*, *\_AAC*,...etc.

### Example

Using the untransposed file from the first example, the command

```
FLIP NEWNAMES=A.
```

uses the values for variable *A* as variable names in the new file. The LIST output for the transposed file is as follows:

CASE_LBL	INCOME	PRICE	YEAR
B	22.00	34.00	1970.00
C	31.00	29.00	1971.00
D	43.00	50.00	1972.00

- Variable *A* does not become an observation in the new file. The string values for *A* are converted to upper case.

The following command transposes this file back to a form resembling its original structure:

```
FLIP.
```

The LIST output for the transposed file is as follows:

CASE_LBL	B	C	D
----------	---	---	---

INCOME	22.00	31.00	43.00
PRICE	34.00	29.00	50.00
YEAR	1970.00	1971.00	1972.00

- Since the `NEWNAMES` subcommand is not used, the values of `CASE_LBL` from the previous `FLIP` (*B*, *C*, and *D*) are used as variable names in the new file.
- The values of `CASE_LBL` are now *INCOME*, *PRICE*, and *YEAR*.

# FORMATS

```
FORMATS varlist(format) [varlist...]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
FORMATS SALARY (DOLLAR8) / HOURLY (DOLLAR7.2) / RAISE BONUS (PCT2).
```

## Overview

FORMATS changes variable print and write formats. In this program, print and write formats are *output* formats. **Print formats**, also called display formats, control the form in which values are displayed by a procedure or by the PRINT command; **write formats** control the form in which values are written by the WRITE command.

FORMATS changes both print and write formats. To change only print formats, use PRINT FORMATS. To change only write formats, use WRITE FORMATS. For information on assigning input formats during data definition, see DATA LIST.

For detailed information on available formats and specifications, see [Variable Types and Formats](#).

## Basic Specification

The basic specification is a variable list followed by a format specification in parentheses. All variables on the list receive the new format.

## Operations

- Unlike most transformations, FORMATS takes effect as soon as it is encountered in the command sequence. Special attention should be paid to its position among commands. [For more information, see the topic Command Order on p. 41.](#)
- Variables not specified on FORMATS retain their current print and write formats in the active dataset. To see the current formats, use the DISPLAY command.
- The new formats are changed only in the active dataset and are in effect for the duration of the current session or until changed again with a FORMATS, PRINT FORMATS, or WRITE FORMATS command. Formats in the original data file (if one exists) are not changed unless the file is resaved with the SAVE or XSAVE command.
- New numeric variables created with transformation commands are assigned default print and write formats of F8.2 (or the format specified on the FORMAT subcommand of SET). The FORMATS command can be used to change the new variable's print and write formats.

- For string variables, you can only use `FORMATS` to switch between `A` and `AHEX` formats, and the `AHEX` length must be exactly twice the `A` length. `FORMATS` cannot be used to change the length of string variables. To change the defined length of a string variable, use the `ALTER TYPE` command.
- If a numeric data value exceeds its width specification, the program attempts to display some value nevertheless. The program first rounds decimal values, then removes punctuation characters, then tries scientific notation, and finally, if there is still not enough space, produces asterisks indicating that a value is present but cannot be displayed in the assigned width.

## Syntax Rules

- You can specify more than one variable or variable list, followed by a format in parentheses. Only one format can be specified after each variable list. For clarity, each set of specifications can be separated by a slash.
- You can use keyword `TO` to refer to consecutive variables in the active dataset.
- The specified width of a format must include enough positions to accommodate any punctuation characters such as decimal points, commas, dollar signs, or date and time delimiters. (This differs from assigning an *input* format on `DATA LIST`, where the program automatically expands the input format to accommodate punctuation characters in output.)
- Custom currency formats (`CCw`, `CCw.d`) must first be defined on the `SET` command before they can be used on `FORMATS`.
- For string variables, you can only use `FORMATS` to switch between `A` and `AHEX` formats. `FORMATS` cannot be used to change the length of string variables. To change the length of a string variable, declare a new variable of the desired length with the `STRING` command and then use `COMPUTE` to copy values from the existing string into the new variable.

## Examples

### Changing Formats for Multiple Variables

```
FORMATS SALARY (DOLLAR8) /HOURLY (DOLLAR7.2)
      /RAISE BONUS (PCT2) .
```

- The print and write formats for `SALARY` are changed to `DOLLAR` format with eight positions, including the dollar sign and comma when appropriate. The value 11550 is displayed as \$11,550. An eight-digit number would require a `DOLLAR11` format: eight characters for the digits, two characters for commas, and one character for the dollar sign.
- The print and write formats for `HOURLY` are changed to `DOLLAR` format with seven positions, including the dollar sign, decimal point, and two decimal places. The value 115 is displayed as \$115.00. If `DOLLAR6.2` had been specified, the value 115 would be displayed as \$115.0. The program would truncate the last 0 because a width of 6 is not enough to display the full value.
- The print and write formats for both `RAISE` and `BONUS` are changed to `PCT` with two positions: one position for the percentage and one position for the percent sign. The value 9 is displayed as 9%. Because the width allows for only two positions, the value 10 is displayed as 10, since the percent sign is truncated.

**Changing Default Variable Formats**

```
COMPUTE V3=V1 + V2.  
FORMATS V3 (F3.1).
```

- `COMPUTE` creates the new numeric variable `V3`. By default, `V3` is assigned an `F8.2` format (or the default format specified on `SET`).
- `FORMATS` changes both the print and write formats for `V3` to `F3.1`.

**Working With Custom Currency Formats**

```
SET CCA='-/- .Df1 ..-' .  
FORMATS COST (CCA14.2).
```

- `SET` defines a European currency format for the custom currency format type `CCA`.
- `FORMATS` assigns format `CCA` to variable `COST`. With the format defined for `CCA` on `SET`, the value 37419 is displayed as `Dfl 37.419,00`. See the `SET` command for more information on custom currency formats.

# FREQUENCIES

FREQUENCIES is available in the Statistics Base option.

```
FREQUENCIES VARIABLES=varlist [varlist...]

[/FORMAT= [{NOTABLE }] [{AVALUE**}
           {LIMIT(n)}  {DVALUE  }
                    {AFREQ   }
                    {DFREQ   }

[/MISSING=INCLUDE]

[/BARCHART=[MINIMUM(n)] [MAXIMUM(n)] [{FREQ(n)  }]]
                                           {PERCENT(n)}

[/PIECHART=[MINIMUM(n)] [MAXIMUM(n)] [{FREQ  }]] [{MISSING  }]]
                                           {PERCENT}  {NOMISSING}

[/HISTOGRAM=[MINIMUM(n)] [MAXIMUM(n)] [{FREQ(n)  }]] [{NONORMAL}]] ]
                                           {NORMAL  }

[/GROUPED=varlist [{(width)  }]]
                  {(boundary list)}

[/NTILES=n]

[/PERCENTILES=value list]

[/STATISTICS=[DEFAULT] [MEAN] [STDDEV] [MINIMUM] [MAXIMUM]
             [SEMEAN] [VARIANCE] [SKWNESS] [SESKEW] [RANGE]
             [MODE] [KURTOSIS] [SEKURT] [MEDIAN] [SUM] [ALL]
             [NONE]]

[/ORDER=[{ANALYSIS}] ] [{VARIABLE}]
```

\*\* Default if subcommand is omitted or specified without keyword.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
FREQUENCIES VARIABLES = RACE.
```

## Overview

FREQUENCIES produces Frequency tables showing frequency counts and percentages of the values of individual variables. You can also use FREQUENCIES to obtain Statistics tables for categorical variables and to obtain Statistics tables and graphical displays for continuous variables.

## Options

**Display Format.** You can suppress tables and alter the order of values within tables using the FORMAT subcommand.

**Statistical Display.** Percentiles and ntiles are available for numeric variables with the `PERCENTILES` and `NTILES` subcommands. The following statistics are available with the `STATISTICS` subcommand: mean, median, mode, standard deviation, variance, skewness, kurtosis, and sum.

**Plots.** Histograms can be specified for numeric variables on the `HISTOGRAM` subcommand. Bar charts can be specified for numeric or string variables on the `BARCHART` subcommand.

**Input Data.** On the `GROUPED` subcommand, you can indicate whether the input data are grouped (or collapsed) so that a better estimate can be made of percentiles.

### **Basic Specification**

The basic specification is the `VARIABLES` subcommand and the name of at least one variable. By default, `FREQUENCIES` produces a Frequency table.

### **Subcommand Order**

Subcommands can be named in any order.

### **Syntax Rules**

- You can specify multiple `NTILES` subcommands.
- `BARCHART` and `HISTOGRAM` are mutually exclusive.
- You can specify numeric variables (with or without decimal values) or string variables. Only the short-string portion of long-string variables are tabulated.
- Keyword `ALL` can be used on `VARIABLES` to refer to all user-defined variables in the active dataset.

### **Operations**

- Variables are tabulated in the order they are mentioned on the `VARIABLES` subcommand.
- If a requested ntile or percentile cannot be calculated, a period (.) is displayed.
- `FREQUENCIES` dynamically builds the table, setting up one cell for each unique value encountered in the data.

### **Limitations**

- Maximum 1,000 variables total per `FREQUENCIES` command.

## **Examples**

### **Including a Statistics Table in the Output**

```
FREQUENCIES  
  VARIABLES=dept  
  /PIECHART  
  /ORDER= ANALYSIS .
```

- The procedure produces a frequency table and pie chart for the variable *dept*.



**Suppressing the Frequency Tables in the Output**

```

FREQUENCIES
  VARIABLES=sale
  /FORMAT=NOTABLE
  /NTILES= 4
  /HISTOGRAM= NORMAL
  /STATISTICS=STDDEV MINIMUM MAXIMUM MEAN MEDIAN
    SKEWNESS KURTOSIS
  /ORDER= ANALYSIS .

```

- The procedure produces a table of statistical summaries and a histogram for *Amount of Last Sale*.
- `FORMAT` suppresses the Frequency tables, which are not useful for continuous variables.

**VARIABLES Subcommand**

`VARIABLES` names the variables to be tabulated and is the only required subcommand.

**FORMAT Subcommand**

`FORMAT` controls various features of the output, including order of categories and suppression of tables.

- The minimum specification is a single keyword.
- By default, `FREQUENCIES` displays the Frequency table and sort categories in ascending order of values for numeric variables and in alphabetical order for string variables.

**Table Order**

<b>AVALUE</b>	<i>Sort categories in ascending order of values (numeric variables) or in alphabetical order (string variables). This is the default.</i>
<b>DVALUE</b>	<i>Sort categories in descending order of values (numeric variables) or in reverse alphabetical order (string variables). This is ignored when HISTOGRAM, NTILES, or PERCENTILES is requested.</i>
<b>AFREQ</b>	<i>Sort categories in ascending order of frequency. This is ignored when HISTOGRAM, NTILES, or PERCENTILES is requested.</i>
<b>DFREQ</b>	<i>Sort categories in descending order of frequency. This is ignored when HISTOGRAM, NTILES, or PERCENTILES is requested.</i>

**Table Suppression**

<b>LIMIT(n)</b>	<i>Suppress frequency tables with more than n categories. The number of missing and valid cases and requested statistics are displayed for suppressed tables.</i>
<b>NOTABLE</b>	<i>Suppress all frequency tables. The number of missing and valid cases are displayed for suppressed tables. NOTABLE overrides LIMIT.</i>

## **BARChart Subcommand**

BARChart produces a bar chart for each variable named on the VARIABLES subcommand. By default, the horizontal axis for each bar chart is scaled in frequencies, and the interval width is determined by the largest frequency count for the variable being plotted. Bar charts are labeled with value labels or with the value itself if no label is defined.

- The minimum specification is the BARChart keyword, which generates default bar charts.
- BARChart cannot be used with HISTOGRAM.

<b>MIN(n)</b>	<i>Lower bound below which values are not plotted.</i>
<b>MAX(n)</b>	<i>Upper bound above which values are not plotted.</i>
<b>FREQ(n)</b>	<i>Vertical axis scaled in frequencies, where optional n is the maximum. If n is not specified or if it is too small, FREQUENCIES chooses 5, 10, 20, 50, 100, 200, 500, 1000, 2000, and so forth, depending on the largest category. This is the default.</i>
<b>PERCENT(n)</b>	<i>Vertical axis scaled in percentages, where optional n is the maximum. If n is not specified or if it is too small, FREQUENCIES chooses 5, 10, 25, 50, or 100, depending on the frequency count for the largest category.</i>

### **Producing a Basic Bar Chart**

```
FREQUENCIES VARIABLES = RACE /BARChart.
```

- FREQUENCIES produces a frequency table and the default bar chart for variable *RACE*.

### **Producing a Custom Bar Chart**

```
FREQUENCIES VARIABLES = V1 V2 /BAR=MAX(10).
```

- FREQUENCIES produces a frequency table and bar chart with values through 10 for each of variables *V1* and *V2*.

## **PIEChart Subcommand**

PIEChart produces a pie chart for each variable named on the VARIABLES subcommand. By default, one slice corresponds to each category defined by the variable with one slice representing all missing values. Pie charts are labeled with value labels or with the value if no label is defined.

- The minimum specification is the PIEChart keyword, which generates default pie charts.
- PIEChart can be requested together with either BARChart or HISTOGRAM.
- FREQ and PERCENT are mutually exclusive. If both are specified, only the first specification is in effect.
- MISSING and NOMISSING are mutually exclusive. If both are specified, only the first specification is in effect.

<b>MIN(n)</b>	<i>Lower bound below which values are not plotted.</i>
<b>MAX(n)</b>	<i>Upper bound above which values are not plotted.</i>
<b>FREQ</b>	<i>The pie charts are based on frequencies. Frequencies are displayed when you request values in the Chart Editor. This is the default.</i>

<b>PERCENT</b>	<i>The pie charts are based on percentage.</i> Percentage is displayed when you request values in the Chart Editor.
<b>MISSING</b>	<i>User-missing and system-missing values are treated as one category.</i> This is the default. Specify <code>INCLUDE</code> on the <code>MISSING</code> subcommand to display system-missing and user-missing values as separate slices.
<b>NOMISSING</b>	<i>Missing values are excluded from the chart.</i> If you specify <code>INCLUDE</code> on the <code>MISSING</code> subcommand, each user-missing value is represented by one slice.

### ***Producing a Basic Pie Chart***

```
FREQUENCIES VARIABLES = RACE /PIECHART.
```

- `FREQUENCIES` produces a frequency table and the default pie chart for variable `RACE`.

### ***Producing a Custom Pie Chart***

```
FREQUENCIES VARIABLES = V1 V2 /PIE=MAX(10).
```

- For each variable `V1` and `V2`, `FREQUENCIES` produces a frequency table and a pie chart with values through 10.

## ***HISTOGRAM Subcommand***

`HISTOGRAM` displays a plot for each numeric variable named on the `VARIABLES` subcommand. By default, the horizontal axis of each histogram is scaled in frequencies and the interval width is determined by the largest frequency count of the variable being plotted.

- The minimum specification is the `HISTOGRAM` keyword, which generates default histograms.
- `HISTOGRAM` cannot be used with `BAR` or `BAR`.

<b>MIN(n)</b>	<i>Lower bound below which values are not plotted.</i>
<b>MAX(n)</b>	<i>Upper bound above which values are not plotted.</i>
<b>FREQ(n)</b>	<i>Vertical axis scaled in frequencies, where optional n is the scale. If n is not specified or if it is too small, FREQUENCIES chooses 5, 10, 20, 50, 100, 200, 500, 1000, 2000, and so forth, depending on the largest category. This is the default.</i>
<b>NORMAL</b>	<i>Superimpose a normal curve.</i> The curve is based on all valid values for the variable, including values excluded by <code>MIN</code> and <code>MAX</code> .
<b>NONORMAL</b>	<i>Suppress the normal curve.</i> This is the default.

### ***Example***

```
FREQUENCIES VARIABLES = V1 /HIST=NORMAL.
```

- `FREQUENCIES` requests a histogram with a superimposed normal curve.

## GROUPED Subcommand

When the values of a variable represent grouped or collapsed data, it is possible to estimate percentiles for the original, ungrouped data from the grouped data. The `GROUPED` subcommand specifies which variables have been grouped. It affects only the output from the `PERCENTILES` and `NTILES` subcommands and the `MEDIAN` statistic from the `STATISTICS` subcommand.

- Multiple `GROUPED` subcommands can be used on a single `FREQUENCIES` command. Multiple variable lists, separated by slashes, can appear on a single `GROUPED` subcommand.
- The variables named on `GROUPED` must have been named on the `VARIABLES` subcommand.
- The value or value list in the parentheses is optional. When it is omitted, the program treats the values of the variables listed on `GROUPED` as midpoints. If the values are not midpoints, they must first be recoded with the `RECODE` command.
- A single value in parentheses specifies the width of each grouped interval. The data values must be group midpoints, but there can be empty categories. For example, if you have data values of 10, 20, and 30 and specify an interval width of 5, the categories are  $10 \pm 2.5$ ,  $20 \pm 2.5$ , and  $30 \pm 2.5$ . The categories  $15 \pm 2.5$  and  $25 \pm 2.5$  are empty.
- A value list in the parentheses specifies interval boundaries. The data values do not have to represent midpoints, but the lowest boundary must be lower than any value in the data. If any data values exceed the highest boundary specified (the last value within the parentheses), they will be assigned to an open-ended interval. In this case, some percentiles cannot be calculated.

### Basic Example

```
RECODE AGE (1=15) (2=25) (3=35) (4=45) (5=55)
        (6=65) (7=75) (8=85) (9=95)
  /INCOME (1=5) (2=15) (3=25) (4=35) (5=45)
        (6=55) (7=65) (8=75) (9=100) .

FREQUENCIES VARIABLES=AGE, SEX, RACE, INCOME
  /GROUPED=AGE, INCOME
  /PERCENTILES=5, 25, 50, 75, 95.
```

- The `AGE` and `INCOME` categories of 1, 2, 3, and so forth are recoded to category midpoints. Note that data can be recoded to category midpoints on any scale; here `AGE` is recoded in years, but `INCOME` is recoded in thousands of dollars.
- The `GROUPED` subcommand on `FREQUENCIES` allows more accurate estimates of the requested percentiles.

### Specifying the Width of Each Grouped Interval

```
FREQUENCIES VARIABLES=TEMP
  /GROUPED=TEMP (0.5)
  /NTILES=10.
```

- The values of `TEMP` (temperature) in this example were recorded using an inexpensive thermometer whose readings are precise only to the nearest half degree.
- The observed values of 97.5, 98, 98.5, 99, and so on, are treated as group midpoints, smoothing out the discrete distribution. This yields more accurate estimates of the deciles.

**Specifying Interval Boundaries**

```
FREQUENCIES VARIABLES=AGE
  /GROUPED=AGE (17.5, 22.5, 27.5, 32.5, 37.5, 42.5, 47.5
                52.5, 57.5, 62.5, 67.5, 72.5, 77.5, 82.5)
  /PERCENTILES=5, 10, 25, 50, 75, 90, 95.
```

- The values of *AGE* in this example have been estimated to the nearest five years. The first category is 17.5 to 22.5, the second is 22.5 to 27.5, and so forth. The artificial clustering of age estimates at multiples of five years is smoothed out by treating *AGE* as grouped data.
- It is not necessary to recode the ages to category midpoints, since the interval boundaries are explicitly given.

**PERCENTILES Subcommand**

PERCENTILES displays the value below which the specified percentage of cases falls. The desired percentiles must be explicitly requested. There are no defaults.

**Example**

```
FREQUENCIES VARIABLES = V1 /PERCENTILES=10 25 33.3 66.7 75.
```

- FREQUENCIES requests the values for percentiles 10, 25, 33.3, 66.7, and 75 for *V1*.

**NTILES Subcommand**

NTILES calculates the percentages that divide the distribution into the specified number of categories and displays the values below which the requested percentages of cases fall. There are no default ntiles.

- Multiple NTILES subcommands are allowed. Each NTILES subcommand generates separate percentiles. Any duplicate percentiles generated by different NTILES subcommands are consolidated in the output.

**Basic Example**

```
FREQUENCIES VARIABLES=V1 /NTILES=4.
```

- FREQUENCIES requests quartiles (percentiles 25, 50, and 75) for *V1*.

**Working With Multiple NTILES Subcommands**

```
FREQUENCIES VARIABLES=V1 /NTILES=4 /NTILES=10.
```

- The first NTILES subcommand requests percentiles 25, 50, and 75.
- The second NTILES subcommand requests percentiles 10 through 90 in increments of 10.
- The 50th percentile is produced by both specifications but is displayed only once in the output.

## STATISTICS Subcommand

STATISTICS controls the display of statistics. By default, cases with missing values are excluded from the calculation of statistics.

- The minimum specification is the keyword STATISTICS, which generates the mean, standard deviation, minimum, and maximum (these statistics are also produced by keyword DEFAULT).

MEAN	<i>Mean.</i>
SEMEAN	<i>Standard error of the mean.</i>
MEDIAN	<i>Median. Ignored when AFREQ or DFREQ are specified on the FORMAT subcommand.</i>
MODE	<i>Mode. If there is more than one mode, only the first mode is displayed.</i>
STDDEV	<i>Standard deviation.</i>
VARIANCE	<i>Variance.</i>
SKEWNESS	<i>Skewness.</i>
SESKEW	<i>Standard error of the skewness statistic.</i>
KURTOSIS	<i>Kurtosis.</i>
SEKURT	<i>Standard error of the kurtosis statistic.</i>
RANGE	<i>Range.</i>
MINIMUM	<i>Minimum.</i>
MAXIMUM	<i>Maximum.</i>
SUM	<i>Sum.</i>
DEFAULT	<i>Mean, standard deviation, minimum, and maximum.</i>
ALL	<i>All available statistics.</i>
NONE	<i>No statistics.</i>

### Specifying a Particular Statistic

```
FREQUENCIES VARIABLES = AGE /STATS=MODE.
```

- STATISTICS requests the mode of AGE.

### Including the Default Statistics

```
FREQUENCIES VARIABLES = AGE /STATS=DEF MODE.
```

- STATISTICS requests the default statistics (mean, standard deviation, minimum, and maximum) plus the mode of AGE.

## MISSING Subcommand

By default, both user-missing and system-missing values are labeled as missing in the table but are not included in the valid and cumulative percentages, in the calculation of descriptive statistics, or in charts and histograms.

INCLUDE	<i>Include cases with user-missing values. Cases with user-missing values are included in statistics and plots.</i>
---------	---

## ***ORDER Subcommand***

You can organize your output by variable or by analysis. Frequencies output that is organized by analysis has a single statistics table for all variables. Output organized by variable has a statistics table and a frequency table for each variable.

- ANALYSIS**      *Organize output by analysis.* Displays a single statistics table for all variables. This is the default.
- VARIABLE**     *Organize output by variable.* Displays a statistics table and a frequency table for each variable.

# GENLIN

GENLIN is available in the Advanced Statistics option.

*Note:* Equals signs (=) used in the syntax chart are required elements. All subcommands are optional.

```

GENLIN
  {dependent-var [(REFERENCE = {LAST**})] [(ORDER = {ASCENDING**})]}
                                {FIRST } {DESCENDING }
                                {value } {DATA }

  {events-var OF {trials-var} }
    {n }

  [BY factor-list [(ORDER = {ASCENDING**})]]
                                {DESCENDING }
                                {DATA }

  [WITH covariate-list]

[/MODEL [effect-list] [INTERCEPT = {YES**}]
                                {NO }

  [OFFSET = {variable}] [SCALEWEIGHT = variable]
    {value }

  [DISTRIBUTION = {BINOMIAL }
                  {GAMMA }
                  {IGAUSS }
                  {MULTINOMIAL }
                  {NEGBIN ({1** })}
                    {value}
                    {MLE }
                  {NORMAL }
                  {POISSON }
                  {TWEEDIE (value) }

  [LINK = {CLOGLOG }
          {IDENTITY }
          {LOG }
          {LOGC }
          {LOGIT }
          {NEGBIN }
          {NLOGLOG }
          {ODDSPOWER (value)}
          {POWER (value) }
          {PROBIT }
          {CUMCAUCHIT }
          {CUMCLOGLOG }
          {CUMLOGIT }
          {CUMNLOGLOG }
          {CUMPROBIT }

  [/CRITERIA [ANALYSISTYPE = {3**} [({WALD**})] [CHECKSEP = {20**}]
              {1 } {LR } {n }
              {ALL}

  [CILEVEL = {95** } [CITYPE = {WALD** }
    {value} {PROFILE [({1E-4**})]}
              {value }

  [COVB = {MODEL**} [HCONVERGE = {0** } [({ABSOLUTE**})]]
    {ROBUST } {value} {RELATIVE }

  [INITIAL = {number-list [NEGBIN({1** })]}
              {value}
    {'savfile' | 'dataset' }

```



```

[LCONVERGE = {0** } [({ABSOLUTE**})] ] [LIKELIHOOD = {FULL**}]
              {value} {RELATIVE }           {KERNEL}

[MAXITERATIONS = {100**}] [MAXSTEPHALVING = {5**}]
              {n }           {n }

[METHOD = {FISHER ({1**})}]
              {n }
              {NEWTON }

[PCONVERGE = {1E-6**} [({ABSOLUTE**})] ]
              {value } {RELATIVE }

[SCALE = {MLE** } ]
              {DEVIANC}
              {PEARSON }
              {1** }
              {value }

[SINGULAR = {1E-12**}]
              {value }

[/REPEATED SUBJECT = variable [* variable...]]

[WITHINSUBJECT = variable [* variable...]]

[SORT = {YES**}]
              {NO }

[CORRTYPE = {INDEPENDENT** } ]
              {AR(1) }
              {EXCHANGEABLE }
              {FIXED (number-list)}
              {MDEPENDENT (n) }
              {UNSTRUCTURED }

[ADJUSTCORR = {YES**}]
              {NO }

[COVB = {ROBUST**}]
              {MODEL }

[HCONVERGE = {0** } ]
              {value}

[MAXITERATIONS = {100**}]
              {n }

[PCONVERGE = {1E-6**} [({ABSOLUTE**})] ]
              {value } {RELATIVE }

[UPDATECORR = {1**}]
              {n }

[/EMMEANS [TABLES = factor [* factor ...]]

[CONTROL = variable (value) [variable (value) ...]]

[SCALE = {ORIGINAL** } ]
              {TRANSFORMED}

[COMPARE = factor [* factor ...]]

[CONTRAST = {PAIRWISE** } ]
              {DEVIATION }
              {DIFFERENCE }
              {HELMERT }
              {POLYNOMIAL ({1,2,3,...**})}
              {number-list}
              {REPEATED }
              {SIMPLE (value) }

[PADJUST = {LSD** } ]
              {BONFERRONI }

```

```

                {SEQBONFERRONI}
                {SIDAK          }
                {SEQSIDAK      }

[/EMMEANS...]

[/MISSING CLASSMISSING = {EXCLUDE**}
                        {INCLUDE  }

[/PRINT  [CORB] [COVB] [CPS**] [FIT**] [GEF] [HISTORY[({1**})]]
                {n }
                [LAGRANGE] [LMATRIX] [MODELINFO**] [SOLUTION**[(EXPONENTIATED)]]
                [SUMMARY**] [DESCRIPTIVES**] [WORKINGCORR] [NONE]]

[/SAVE [XBPRED[({varname          })]] [XBSTDERROR[({varname          })]]]
                {rootname[:{25**  }]}                {rootname[:{25**  }]}
                {integer}                {integer}
                [MEANPRED[({varname          })]]
                {rootname[:{25**  }]}
                {integer}
                [CIMEANPREDL[({varname          })]] CIMEANPREDU[({varname          })]]]
                {rootname[:{25**  }]}                {rootname[:{25**  }]}
                {integer}                {integer}
                [PREDFVAL[({varname})]] [LEVERAGE[({varname})]] [RESID[({varname})]]
                [PEARSONRESID[({varname})]] [DEVIANCERESID[({varname})]]
                [STDPEARSONRESID[({varname})]] [STDDEVIANCERESID[({varname})]]
                [LIKELIHOODRESID[({varname})]] [COOK[({varname})]]

[/OUTFILE {CORB = 'savfile' | 'dataset'} {MODEL = 'file'   }
          {COVB = 'savfile' | 'dataset'} {PARAMETER = 'file'}

```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### Release History

#### Release 15.0

- Command introduced.

#### Release 16.0

- Added multinomial and tweedie distributions; added MLE estimation option for ancillary parameter of negative binomial distribution (MODEL subcommand, DISTRIBUTION keyword). Notes related to the addition of the new distributions added throughout.
- Added cumulative Cauchit, cumulative complementary log-log, cumulative logit, cumulative negative log-log, and cumulative probit link functions (MODEL subcommand, LINK keyword).
- Added likelihood-ratio chi-square statistics as an alternative to Wald statistics (CRITERIA subcommand, ANALYSISTYPE keyword).
- Added profile likelihood confidence intervals as an alternative to Wald confidence intervals (CRITERIA subcommand, CITYPE keyword).
- Added option to specify initial value for ancillary parameter of negative binomial distribution (CRITERIA subcommand, INITIAL keyword).
- Changed default display of the likelihood function for GEEs to show the full value instead of the kernel (CRITERIA subcommand, LIKELIHOOD keyword).

### Example

```
GENLIN mydepvar BY a b c WITH x y z  
  /MODEL a b c x y z.
```

## Overview

The GENLIN procedure fits the generalized linear model and generalized estimating equations.

The generalized linear model includes one dependent variable and usually one or more independent effects. Subjects are assumed to be independent. The generalized linear model covers not only widely used statistical models such as linear regression for normally distributed responses, logistic models for binary data, and loglinear models for count data, but also many other statistical models via its very general model formulation. However, the independence assumption prohibits the model from being applied to correlated data.

Generalized estimating equations extend the generalized linear model to correlated longitudinal data and clustered data. More particularly, generalized estimating equations model correlations within subjects. Data across subjects are still assumed independent.

### Options

**Independence Assumption.** The GENLIN procedure fits either the generalized linear model assuming independence across subjects, or generalized estimating equations assuming correlated measurements within subjects but independence across subjects.

**Events/Trials Specification for Binomial Distribution.** The typical dependent variable specification will be a single variable, but for the binomial distribution the dependent variable may be specified using a number-of-events variable and a number-of-trials variable. Alternatively, if the number of trials is the same across all subjects, then trials may be specified using a fixed number instead of a variable.

**Probability Distribution of Dependent Variable.** The probability distribution of the dependent variable may be specified as normal, binomial, gamma, inverse Gaussian, multinomial, negative binomial, Poisson, or Tweedie.

**Link Function.** GENLIN offers the following link functions: Identity, complementary log-log, log, log-complement, logit, negative binomial, negative log-log, odds power, power, and probit. For the multinomial distribution, the following link functions are available: cumulative Cauchit, cumulative complementary log-log, cumulative logit, cumulative negative log-log, and cumulative probit.

**Correlation Structure for Generalized Estimating Equations.** When measurements within subjects are assumed correlated, the correlation structure may be specified as independent, AR(1), exchangeable, fixed,  $m$ -dependent, or unstructured.

**Estimated Marginal Means.** Estimated marginal means may be computed for one or more crossed factors and may be based on either the response or the linear predictor.

**Basic Specification**

- The basic specification is a `MODEL` subcommand with one or more model effects and a variable list identifying the dependent variable, the factors (if any), and the covariates (if any).
- If the `MODEL` subcommand is not specified, or is specified with no model effects, then the default model is the intercept-only model using the normal distribution and identity link.
- If the `REPEATED` subcommand is not specified, then subjects are assumed to be independent.
- If the `REPEATED` subcommand is specified, then generalized estimating equations, which model correlations within subjects, are fit. By default, generalized estimating equations use the independent correlation structure.
- The basic specification displays default output, including a case processing summary table, variable information, model information, goodness of fit statistics, model summary statistics, and parameter estimates and related statistics.

**Syntax Rules**

- The dependent variable, or an events/trials specification is required. All other variables and subcommands are optional.
- It is invalid to specify a dependent variable and an events/trials specification in the same `GENLIN` command.
- Multiple `EMMEANS` subcommands may be specified; each is treated independently. All other subcommands may be specified only once.
- The `EMMEANS` subcommand may be specified without options. All other subcommands must be specified with options.
- Each keyword may be specified only once within a subcommand.
- The command name, all subcommand names, and all keywords must be spelled in full.
- Subcommands may be specified in any order.
- Within subcommands, keyword settings may be specified in any order.
- The following variables, if specified, must be numeric: events and trials variables, covariates, `OFFSET` variable, and `SCALEWEIGHT` variable. The following, if specified, may be numeric or string variables: the dependent variable, factors, `SUBJECT` variables, and `WITHINSUBJECT` variables.
- All variables must be unique within and across the following variables or variable lists: the dependent variable, events variable, trials variable, factor list, covariate list, `OFFSET` variable, and `SCALEWEIGHT` variable.
- The dependent variable, events variable, trials variable, and covariates may not be specified as `SUBJECT` or `WITHINSUBJECT` variables.
- `SUBJECT` variables may not be specified as `WITHINSUBJECT` variables.
- The minimum syntax is a dependent variable. This specification fits an intercept-only model.

**Case Frequency**

- If an `WEIGHT` variable is specified, then its values are used as frequency weights by the `GENLIN` procedure.

- Weight values are rounded to the nearest whole numbers before use. For example, 0.5 is rounded to 1, and 2.4 is rounded to 2.
- The WEIGHT variable may not be specified on any subcommand in the GENLIN procedure.
- Cases with missing weights or weights less than 0.5 are not used in the analyses.

## Examples

### Poisson Regression

```
* Generalized Linear Models.
GENLIN damage_incidents BY type construction operation (ORDER=DESCENDING)
  /MODEL type construction operation INTERCEPT=YES OFFSET=log_months_service
    DISTRIBUTION=POISSON LINK=LOG
  /CRITERIA METHOD=FISHER(1) SCALE=PEARSON COVB=MODEL MAXITERATIONS=100
    MAXSTEPHALVING=5 PCONVERGE=1E-006 (ABSOLUTE) SINGULAR=1E-012
    ANALYSISTYPE=3 (WALD) CILEVEL=95 CITYPE=WALD LIKELIHOOD=FULL
  /EMMEANS TABLES=type SCALE=TRANSFORMED COMPARE=type CONTRAST=PAIRWISE
    PADJUST=SEQSIDAK
  /EMMEANS TABLES=construction SCALE=TRANSFORMED COMPARE=construction
    CONTRAST=PAIRWISE PADJUST=SEQSIDAK
  /MISSING CLASSMISSING=EXCLUDE
  /PRINT CPS DESCRIPTIVES MODELINFO FIT SUMMARY SOLUTION
  /SAVE XBPRED STDDEVIANCERESID.
```

- The procedure fits a model for the dependent variable *damage\_incidents*, using *type*, *construction*, and *operation* as factors.
- The model specification assumes that *damage\_incidents* has a Poisson distribution. A log link function relates the distribution of *damage\_incidents* to a linear combination of the predictors, including an intercept term, and an offset equal to the values *log\_months\_service*.
- The Pearson chi-square method is used to estimate the scale parameter. All other model fitting criteria are set to their default values.
- Estimated marginal means are computed on the scale of the linear predictor for *type* and *construction* using pairwise contrasts. The sequential Sidak method for multiple comparisons is used to adjust *p*-values.
- Print outputs are set to their default values.
- The model-estimated values of the linear predictor and the standardized deviance residual are saved to the active dataset.

### Gamma Regression

```
* Generalized Linear Models.
GENLIN claimamt BY holderage vehiclegroup vehicleage (ORDER=DESCENDING)
  /MODEL holderage vehiclegroup vehicleage INTERCEPT=YES
    SCALEWEIGHT=nclaims DISTRIBUTION=GAMMA LINK=POWER (-1)
  /CRITERIA METHOD=FISHER(1) SCALE=PEARSON COVB=MODEL MAXITERATIONS=100
    MAXSTEPHALVING=5 PCONVERGE=1E-006 (ABSOLUTE) SINGULAR=1E-012
    ANALYSISTYPE=3 (WALD) CILEVEL=95 CITYPE=WALD LIKELIHOOD=FULL
  /EMMEANS TABLES=holderage SCALE=ORIGINAL COMPARE=holderage
    CONTRAST=REPEATED PADJUST=SEQSIDAK
  /EMMEANS TABLES=vehiclegroup SCALE=ORIGINAL COMPARE=vehiclegroup
    CONTRAST=PAIRWISE PADJUST=SEQSIDAK
  /EMMEANS TABLES=vehicleage SCALE=ORIGINAL COMPARE=vehicleage
    CONTRAST=REPEATED PADJUST=SEQSIDAK
```

```

/MISSING CLASSMISSING=EXCLUDE
/PRINT CPS DESCRIPTIVES MODELINFO FIT SUMMARY SOLUTION
/SAVE XBPRED STDDEVIANCERESID.

```

- The procedure fits a model for the dependent variable *claimamt*, using *holderage*, *vehiclegroup*, and *vehicleage* as factors. The category order for all factors is descending values of factor levels.
- The model specification assumes that *claimamt* has a gamma distribution. A power link function with -1 as the exponent relates the distribution of *claimamt* to a linear combination of the predictors (including an intercept term).
- The Pearson chi-square method is used to estimate the scale parameter, with *nclaims* providing scale weights. All other model fitting criteria are set to their default values.
- Estimated marginal means are computed for *holderage*, using repeated contrasts; *vehiclegroup*, using pairwise contrasts; and *vehicleage*, using repeated contrasts. All tests are adjusted using the sequential Sidak method.
- Print outputs are set to their default values.
- The model-estimated values of the linear predictor and the standardized deviance residual are saved to the active dataset.

### **Complementary Log-log Regression for Interval-Censored Survival Data**

```

* Generalized Linear Models.
GENLIN result2 (REFERENCE=FIRST) BY id duration treatment period
  (ORDER=DESCENDING) WITH age
  /MODEL period duration treatment age INTERCEPT=NO
  DISTRIBUTION=BINOMIAL LINK=CLOGLOG
  /CRITERIA METHOD=FISHER(1) SCALE=1 COVB=MODEL MAXITERATIONS=100
  MAXSTEPHALVING=5 PCONVERGE=1E-006 (ABSOLUTE) SINGULAR=1E-012
  ANALYSISTYPE=3 (WALD) CILEVEL=95 CITYPE=WALD LIKELIHOOD=FULL
  /MISSING CLASSMISSING=EXCLUDE
  /PRINT CPS DESCRIPTIVES MODELINFO FIT SUMMARY SOLUTION.

```

- The procedure fits a model for the dependent variable *result2*, using *id* as a factor to determine subpopulations and *duration*, *treatment*, and *period* as factors to predict values, with *age* as a covariate. The first category of *result2* is used as the reference category, and the category order for all factors is descending values of factor levels.
- The model specification assumes that *result2* has a binomial distribution. A complementary log-log link function relates the probability of *result2* to a linear combination of the predictors, excluding an intercept term.
- Model fitting criteria and print output are set to their default values.

### **Repeated Measures Logistic Regression (Generalized Estimating Equation)**

```

* Generalized Estimating Equations.
GENLIN
  wheeze (REFERENCE=LAST) BY age smoker (ORDER=ASCENDING)
  /MODEL age smoker INTERCEPT=YES
  DISTRIBUTION=BINOMIAL LINK=LOGIT
  /CRITERIA METHOD=FISHER(1) SCALE=1 MAXITERATIONS=100 MAXSTEPHALVING=5
  PCONVERGE=1E-006 (ABSOLUTE) SINGULAR=1E-012 ANALYSISTYPE=3 (WALD)
  CILEVEL=95
  /REPEATED SUBJECT=id WITHINSUBJECT=age SORT=YES CORRTYPE=UNSTRUCTURED
  ADJUSTCORR=YES COVB=ROBUST MAXITERATIONS=100 PCONVERGE=1e-006 (ABSOLUTE)

```

```

UPDATECORR=1
/MISSING CLASSMISSING=EXCLUDE
/PRINT CPS DESCRIPTIVES MODELINFO FIT SUMMARY SOLUTION WORKINGCORR.

```

- The procedure fits a model for the dependent variable *wheeze*, using *smoker* and *age* as factors. The first category of *wheeze* is used as the reference category.
- The model specification assumes that *wheeze* has a binomial distribution. A logit link function relates the probability of *wheeze* to a linear combination of the predictors, including an intercept term.
- Clusters of correlated observations are defined by values of the subject variable *id*. Repeated measurements are ordered within subjects by values of *age*. An unstructured working correlation matrix is estimated.
- Model fitting criteria are set to their default values.
- The working correlation matrix is requested as output in addition to the default output.

## Variable List

The GENLIN command variable list specifies the dependent variable using either a single variable or events and trials variables. Alternatively, the number of trials may be specified as a fixed number. The variable list also specifies any factors and covariates.

If an events/trials specification is used for the dependent variable, then the GENLIN procedure automatically computes the ratio of the events variable over the trials variable or number. Technically, the procedure treats the events variable as the dependent variable in the sense that predicted values and residuals are based on the events variable rather than the events/trials ratio.

- The first specification on GENLIN must be a single dependent variable name or an events/trials specification.
- If the dependent variable is specified as a single variable, then it may be scale, an integer-valued count variable, binary, or ordinal.
- If the dependent variable is binary, then it may be numeric or string and there may be only two distinct valid data values.
- If the dependent variable is categorical, then it may be numeric or string and must have at least two distinct valid data values.
- If the dependent variable is not binary or categorical, then it must be numeric.
- The REFERENCE keyword specifies the dependent variable value to use as the reference category for parameter estimation. No model parameters are assigned to the reference category.

### REFERENCE = LAST

*The last dependent variable value is the reference category.* The last dependent variable value is defined based on the ascending order of the data values. This is the default.

If REFERENCE = LAST, then the procedure models the first value as the response, treating the last value as the reference category.

### REFERENCE = FIRST

*The first dependent variable value is the reference category.* The first dependent variable value is defined based on the ascending order of the data values.

If `REFERENCE = FIRST`, then the procedure models the last value as the response, treating the first value as the reference category.

**REFERENCE = value**

*The specified dependent variable value is the reference category.* Put the value inside a pair of quotes if it is formatted (such as date or time) or if the dependent variable is of string type; note, however, that this does not work for custom currency formats.

If `REFERENCE = value`, then the procedure models the unspecified value as the response, treating the specified value as the reference category.

- The `REFERENCE` specification is honored only if the dependent variable is binary and the binomial distribution is used (that is, `DISTRIBUTION = BINOMIAL` is specified on the `MODEL` subcommand). Otherwise, this specification is silently ignored.
- If the dependent variable is a string variable, then the value at the highest or lowest level is locale-dependent.
- If a value is specified as the reference category of the dependent variable, then the value must exist in the data.
- The `ORDER` keyword following the dependent variable is honored only if the dependent variable is categorical and the multinomial distribution is used (`/MODEL DISTRIBUTION = MULTINOMIAL`). Otherwise, this specification is silently ignored.
- `ORDER` determines the sort order of the dependent variable's values. Cumulative link functions are applied based on this order.

**ORDER = ASCENDING**

*Dependent variable values are sorted in ascending order, from the lowest value to the highest value.* This is the default.

**ORDER = DATA**

*Dependent variable values are not sorted.* The first value encountered in the data defines the first category, the last value encountered defines the last category. This option may not be specified if splits are defined on the `SPLIT FILE` command.

**ORDER = DESCENDING**

*Dependent variable values are sorted in descending order, from the highest value to the lowest value.*

- If the dependent variable is a string variable, then ascending and descending order are locale-dependent.
- If an events/trials specification is used, then the events variable must be specified first, followed by the `OF` keyword, and then the trials variable or number.
- If an events/trials specification is used, then `DISTRIBUTION = BINOMIAL` must be specified on the `MODEL` subcommand. In this case, the procedure automatically computes the ratio of the events variable over the trials variable or number.
- The events and trials variables must be numeric.
- The events variable is usually the number of successes for each case. Data values must be nonnegative integers. Cases with invalid values are not used in the analysis.



- If a trials variable is specified, data values must be positive integers, and each value must be greater than or equal to the corresponding events value for a case. Cases with invalid values are not used in the analysis. If a number is specified, then it must be a positive integer, and it must be greater than or equal to the events value for each case. Cases with invalid values are not used in the analysis.
- The events and trials options are invalid if a dependent variable name is specified.
- The names of the factors and covariates, if any, follow the dependent variable or events/trials specification. Names of factors are specified following the keyword `BY`. Names of covariates are specified following the keyword `WITH`.
- The `ORDER` specification following a list of factor variable names determines the sort order of factor values. This order is relevant for determining a factor's last level, which may be associated with a redundant parameter in the estimation algorithm.

**ORDER = ASCENDING**

*Factor variable values are sorted in ascending order, from the lowest value to the highest value. This is the default order.*

**ORDER = DATA**

*Factor variable values are not sorted.* The first value encountered in the data defines the first category; the last value encountered defines the last category.

This option may not be specified if splits are defined on the `SPLIT FILE` command.

**ORDER = DESCENDING**

*Factor variable values are sorted in descending order, from the highest value to the lowest value.*

- Covariates must be numeric, but factors can be numeric or string variables.
- Each variable may be specified only once on the variable list.
- The `OFFSET` and `SCALEWEIGHT` variables may not be specified on the `GENLIN` command variable list.
- The `SUBJECT` and `WITHINSUBJECT` variables may not be specified as dependent, events, or trials variables on the `GENLIN` command variable list.
- Cases with missing values on the dependent variable, the events or trials variable, or any covariate are not used in the analysis.

## ***MODEL Subcommand***

The `MODEL` subcommand is used to specify model effects, an offset or scale weight variable if either exists, the probability distribution, and the link function.

- The effect list includes all effects to be included in the model except for the intercept, which is specified using the `INTERCEPT` keyword. Effects must be separated by spaces or commas.
- If the multinomial distribution is used (`DISTRIBUTION = MULTINOMIAL`), then the intercept is inapplicable. The multinomial model always includes threshold parameters. The number of threshold parameters is one less than the number of dependent variable categories.
- To include a term for the main effect of a factor, enter the variable name of the factor.

- To include a term for an interaction between factors, use the keyword `BY` or an asterisk (\*) to join the factors involved in the interaction. For example, `A*B` means a two-way interaction effect of `A` and `B`, where `A` and `B` are factors. `A*A` is not allowed because factors in an interaction effect must be distinct.
- To include a term for nesting one effect within another, use a pair of parentheses. For example, `A(B)` means that `A` is nested within `B`.
- Multiple nesting is allowed. For example, `A(B(C))` means that `B` is nested within `C`, and `A` is nested within `B(C)`. When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.
- Interactions between nested effects are not valid. For example, neither `A(C)*B(C)` nor `A(C)*B(D)` is valid.
- To include a covariate term in the design, enter the variable name of the covariate.
- Covariates can be connected, but not nested, through the \* operator to form another covariate effect. Interactions among covariates such as `X1*X1` and `X1*X2` are valid, but `X1(X2)` is not.
- Factor and covariate effects can be connected only by the \* operator. Suppose `A` and `B` are factors, and `X1` and `X2` are covariates. Examples of valid factor-by-covariate interaction effects are `A*X1`, `A*B*X1`, `X1*A(B)`, `A*X1*X1`, and `B*X1*X2`.
- If the `MODEL` subcommand is not specified, or if it is specified with no model effects, then the `GENLIN` procedure fits the intercept-only model (unless the intercept is excluded on the `INTERCEPT` keyword). If the multinomial distribution is being used, then the `GENLIN` procedure fits the thresholds-only model.

### ***INTERCEPT Keyword***

The `INTERCEPT` keyword controls whether an intercept term is included in the model.

- If the multinomial distribution is in effect (`DISTRIBUTION = MULTINOMIAL`), then the `INTERCEPT` keyword is silently ignored.

<b>YES</b>	<i>The intercept is included in the model. This is the default.</i>
<b>NO</b>	<i>The intercept is not included in the model. If no model effects are defined and <code>INTERCEPT = NO</code> is specified, then a null model is fit.</i>

### ***OFFSET Keyword***

The `OFFSET` keyword specifies an offset variable or fixes the offset at a number.

- The offset variable, if specified, must be numeric.
- The offset variable may not be a dependent variable, events or trials variable, factor, covariate, `SCALEWEIGHT`, `SUBJECT`, or `WITHINSUBJECT` variable.
- Cases with missing values on the `OFFSET` variable are not used in the analysis.
- Specifying a number when `INTERCEPT = YES` is equivalent to adding a constant to the intercept.
- Specifying a number when `INTERCEPT = NO` is equivalent to fixing the intercept at the specified number.

**SCALEWEIGHT Keyword**

The `SCALEWEIGHT` keyword specifies a variable that contains omega weight values for the scale parameter.

- The scale weight variable must be numeric.
- The scale weight variable may not be a dependent variable, events or trials variable, factor, covariate, `OFFSET`, `SUBJECT`, or `WITHINSUBJECT` variable.
- Cases with scale weight values that are less than or equal to 0, or missing, are not used in the analysis.

**DISTRIBUTION Keyword**

The `DISTRIBUTION` keyword specifies the probability distribution of the dependent variable.

- The default probability distribution depends on the specification format of the dependent variable. If an events/trials specification is used, then the default distribution is `BINOMIAL`. If a single variable specification is used, then the default distribution is `NORMAL`.
- Caution must be exercised when the dependent variable has events/trials format, and the `LINK` but not the `DISTRIBUTION` keyword is used. In this condition, depending on the `LINK` specification, the default `DISTRIBUTION = BINOMIAL` may yield an improper combination of `DISTRIBUTION` and `LINK` settings.
- Also, caution must be exercised when the dependent variable has single variable format, and the `LINK` but not the `DISTRIBUTION` keyword is used. In this condition, if the dependent variable is a string then an error will result because a string variable cannot have a normal probability distribution. Moreover, depending on the `LINK` specification, the default `DISTRIBUTION = NORMAL` may yield an improper combination of `DISTRIBUTION` and `LINK` settings.
- The discussion of the `LINK` keyword below gives details about proper and improper combinations of `DISTRIBUTION` and `LINK` settings.

**BINOMIAL**     *Binomial probability distribution.* If the dependent variable is specified as a single variable, then it may be numeric or string and there may be only two distinct valid data values.

If the events and trials options are specified on the `GENLIN` command, then the procedure automatically computes the ratio of the events variable over the trials variable or number. The events variable—and the trials variable if specified—must be numeric. Data values for the events variable must be integers greater than or equal to zero. Data values for the trials variable must be integers greater than zero. For each case, the trials value must be greater than or equal to the events value. If an events value is noninteger, less than zero, or missing, then the corresponding case is not used in the analysis. If a trials value is noninteger, less than or equal to zero, less than the events value, or missing, then the corresponding case is not used in the analysis.

If the trials option specifies a number, then it must be a positive integer, and it must be greater than or equal to the events value for each case. Cases with invalid values are not used in the analysis.

This is the default probability distribution if the dependent variable is specified using events/trials format.

<b>GAMMA</b>	<i>Gamma probability distribution.</i> The dependent variable must be numeric, with data values greater than zero. If a data value is less than or equal to zero, or missing, then the corresponding case is not used in the analysis.
<b>IGAUSS</b>	<i>Inverse Gaussian probability distribution.</i> The dependent variable must be numeric, with data values greater than zero. If a data value is less than or equal to zero, or missing, then the corresponding case is not used in the analysis.
<b>MULTINO- MIAL</b>	<i>Multinomial probability distribution.</i> The dependent variable must be specified as a single variable, it may be numeric or string, and it must have at least two distinct valid data values. The dependent variable is assumed to be ordinal with values having an intrinsic ordering.
<b>NEGBIN(number   MLE)</b>	<p><i>Negative binomial probability distribution.</i> The dependent variable must be numeric, with data values that are integers greater than or equal to zero. If a data value is noninteger, less than zero, or missing, then the corresponding case is not used in the analysis.</p> <p>The option specifies the negative binomial distribution's ancillary parameter. Specify a number greater than or equal to zero to fix the parameter at the number. Specify <code>MLE</code> to use the maximum likelihood estimate of the parameter. The default value is 1.</p> <p>If the <code>REPEATED</code> subcommand is specified, then the ancillary parameter is treated as follows. (1) If the ancillary parameter is specified as a number, then it is fixed at that number for the initial generalized linear model and the generalized estimating equations. (2) If the ancillary parameter is estimated using maximum likelihood in the initial generalized linear model, then the estimate is passed to the generalized estimating equations, where it is treated as a fixed number. (3) If <code>NEGBIN(MLE)</code> is specified but the initial generalized linear model is bypassed and initial values are directly input to the generalized estimating equations (see the discussion of Initial Values and Generalized Estimating Equations in <a href="#">REPEATED Subcommand</a>), then the initial value of the ancillary parameter is passed to the generalized estimating equations, where it is treated as a fixed number.</p>
<b>NORMAL</b>	<i>Normal probability distribution.</i> The dependent variable must be numeric. This is the default probability distribution if the dependent variable is specified using single-variable format.
<b>POISSON</b>	<i>Poisson probability distribution.</i> The dependent variable must be numeric, with data values that are integers greater than or equal to zero. If a data value is noninteger, less than zero, or missing, then the corresponding case is not used in the analysis.
<b>TWEEDIE(number)</b>	<i>Tweedie probability distribution.</i> The dependent variable must be numeric, with data values greater than or equal to zero. If a data value is less than zero or missing, then the corresponding case is not used in the analysis. The required number specification is the fixed value of the Tweedie distribution's parameter. Specify a number greater than one and less than two. There is no default value.

### **LINK Keyword**

The `LINK` keyword specifies the link function. The following link functions are available.

- If the multinomial distribution is in effect (`DISTRIBUTION = MULTINOMIAL`), then only the the cumulative link functions are available. Keywords corresponding to these functions have prefix `CUM`. If a non-cumulative link function is specified for the multinomial distribution, then an error is issued.

<b>IDENTITY</b>	<i>Identity link function. <math>f(x)=x</math></i>
<b>CLOGLOG</b>	<i>Complementary log-log link function. <math>f(x)=\ln(-\ln(1-x))</math></i>
<b>LOG</b>	<i>Log link function. <math>f(x)=\ln(x)</math></i>
<b>LOGC</b>	<i>Log complement link function. <math>f(x)=\ln(1-x)</math></i>
<b>LOGIT</b>	<i>Logit link function. <math>f(x)=\ln(x / (1-x))</math></i>
<b>NEGBIN</b>	<i>Negative binomial link function. <math>f(x)=\ln(x / (x+k^{-1}))</math></i>
<b>NLOGLOG</b>	<i>Negative log-log link function. <math>f(x)=-\ln(-\ln(x))</math></i>
<b>ODDSPOWER(number)</b>	<i>Odds power link function. <math>f(x)=[(x/(1-x))^{\alpha}-1]/\alpha</math>, if <math>\alpha \neq 0</math>. <math>f(x)=\ln(x)</math>, if <math>\alpha=0</math>. <math>\alpha</math> is the required number specification and must be a real number. There is no default value.</i>
<b>POWER(number)</b>	<i>Power link function. <math>f(x)=x^{\alpha}</math>, if <math>\alpha \neq 0</math>. <math>f(x)=\ln(x)</math>, if <math>\alpha=0</math>. <math>\alpha</math> is the required number specification and must be a real number. If <math> \alpha  &lt; 2.2e-16</math>, <math>\alpha</math> is treated as 0. There is no default value.</i>
<b>PROBIT</b>	<i>Probit link function. <math>f(x)=\Phi^{-1}(x)</math>, where <math>\Phi^{-1}</math> is the inverse standard normal cumulative distribution function.</i>
<b>CUMCAUCHIT</b>	<i>Cumulative Cauchit link function. <math>f(x) = \tan(\pi (x - 0.5))</math>. May be specified only if <code>DISTRIBUTION = MULTINOMIAL</code>.</i>
<b>CUMCLOGLOG</b>	<i>Cumulative complementary log-log link function. <math>f(x)=\ln(-\ln(1-x))</math>. May be specified only if <code>DISTRIBUTION = MULTINOMIAL</code>.</i>
<b>CUMLOGIT</b>	<i>Cumulative logit link function. <math>f(x)=\ln(x / (1-x))</math>. May be specified only if <code>DISTRIBUTION = MULTINOMIAL</code>.</i>
<b>CUMNLOGLOG</b>	<i>Cumulative negative log-log link function. <math>f(x)=-\ln(-\ln(x))</math>. May be specified only if <code>DISTRIBUTION = MULTINOMIAL</code>.</i>
<b>CUMPROBIT</b>	<i>Cumulative probit link function. <math>f(x)=\Phi^{-1}(x)</math>, where <math>\Phi^{-1}</math> is the inverse standard normal cumulative distribution function. May be specified only if <code>DISTRIBUTION = MULTINOMIAL</code>.</i>

- If neither the `DISTRIBUTION` nor the `LINK` keyword is specified, then the default link function is `IDENTITY`.
- If `DISTRIBUTION` is specified but `LINK` is not, then the default setting for `LINK` depends on the `DISTRIBUTION` setting as shown in the following table.

<b>DISTRIBUTION Setting</b>	<b>Default LINK Setting</b>
NORMAL	IDENTITY
BINOMIAL	LOGIT
GAMMA	POWER(-1)
IGAUSS	POWER(-2)
MULTINOMIAL	CUMLOGIT
NEGBIN	LOG
POISSON	LOG
TWEEDIE	POWER(1-p), where p is the Tweedie distribution's parameter

- The GENLIN procedure will fit a model if a permissible combination of LINK and DISTRIBUTION specifications is given. The table below indicates the permissible LINK and DISTRIBUTION combinations. Specifying an improper combination will yield an error message.
- Note that the default setting for DISTRIBUTION is NORMAL irrespective of the LINK specification, and that not all LINK specifications are valid for DISTRIBUTION = NORMAL. Thus, if LINK is specified but DISTRIBUTION is not, then the default DISTRIBUTION = NORMAL may yield an improper combination of DISTRIBUTION and LINK settings.

**Table 94-1**  
Valid combinations of distribution and link function

Link	Distribution						
	NORMAL	BINOMIAL	GAMMA	IGAUSS	NEGBIN	POISSON	TWEEDIE
IDENTITY	X	X	X	X	X	X	X
CLOGLOG		X					
LOG	X	X	X	X	X	X	X
LOGC		X					
LOGIT		X					
NEGBIN					X		
NLOGLOG		X					
ODDSPOWER		X					
PROBIT		X					
POWER	X	X	X	X	X	X	X

*Note:* The NEGBIN link function is not available if DISTRIBUTION = NEGBIN(0) is specified.

## **CRITERIA Subcommand**

The CRITERIA subcommand controls statistical criteria for the generalized linear model and specifies numerical tolerance for checking singularity.

Note that if the REPEATED subcommand is used, then the GENLIN procedure fits generalized estimating equations, which comprise a generalized linear model and a working correlation matrix that models within-subject correlations. In this case, the GENLIN procedure first fits a generalized linear model assuming independence and uses the final parameter estimates as the initial values for the linear model part of the generalized estimating equations. (For more information, see the topic REPEATED Subcommand on p. 781.) The description of each CRITERIA subcommand keyword below is followed by a statement indicating how the keyword is affected by specification of the REPEATED subcommand.

### **ANALYSISTYPE = 3 | 1 | ALL (WALD | LR)**

*Type of analysis for each model effect.* Specify 1 for a type I analysis, 3 for type III analysis, or ALL for both. Each of these specifications computes chi-square statistics for each model effect.

Optionally, 1, 3, or ALL may be followed by WALD or LR in parentheses to specify the type of chi-square statistics to compute. WALD computes Wald statistics, LR computes likelihood-ratio statistics.

If likelihood-ratio statistics are computed, then the log-likelihood convergence criterion is used in all reduced models if type I analysis is in effect, or in all constrained models if type III analysis is in effect, irrespective of the convergence criteria used for parameter estimation in the full model. That is, for reduced or constrained models, any HCONVERGE and PCONVERGE specifications are not used, but all LCONVERGE specifications are used. (See the discussions of the HCONVERGE, PCONVERGE, and LCONVERGE keywords below.) If the log-likelihood convergence criterion is not in effect for the full model, then the reduced or constrained models use the log-likelihood convergence criterion with tolerance level 1E-4 and absolute change. The maximum number of iterations (MAXITERATIONS), maximum number of step-halvings (MAXSTEPHALVING), and starting iteration for checking complete and quasi-complete separation (CHECKSEP) are the same for reduced or constrained models as for the full model.

The default value is 3 (WALD).

If the REPEATED subcommand is specified, then the option on the ANALYSISTYPE keyword is used for the generalized estimating equations. In this case, the WALD option computes Wald statistics, but the LR option computes generalized score statistics instead of likelihood-ratio statistics. For generalized score statistics, the convergence criteria for reduced or constrained models are the same as for the full model; that is, HCONVERGE or PCONVERGE as specified on the REPEATED subcommand.

#### **CHECKSEP = integer**

*Starting iteration for checking complete and quasi-complete separation.* Specify an integer greater than or equal to zero. This criterion is not used if the value is 0. The default value is 20. This criterion is used only for the binomial or multinomial probability distributions (that is, if DISTRIBUTION = BINOMIAL or MULTINOMIAL is specified on the MODEL subcommand). For all other probability distributions, it is silently ignored.

If the CHECKSEP value is greater than 0 and the binomial or multinomial probability distribution is being used, then separation is always checked following the final iteration.

If the REPEATED subcommand is specified, then the CHECKSEP keyword is applicable only to the initial generalized linear model.

#### **CILEVEL = number**

*Confidence interval level for coefficient estimates and estimated marginal means.* Specify a number greater than or equal to 0, and less than 100. The default value is 95.

If the REPEATED subcommand is specified, then the CILEVEL keyword is applicable to any parameter that is fit in the process of computing the generalized estimating equations.

#### **CITYPE = WALD | PROFILE(number)**

*Confidence interval type.* Specify WALD for Wald confidence intervals, or PROFILE for profile likelihood confidence intervals. The default value is WALD.

PROFILE may be followed optionally by parentheses containing the tolerance level used by the two convergence criteria. The default value is 1E-4.

If the REPEATED subcommand is specified, then the CITYPE keyword is applicable only to the initial generalized linear model. For the linear model part of the generalized estimating equations, Wald confidence intervals are always used.

#### **COVB = MODEL | ROBUST**

*Parameter estimate covariance matrix.* Specify MODEL to use the model-based estimator of the parameter estimate covariance matrix, or ROBUST to use the robust estimator. The default value is MODEL.

If the REPEATED subcommand is specified, then the CRITERIA subcommand COVB keyword is silently ignored. The REPEATED subcommand COVB keyword is applicable to the linear model part of the generalized estimating equations.

**HCONVERGE = number (ABSOLUTE | RELATIVE)**

*Hessian convergence criterion.* Specify a number greater than or equal to 0, and the ABSOLUTE or RELATIVE keyword in parentheses to define the type of convergence. The number and keyword may be separated by a space character or a comma. If HCONVERGE = 0, the absolute Hessian convergence criterion will be checked with value 1E-4 after any specified convergence criteria have been satisfied. If it is not met, a warning is displayed. The default value is 0 (ABSOLUTE).

At least one of the CRITERIA subcommand keywords HCONVERGE, LCONVERGE, PCONVERGE must specify a nonzero number.

For a model with a normal distribution and identity link function, an iterative process is not used for parameter estimation. Thus, if DISTRIBUTION = NORMAL and LINK = IDENTITY on the MODEL subcommand, then the HCONVERGE keyword is silently ignored.

If the REPEATED subcommand is specified, then the CRITERIA subcommand HCONVERGE keyword is applicable only to the initial generalized linear model. The REPEATED subcommand HCONVERGE keyword is applicable to the linear model part of the generalized estimating equations.

**INITIAL = number-list | 'savfile' | 'dataset'**

*Initial values for parameter estimates.* Specify a list of numbers, an external IBM® SPSS® Statistics data file, or a currently open dataset. If a list of numbers is specified, then each number must be separated by a space character or a comma. If an external data file is specified, then the full path and filename must be given in quotes.

If the INITIAL keyword is specified, then initial values must be supplied for all parameters (including redundant parameters) in the generalized linear model. The ordering of the initial values should correspond to the ordering of the model parameters used by the GENLIN procedure. One way to determine how parameters are ordered for a given model is to run the GENLIN procedure for the model – without the INITIAL keyword – and examine the PRINT subcommand SOLUTION output.

If INITIAL is not specified, then the GENLIN procedure automatically determines the initial values.

If DISTRIBUTION = NORMAL and LINK = IDENTITY on the MODEL subcommand, then the INITIAL keyword is ignored with a warning.

If the REPEATED subcommand is specified, then the CRITERIA subcommand INITIAL keyword is applicable only to the initial generalized linear model. See the REPEATED subcommand below for a detailed discussion of initial values and generalized estimating equations.

**Initial Values Specified using a List of Numbers**

For all distributions except multinomial, if MODEL INTERCEPT = YES, then the initial values must begin with the initial value for the intercept parameter. If MODEL INTERCEPT = NO, then the initial values must begin with the initial value for the first regression parameter.

If SCALE = MLE, then the initial values must continue with the initial value for the scale parameter. If SCALE = DEVIANCE, PEARSON, or a fixed number, then a value may be given for the scale parameter but it is optional and always silently ignored.

Finally, if DISTRIBUTION = NEGBIN(MLE), then the initial values may end with an initial value for the negative binomial distribution's ancillary parameter. The initial value for this parameter must be specified as NEGBIN(number), where number is a number greater than or equal to zero. The default value is 1. If DISTRIBUTION = NEGBIN(MLE) is not in effect, then NEGBIN(number) is silently ignored.

For the multinomial distribution, the ordering of initial values is: threshold parameters, regression parameters.



Any additional unused numbers at the end of the list; that is, any numbers beyond those that are mapped to parameters, are silently ignored.

If the `SPLIT FILE` command is in effect, then the exact same list is applied to all splits. That is, each split must have the same set of parameters, and the same list is applied to each split. If the list contains too few or too many numbers for any split, then an error message is displayed.

#### Initial Values Specified using a Dataset or External SPSS Statistics Data File

If a currently open dataset or external SPSS Statistics data file is specified, then the file structure must be the same as that used in the `OUTFILE` subcommand `CORB` and `COVB` files. This structure allows the final values from one run of the `GENLIN` procedure to be saved in a `CORB` or `COVB` file and input as initial values in a subsequent run of the procedure.

In the dataset, the ordering of variables from left to right must be: `RowType_`, `VarName_`, `P1`, `P2`, ... The variables `RowType_` and `VarName_` are string variables. `P1`, `P2`, ... are numeric variables corresponding to an ordered list of the parameters. (Variable names `P1`, `P2`, ... are not required; the procedure will accept any valid variable names for the parameters. The mapping of variables to parameters is based on variable position, not variable name.) Any variables beyond the last parameter are ignored.

Initial values are supplied on a record with value 'EST' for variable `RowType_`; the actual initial values are given under variables `P1`, `P2`, ... The `GENLIN` procedure ignores all records for which `RowType_` has a value other than 'EST', as well as any records beyond the first occurrence of `RowType_` equal to 'EST'.

The required order of the intercept (if any) or threshold parameters, and regression parameters, is the same as for the list of numbers specification. However, when initial values are entered via a dataset or external data file, these parameters must always be followed by the scale parameter and then, if `DISTRIBUTION = NEGBIN`, by the negative binomial parameter.

If `SPLIT FILE` is in effect, then the variables must begin with the split-file variable or variables in the order specified on the `SPLIT FILE` command, followed by `RowType_`, `VarName_`, `P1`, `P2`, ... as above. Splits must occur in the specified dataset in the same order as in the original dataset.

#### Examples.

The following example specifies initial values using a list of numbers. Suppose factor A has three levels. The `INITIAL` keyword supplies initial value 1 for the intercept, 1.5 for the first level of factor A, 2.5 for the second level, 0 for the last level, and 3 for the covariate X.

```
GENLIN depvar BY a WITH x
      /MODEL a x
      /CRITERIA INITIAL = 1 1.5 2.5 0 3.
```

The next example outputs the final estimates from one run of the `GENLIN` procedure and inputs these estimates as the initial values in the second run.

```
GENLIN depvar BY a WITH x
      /MODEL a x
      /OUTFILE COVB = '/work/estimates.sav'.
GENLIN depvar BY a WITH x
      /MODEL a x
      /CRITERIA INITIAL = '/work/estimates.sav'.
```

#### LCONVERGE = number (ABSOLUTE | RELATIVE)

*Log-likelihood convergence criterion.* Specify a number greater than or equal to 0, and the `ABSOLUTE` or `RELATIVE` keyword in parentheses to define the type of convergence. The number and keyword may be separated by a space character or a comma. The log-likelihood convergence criterion is not used if the number is 0. The default value is 0 (`ABSOLUTE`).

At least one of the CRITERIA subcommand keywords HCONVERGE, LCONVERGE, PCONVERGE must specify a nonzero number.

If DISTRIBUTION = NORMAL and LINK = IDENTITY on the MODEL subcommand, then the LCONVERGE keyword is silently ignored.

If the REPEATED subcommand is specified, then the LCONVERGE keyword is applicable only to the initial generalized linear model.

#### **LIKELIHOOD = FULL | KERNEL**

*Form of the log-likelihood or log-quasi-likelihood function.* Specify FULL for the full function, or KERNEL for the kernel of the function. The default value is FULL.

For generalized linear models, the LIKELIHOOD keyword specifies the form of the log likelihood function. If the REPEATED subcommand is specified, then it specifies the form of the log quasi-likelihood function.

#### **MAXITERATIONS = integer**

*Maximum number of iterations.* Specify an integer greater than or equal to 0. The default value is 100. If DISTRIBUTION = NORMAL and LINK = IDENTITY on the MODEL subcommand, then the MAXITERATIONS keyword is silently ignored.

If the REPEATED subcommand is specified, then the CRITERIA subcommand MAXITERATIONS keyword is applicable only to the initial generalized linear model. The REPEATED subcommand MAXITERATIONS keyword is applicable to the linear model part of the generalized estimating equations.

#### **MAXSTEPHALVING = integer**

*Maximum number of steps in step-halving method.* Specify an integer greater than 0. The default value is 5. If DISTRIBUTION = NORMAL and LINK = IDENTITY on the MODEL subcommand, then the MAXSTEPHALVING keyword is silently ignored.

If the REPEATED subcommand is specified, then the MAXSTEPHALVING keyword is applicable only to the initial generalized linear model.

#### **METHOD = FISHER | NEWTON | FISHER(integer)**

*Model parameters estimation method.* Specify FISHER to use the Fisher scoring method, NEWTON to use the Newton-Raphson method, or FISHER(integer) to use a hybrid method.

In the hybrid method option, *integer* is an integer greater than 0 and specifies the maximum number of Fisher scoring iterations before switching to the Newton-Raphson method. If convergence is achieved during the Fisher scoring phase of the hybrid method, then additional Newton-Raphson steps are performed until convergence is achieved for Newton-Raphson too.

The default algorithm for the generalized linear model uses Fisher scoring in the first iteration and Newton-Raphson thereafter; the default value for the METHOD keyword is FISHER(1).

If DISTRIBUTION = NORMAL and LINK = IDENTITY on the MODEL subcommand, then the METHOD keyword is silently ignored.

If the REPEATED subcommand is specified, then the METHOD keyword is applicable only to the initial generalized linear model.

#### **PCONVERGE = number (ABSOLUTE | RELATIVE)**

*Parameter convergence criterion.* Specify a number greater than or equal to 0, and the ABSOLUTE or RELATIVE keyword in parentheses to define the type of convergence. The number and keyword may be separated by a space character or a comma. The parameter convergence criterion is not used if the number is 0. The default value is 1E-6 (ABSOLUTE).

At least one of the CRITERIA subcommand keywords HCONVERGE, LCONVERGE, PCONVERGE must specify a nonzero number.

If DISTRIBUTION = NORMAL and LINK = IDENTITY on the MODEL subcommand, then the PCONVERGE keyword is silently ignored.

If the REPEATED subcommand is specified, then the CRITERIA subcommand PCONVERGE keyword is applicable only to the initial generalized linear model. The REPEATED subcommand PCONVERGE keyword is applicable to the linear model part of the generalized estimating equations.

**SCALE = MLE | DEVIANCE | PEARSON | number**

*Method of fitting the scale parameter.* Specify MLE to compute a maximum likelihood estimate, DEVIANCE to compute the scale parameter using the deviance, PEARSON to compute it using the Pearson chi-square, or a number greater than 0 to fix the scale parameter.

If the MODEL subcommand specifies DISTRIBUTION = NORMAL, IGAUSS, GAMMA, or TWEEDIE then any of the SCALE options may be used. For these distributions, the default value is MLE.

If the MODEL subcommand specifies DISTRIBUTION = NEGBIN, POISSON, BINOMIAL, or MULTINOMIAL, then DEVIANCE, PEARSON, or a fixed number may be used. For these distributions, the default value is the fixed number 1.

If the REPEATED subcommand is specified, then the SCALE keyword is directly applicable only to the initial generalized linear model. For the linear model part of the generalized estimating equations, the scale parameter is treated as follows:

- If SCALE = MLE, then the scale parameter estimate from the initial generalized linear model is passed to the generalized estimating equations, where it is updated by the Pearson chi-square divided by its degrees of freedom.
- If SCALE = DEVIANCE or PEARSON, then the scale parameter estimate from the initial generalized linear model is passed to the generalized estimating equations, where it is treated as a fixed number.
- If SCALE is specified with a fixed number, then the scale parameter is also held fixed at the same number in the generalized estimating equations.

**SINGULAR = number**

*Tolerance value used to test for singularity.* Specify a number greater than 0. The default value is 1E-12.

If the REPEATED subcommand is specified, then the SINGULAR keyword is applicable to any linear model that is fit in the process of computing the generalized estimating equations.

## ***REPEATED Subcommand***

The REPEATED subcommand specifies the correlation structure used by generalized estimating equations to model correlations within subjects and controls statistical criteria in the nonlikelihood-based iterative fitting algorithm. If the REPEATED subcommand is not specified, then the GENLIN procedure fits a generalized linear model assuming independence.

### ***Initial Values and Generalized Estimating Equations***

Generalized estimating equations require initial values for the parameter estimates in the linear model. Initial values are not needed for the working correlation matrix because matrix elements are based on the parameter estimates.

The GENLIN procedure automatically supplies initial values to the generalized estimating equations algorithm. The default initial values are the final parameter estimates from the ordinary generalized linear model, assuming independence, that is fit based on the MODEL and CRITERIA subcommand specifications.

Recall that if the `REPEATED` subcommand is specified, then the `CRITERIA` subcommand `SCALE` keyword is directly applicable only to the initial generalized linear model. For the linear model part of the generalized estimating equations, the scale parameter is treated as follows.

- If `SCALE = MLE`, then the scale parameter estimate from the initial generalized linear model is passed to the generalized estimating equations, where it is updated by the Pearson chi-square divided by its degrees of freedom. Pearson chi-square is used because generalized estimating equations do not have the concept of likelihood, and hence the scale estimate cannot be updated by methods related to maximum likelihood estimation.
- If `SCALE = DEVIANCE` or `PEARSON`, then the scale parameter estimate from the initial generalized linear model is passed to the generalized estimating equations, where it is treated as a fixed number.
- If `SCALE` is specified with a fixed number, then the scale parameter is also held fixed in the generalized estimating equations.

It is possible to bypass fitting the generalized linear model and directly input initial values to the generalized estimating equations algorithm. To do this, specify the linear model as usual on the `MODEL` subcommand. Then, on the `CRITERIA` subcommand, specify initial values for the linear model on the `INITIAL` keyword and set `MAXITERATIONS = 0`.

For example, suppose factor A has three levels. The `INITIAL` keyword supplies initial value 1 for the intercept, 1.5 for the first level of factor A, 2.5 for the second level, 0 for the last level, and 3 for the covariate X. Because `MAXITERATIONS = 0`, no iterations are performed for the generalized linear model and the specified initial values are passed directly to the generalized estimating equations algorithm.

```
GENLIN depvar BY a WITH x
  /MODEL a x
    DISTRIBUTION = BINOMIAL
    LINK = LOGIT
    INITIAL = 1 1.5 2.5 0 3
    MAXITERATIONS = 0
  /REPEATED
    SUBJECT=idvar.
```

It is also possible to use a maximum likelihood estimate of the scale parameter as the initial value and to fix the scale parameter at this initial value for the generalized estimating equations. That is, we can override the default updating by the Pearson chi-square divided by its degrees of freedom. To do this, first fit a generalized linear model, estimating the scale parameter via maximum likelihood, and save the final parameter estimates in an external file (using the `OUTFILE` subcommand `CORB` or `COVB` option). Next, open this external file and copy the scale parameter estimate in full precision. Finally, fit the generalized estimating equations, using the final parameter estimates from the generalized linear model as the initial values, with `MAXITERATIONS = 0` on the `CRITERIA` subcommand and `SCALE` fixed at the scale parameter estimate on the `CRITERIA` subcommand. The following example syntax assumes that the maximum likelihood estimate of the scale parameter is 0.1234567890123456.

```
GENLIN depvar BY a WITH x
  /MODEL a x
    DISTRIBUTION = NORMAL
    LINK = LOG
```

```

/CRITERIA
  SCALE = MLE
/OUTFILE
  COVB = '/work/estimates.sav' .

GENLIN depvar BY a WITH x
/MODEL a x
  DISTRIBUTION = NORMAL
  LINK = LOG
/CRITERIA
  INITIAL = '/work/estimates.sav'
  MAXITERATIONS = 0
  SCALE = 0.1234567890123456
/REPEATED
  SUBJECT=idvar.

```

When the negative binomial distribution is used (`/MODEL DISTRIBUTION = NEGBIN`), the distribution's ancillary parameter is treated as follows:

1. If the ancillary parameter is specified as a number, then it is fixed at that number for the initial generalized linear model and the generalized estimating equations.
2. If the ancillary parameter is estimated using maximum likelihood in the initial generalized linear model, then the estimate is passed to the generalized estimating equations, where it is treated as a fixed number.
3. If `NEGBIN(MLE)` is specified but the initial generalized linear model is bypassed and initial values are directly input to the generalized estimating equations, then the initial value of the ancillary parameter is passed to the generalized estimating equations, where it is treated as a fixed number.

### ***SUBJECT* Keyword**

The `SUBJECT` keyword identifies subjects in the active dataset. Complete independence is assumed *across* subjects, but responses *within* subjects are assumed to be correlated.

- Specify a single variable or a list of variables connected by asterisks (\*) or the keyword `BY`.
- Variables may be numeric or string variables.
- The number of subjects equals the number of distinct combinations of values of the variables.
- If the active dataset is sorted by the subject variables, then all records with equal values on the subject variables are contiguous and define the measurements for one subject.
- In contrast, if the active dataset is not sorted, then the `GENLIN` procedure reads the data record by record. Each block of equal values on the subject variables defines a new subject. Please be aware that this approach may produce invalid results if all records for a subject are not contiguous in the active dataset.
- By default, the `GENLIN` procedure automatically sorts the active dataset by subject and any within-subject variables before performing analyses. See the `SORT` keyword below for more information.
- All specified variables must be unique.
- The dependent, events, trials, and `WITHINSUBJECT` variables may not be specified as `SUBJECT` variables.

- The `SUBJECT` keyword is required if the `REPEATED` subcommand is used.
- Cases with missing values for any of the subject variables are not used in the analysis.

### ***WITHINSUBJECT Keyword***

The `WITHINSUBJECT` keyword gives the within-subject or time effect. This effect defines the ordering of measurements within subjects. If some measurements do not appear in the data for some subjects, then the existing measurements are ordered and the omitted measurements are treated as missing values. If `WITHINSUBJECT` is not used, then measurements may be improperly ordered and missing values assumed for the last measurements within subjects.

- Specify a single variable or a list of variables connected by asterisks (\*) or the keyword `BY`.
- Variables may be numeric or string variables.
- The `WITHINSUBJECT` keyword is honored only if the default `SORT = YES` is in effect. The number of measurements within a subject equals the number of distinct combinations of values of the `WITHINSUBJECT` variables.
- The `WITHINSUBJECT` keyword is ignored and a warning is issued if `SORT = NO` is in effect. In this case, the `GENLIN` procedure reads the records for a subject in the order given in the active dataset.
- By default, the `GENLIN` procedure automatically sorts the active dataset by subject and any within-subject variables before performing analyses. See the `SORT` keyword below for more information.
- All specified variables must be unique.
- The dependent, events, trials, and `SUBJECT` variables may not be specified as `WITHINSUBJECT` variables.
- The `WITHINSUBJECT` keyword is not required if the data are properly ordered within each subject.
- Cases with missing values for any of the within-subject variables are not used in the analysis.

### ***SORT Keyword***

The `SORT` keyword indicates whether to sort cases in the working dataset by the subject effect and the within-subject effect.

- |            |   |
|------------|---|
| <b>YES</b> | <i>Sort cases by subject and any within-subject variables.</i> The <code>GENLIN</code> procedure sorts the active dataset before performing analyses. The subject and any within-subject variables are sorted based on the ascending sort order of their data values. If any of the variables are strings, then their sort order is locale-dependent. This is the default. This sort is temporary—it is in effect only for the duration of the <code>GENLIN</code> procedure. |
| <b>NO</b>  | <i>Do not sort cases by subject and any within-subject variables.</i> If <code>SORT = NO</code> is specified, then the <code>GENLIN</code> procedure does not sort the active dataset before performing analyses.   |

**CORRTYPE Keyword**

The CORRTYPE keyword specifies the working correlation matrix structure.

- INDEPENDENT** *Independent working correlation matrix.* This is the default working correlation matrix structure.
- AR(1)** *AR(1) working correlation matrix.*
- EXCHANGE-  
ABLE** *Exchangeable working correlation matrix.*
- FIXED(list)** *Fixed working correlation matrix.* Specify a list of numbers, with each number separated by a space character or a comma.
- The list of numbers must define a valid working correlation matrix. The number of rows and the number of columns must equal the dimension of the working correlation matrix. This dimension depends on the subject effect, the within-subject effect, whether the active dataset is sorted, and the data. The simplest way to determine the working correlation matrix dimension is to run the GENLIN procedure first for the model using the default working correlation matrix structure (instead of the FIXED structure) and examine the PRINT MODELINFO output for the working correlation matrix dimension. Then, rerun the procedure with the FIXED specification.
- Specify only the lower triangular portion of the matrix. Matrix elements must be specified row-by-row. All elements must be between 0 and 1 inclusive.
- For example, if there are three measurements per subject, then the following specification defines a 3 \* 3 working correlation matrix.

```
CORRTYPE = FIXED(0.84 0.65 0.75)
```

```
1.00 0.84 0.65
0.84 1.00 0.75
0.65 0.75 1.00
```

There is no default value for the fixed working correlation matrix.

**MDEPENDENT(integer)**

*m-dependent working correlation matrix.* Specify the value of  $m$  in parentheses as an integer greater than or equal to 0. The specified  $m$  should be less than the number of row or column levels in the working correlation matrix. If the specified  $m$  is greater than the dimension of the working correlation matrix, then  $m$  is set equal to the number of row or column levels minus 1. For example, if the dimension of the working correlation matrix is 4, then  $m$  should be 3 or less. In this case, if you specify  $m > 3$ , then  $m$  will be set equal to 3. There is no default value.

**UNSTRUC-  
TURED** *Unstructured working correlation matrix.*

**ADJUSTCORR Keyword**

The ADJUSTCORR keyword indicates whether to adjust the working correlation matrix estimator by the number of nonredundant parameters.

- YES** *Adjust the working correlation matrix estimator.* This is the default.
- NO** *Compute the working correlation matrix estimator without the adjustment.*

**COVB Keyword**

The COVB keyword specifies whether to use the robust or the model-based estimator of the parameter estimate covariance matrix for generalized estimating equations.

- |               |   |
|---------------|---|
| <b>ROBUST</b> | <i>Robust estimator of the parameter estimate covariance matrix. This is the default.</i> |
| <b>MODEL</b>  | <i>Model-based estimator of the parameter estimate covariance matrix.</i>                 |

**HCONVERGE Keyword**

The HCONVERGE keyword specifies the Hessian convergence criterion for the generalized estimating equations algorithm. For generalized estimating equations, the Hessian convergence criterion is always absolute.

- Specify a number greater than or equal to 0. If HCONVERGE = 0, the Hessian convergence criterion will be checked value 1E-4 after any specified convergence criteria have been satisfied. If it is not met, a warning is displayed. The default value is 0.
- At least one of the REPEATED subcommand keywords HCONVERGE, PCONVERGE must specify a nonzero number.

**MAXITERATIONS Keyword**

The MAXITERATIONS keyword specifies the maximum number of iterations for the generalized estimating equations algorithm.

- Specify an integer greater than or equal to 0. The default value is 100.

**PCONVERGE Keyword**

The PCONVERGE keyword specifies the parameter convergence criterion for the generalized estimating equations algorithm.

- Specify a number greater than or equal to 0, and the ABSOLUTE or RELATIVE keyword in parentheses to define the type of convergence. The number and keyword may be separated by a space character or a comma. The parameter convergence criterion is not used if the number is 0. The default value is 1E-6 (ABSOLUTE).
- At least one of the REPEATED subcommand keywords HCONVERGE, PCONVERGE must specify a nonzero number.

**UPDATECORR Keyword**

The UPDATECORR keyword specifies the number of iterations between updates of the working correlation matrix. Elements in the working correlation matrix are based on the parameter estimates, which are updated in each iteration of the algorithm. The UPDATECORR keyword specifies the iteration interval at which to update working correlation matrix elements. Specifying a value greater than 1 may reduce processing time.

- Specify an integer greater than 0.



- The working correlation matrix is not updated at all if the value is 0. In this case, the initial working correlation matrix is used throughout the estimation process.
- The default value is 1. By default, the working correlation matrix is updated after every iteration, beginning with the first.
- The `UPDATECORR` value must be less than or equal to the `REPEATED MAXITERATIONS` value.

## ***EMMEANS Subcommand***

The `EMMEANS` subcommand displays estimated marginal means of the dependent variable for all level combinations of a set of factors. Note that these are predicted, not observed, means. Estimated marginal means can be computed based on the original scale of the dependent variable or the based on the link function transformation.

- Multiple `EMMEANS` subcommands are allowed. Each is treated independently.
- The `EMMEANS` subcommand may be specified with no additional keywords. The output for an empty `EMMEANS` subcommand is the overall estimated marginal mean of the response, collapsing over any factors and holding any covariates at their overall means.
- Estimated marginal means are not available if the multinomial distribution is used. If `DISTRIBUTION = MULTINOMIAL` on the `MODEL` subcommand and the `EMMEANS` subcommand is specified, then `EMMEANS` is ignored and a warning is issued.

### ***TABLES Keyword***

The `TABLES` keyword specifies the cells for which estimated marginal means are displayed.

- Valid options are factors appearing on the `GENLIN` command factor list, and crossed factors constructed of factors on the factor list. Crossed factors can be specified using an asterisk (\*) or the keyword `BY`. All factors in a crossed factor specification must be unique.
- If the `TABLES` keyword is specified, then the `GENLIN` procedure collapses over any factors on the `GENLIN` command factor list but not on the `TABLES` keyword before computing the estimated marginal means for the dependent variable.
- If the `TABLES` keyword is not specified, then the overall estimated marginal mean of the dependent variable, collapsing over any factors, is computed.

### ***CONTROL Keyword***

The `CONTROL` keyword specifies the covariate values to use when computing the estimated marginal means.

- Specify one or more covariates appearing on the `GENLIN` command covariate list, each of which must be followed by a numeric value or the keyword `MEAN` in parentheses.
- If a numeric value is given for a covariate, then the estimated marginal means will be computed by holding the covariate at the supplied value. If the keyword `MEAN` is used, then the estimated marginal means will be computed by holding the covariate at its overall mean. If a covariate is not specified on the `CONTROL` option, then its overall mean will be used in estimated marginal means calculations.
- Any covariate may occur only once on the `CONTROL` keyword.

**SCALE Keyword**

The SCALE keyword specifies whether to compute estimated marginal means based on the original scale of the dependent variable or based on the link function transformation.

**ORIGINAL** *Estimated marginal means are based on the original scale of the dependent variable.* Estimated marginal means are computed for the response. This is the default. Note that when the dependent variable is specified using the events/trials option, ORIGINAL gives the estimated marginal means for the events/trials proportion rather than for the number of events.

**TRANSFORMED** *Estimated marginal means are based on the link function transformation.* Estimated marginal means are computed for the linear predictor.

**Example**

The following syntax specifies a logistic regression model with binary dependent variable Y and categorical predictor A. Estimated marginal means are requested for each level of A. Because SCALE = ORIGINAL is used, the estimated marginal means are based on the original response. Thus, the estimated marginal means are real numbers between 0 and 1. If SCALE = TRANSFORMED had been used instead, then the estimated marginal means would be based on the logit-transformed response and would be real numbers between negative and positive infinity.

```
GENLIN y BY a
  /MODEL a
    DISTRIBUTION=BINOMIAL
    LINK=LOGIT
  /EMMEANS TABLES=a SCALE=ORIGINAL.
```

**COMPARE Keyword**

The COMPARE keyword specifies a factor or a set of crossed factors, the levels or level combinations of which are compared using the contrast type specified on the CONTRAST keyword.

- Valid options are factors appearing on the TABLES keyword. Crossed factors can be specified using an asterisk (\*) or the keyword BY. All factors in a crossed factor specification must be unique.
- The COMPARE keyword is valid only if the TABLES keyword is also specified.
- If a single factor is specified, then levels of the factor are compared for each level combination of any other factors on the TABLES keyword.
- If a set of crossed factors is specified, then level combinations of the crossed factors are compared for each level combination of any other factors on the TABLES keyword. Crossed factors may be specified only if PAIRWISE is specified on the CONTRAST keyword.
- By default, the GENLIN procedure sorts levels of the factors in ascending order and defines the highest level as the last level. (If the factor is a string variable, then the value of the highest level is locale-dependent.) However, the sort order can be modified using the ORDER keyword following the factor list on the GENLIN command.
- Only one COMPARE keyword is allowed on a given EMMEANS subcommand.

**CONTRAST Keyword**

The **CONTRAST** keyword specifies the type of contrast to use for the levels of the factor, or level combinations of the crossed factors, on the **COMPARE** keyword. The **CONTRAST** keyword creates an **L** matrix (that is, a coefficient matrix) such that the columns corresponding to the factor(s) match the contrast given. The other columns are adjusted so that the **L** matrix is estimable.

- The **CONTRAST** keyword is valid only if the **COMPARE** keyword is also specified.
- If a single factor is specified on the **COMPARE** keyword, then any contrast type may be specified on the **CONTRAST** keyword.
- If a set of crossed factors is specified on the **COMPARE** keyword, then only the **PAIRWISE** keyword may be specified on the **CONTRAST** keyword.
- Only one **CONTRAST** keyword is allowed on a given **EMMEANS** subcommand.
- If the **COMPARE** keyword is specified without **CONTRAST**, then pairwise comparisons are performed for the factor(s) on **COMPARE**.
- **DIFFERENCE**, **HELMERT**, **REPEATED**, and **SIMPLE** contrasts are defined with respect to a first or last level. The first or last level is determined by the **ORDER** specification following the factors on the **GENLIN** command line. By default, **ORDER = ASCENDING** and the last level corresponds to the last level.

The following contrast types are available.

**PAIRWISE** *Pairwise comparisons are computed for all level combinations of the specified or implied factors. This is the default contrast type.*

For example,

```
GENLIN y BY a b c
...
/EMMEANS TABLES=a*b*c COMPARE a*b CONTRAST=PAIRWISE.
```

The specified contrast performs pairwise comparisons of all level combinations of factors A and B, for each level of factor C.

Pairwise contrasts are not orthogonal.

**DEVIATION (value)**

*Each level of the factor is compared to the grand mean. Deviation contrasts are not orthogonal.*

**DIFFERENCE** *Each level of the factor except the first is compared to the mean of previous levels. In a balanced design, difference contrasts are orthogonal.*

**HELMERT** *Each level of the factor except the last is compared to the mean of subsequent levels. In a balanced design, Helmert contrasts are orthogonal.*

**POLYNOMIAL (number list)**

*Polynomial contrasts.* The first degree of freedom contains the linear effect across the levels of the factor, the second contains the quadratic effect, and so on. By default, the levels are assumed to be equally spaced; the default metric is  $(1\ 2\ \dots\ k)$ , where  $k$  levels are involved. The `POLYNOMIAL` keyword may be followed optionally by parentheses containing a number list. Numbers in the list must be separated by spaces or commas. Unequal spacing may be specified by entering a metric consisting of one number for each level of the factor. Only the relative differences between the terms of the metric matter. Thus, for example,  $(1\ 2\ 4)$  is the same metric as  $(2\ 3\ 5)$  or  $(20\ 30\ 50)$  because, in each instance, the difference between the second and third numbers is twice the difference between the first and second. All numbers in the metric must be unique; thus,  $(1\ 1\ 2)$  is not valid.

A user-specified metric must supply at least as many numbers as there are levels of the compared factor. If too few numbers are specified, then a warning is issued and hypothesis tests are not performed. If too many numbers are specified, then a warning is issued but hypothesis tests are still performed. In the latter case, the contrast is created based on the specified numbers beginning with the first and using as many numbers as there are levels of the compared factor. In any event, we recommend printing the L matrix (`/PRINT LMATRIX`) to confirm that the proper contrast is being constructed.

For example,

```
GENLIN y BY a
...
/EMMEANS TABLES=a CONTRAST=POLYNOMIAL(1 2 4).
```

Suppose that factor A has three levels. The specified contrast indicates that the three levels of A are actually in the proportion 1:2:4.

Alternatively, suppose that factor A has two levels. In this case, the specified contrast indicates that the two levels of A are in the proportion 1:2.

In a balanced design, polynomial contrasts are orthogonal.

**REPEATED** *Each level of the factor except the last is compared to the next level.*  
Repeated contrasts are not orthogonal.

**SIMPLE (value)**

*Each level of the factor except the last is compared to the last level.*  
The `SIMPLE` keyword may be followed optionally by parentheses containing a value. Put the value inside a pair of quotes if it is formatted (such as date or currency) or if the factor is of string type. If a value is specified, then the factor level with that value is used as the omitted reference category. If the specified value does not exist in the data, then a warning is issued and the last level is used.

For example,

```
GENLIN y BY a
...
/EMMEANS TABLES=a CONTRAST=SIMPLE(1).
```

The specified contrast compares all levels of factor A (except level 1) to level 1.

Simple contrasts are not orthogonal.

**PADJUST Keyword**

The PADJUST keyword indicates the method of adjusting the significance level.

- LSD**                    *Least significant difference.* This method does not control the overall probability of rejecting the hypotheses that some linear contrasts are different from the null hypothesis value(s). This is the default.
- BONFERRONI**            *Bonferroni.* This method adjusts the observed significance level for the fact that multiple contrasts are being tested.
- SEQBONFERRONI**        *Sequential Bonferroni.* This is a sequentially step-down rejective Bonferroni procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.
- SIDAK**                    *Sidak.* This method provides tighter bounds than the Bonferroni approach.
- SEQSIDAK**              *Sequential Sidak.* This is a sequentially step-down rejective Sidak procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.

**MISSING Subcommand**

The MISSING subcommand specifies how missing values are handled.

- Cases with system missing values on any variable used by the GENLIN procedure are excluded from the analysis.
- Cases must have valid data for the dependent variable or the events and trials variables, any covariates, the OFFSET variable if it exists, the SCALEWEIGHT variable if it exists, and any SUBJECT and WITHINSUBJECT variables. Cases with missing values for any of these variables are not used in the analysis.
- The CLASSMISSING keyword specifies whether user-missing values of any factors are treated as valid.

- EXCLUDE**                *Exclude user-missing values among any factor or subpopulation variables.* Treat user-missing values for these variables as invalid data. This is the default.
- INCLUDE**                *Include user-missing values among any factor or subpopulation variables.* Treat user-missing values for these variables as valid data.

**PRINT Subcommand**

The PRINT subcommand is used to display optional output.

- If the `PRINT` subcommand is not specified, then the default output indicated below is displayed.
- If the `PRINT` subcommand is specified, then the `GENLIN` procedure displays output only for those keywords that are specified.

**CORB**            *Correlation matrix for parameter estimates.*  
**COVB**            *Covariance matrix for parameter estimates.*  
**CPS**             *Case processing summary.* For generalized estimating equations, this keyword also displays the Correlated Data Summary table. This is the default output if the `PRINT` subcommand is not specified.

#### **DESCRIPTIVES**

*Descriptive statistics.* Displays descriptive statistics and summary information about the dependent variable, covariates, factors. This is the default output if the `PRINT` subcommand is not specified.

**FIT**             *Goodness of fit.* For generalized linear models, displays deviance and scaled deviance, Pearson chi-square and scaled Pearson chi-square, log likelihood, Akaike's information criterion (AIC), finite sample corrected AIC (AICC), Bayesian information criterion (BIC), and consistent AIC (CAIC).

Note that when the scale parameter is fit using the deviance (`/CRITERIA SCALE = DEVIANCE`) or Pearson chi-square (`/CRITERIA SCALE = PEARSON`), the algorithm begins by assuming the scale parameter equals 1. Following estimation of the regression coefficients, the estimated scale parameter is calculated. Finally, estimated standard errors, Wald confidence intervals, and significance tests are adjusted based on the estimated scale parameter. However, in order to ensure fair comparison in the information criteria and the model fit omnibus test (see the `SUMMARY` keyword below), the log likelihood is not revised by the estimated scale parameter. Instead, when the scale parameter is fit using the deviance or Pearson chi-square, the log likelihood is computed with the scale parameter set equal to 1.

For generalized estimating equations, displays two extensions of AIC for model selection: Quasi-likelihood under the independence model criterion (QIC) for choosing the best correlation structure, and corrected quasi-likelihood under the independence model criterion (QICC) for choosing the best subset of predictors.

The quasi-likelihood functions are computed with the scale parameter set equal to a fixed value if a fixed value is specified on the `/CRITERIA SCALE` keyword. Otherwise, if `/CRITERIA SCALE = MLE, DEVIANCE, or PEARSON`, then the quasi-likelihood functions are computed with the scale parameter set equal to 1.

Goodness of fit statistics are not available for generalized estimating equations when the multinomial distribution is used. Thus, if the `REPEATED` subcommand and `/MODEL DISTRIBUTION = MULTINOMIAL` are specified, then the `FIT` keyword is silently ignored. This is the default output if the `PRINT` subcommand is not specified.

**GEF**             *General estimable function.*

**HISTORY (integer)**

*Iteration history.* For generalized linear models, displays the iteration history for the parameter estimates and log-likelihood, and prints the last evaluation of the gradient vector and the Hessian matrix. Also displays the iteration history for the profile likelihood confidence intervals (if requested via `CRITERIA CTYPE = PROFILE`) and for type I or III analyses (if requested via `PRINT SUMMARY`).

For generalized estimating equations, displays the iteration history for the parameter estimates, and prints the last evaluation of the generalized gradient and the Hessian matrix. Also displays the iteration history for type III analyses (if requested via `PRINT SUMMARY`).

The `HISTORY` keyword may be followed optionally by an integer  $n$  in parentheses, where the integer is greater than zero. The iteration history table displays parameter estimates for every  $n$  iterations beginning with the 0th iteration (the initial estimates). The default is to print every iteration ( $n = 1$ ). If `HISTORY` is specified, then the last iteration is always displayed regardless of the value of  $n$ .

**LAGRANGE** *Lagrange multiplier test.* For the normal, gamma, inverse Gaussian, and Tweedie distributions, displays Lagrange multiplier test statistics for assessing the validity of a scale parameter that is computed using the deviance or Pearson chi-square, or set at a fixed number. For the negative binomial distribution, tests the fixed ancillary parameter.

The `LAGRANGE` keyword is honored if `MODEL DISTRIBUTION = NORMAL, GAMMA, IGAUSS, or TWEEDIE` and `CRITERIA SCALE = DEVIANCE, PEARSON, or number`; or if `MODEL DISTRIBUTION = NEGBIN(number)` is specified. Otherwise the keyword is ignored and a warning is issued.

If the `REPEATED` subcommand is specified, then the `LAGRANGE` keyword is silently ignored.

**LMATRIX** *Set of contrast coefficient (L) matrices.* Displays contrast coefficients for the default effects and for the estimated marginal means if requested.

**MODELINFO** *Model information.* Displays the dataset name, dependent variable or events and trials variables, offset variable, scale weight variable, probability distribution, and link function.

For generalized estimating equations, also displays the subject variables, within-subject variables, and working correlation matrix structure.

This is the default output if the `PRINT` subcommand is not specified.

**SOLUTION** *Parameter estimates and corresponding statistics.* This is the default output if the `PRINT` subcommand is not specified.

The `SOLUTION` keyword may be followed optionally by the keyword `EXPONENTIATED` in parentheses to display exponentiated parameter estimates in addition to the raw parameter estimates.

**SUMMARY** *Model summary statistics.* Displays model fit tests, including likelihood ratio statistics for the model fit omnibus test, and statistics for the type I or III contrasts for each effect (depending on the `CRITERIA ANALYSISTYPE` specification). This is default output if the `PRINT` subcommand is not specified.

If the `REPEATED` subcommand is specified, then only the statistics for each effect are displayed.

**WORKINGCORR**

- NONE** *Working correlation matrix.* This keyword is honored only if the REPEATED is in effect. Otherwise it is silently ignored.
- No PRINT subcommand output.* None of the PRINT subcommand output is displayed. If NONE is specified, then no other keywords are allowed on the PRINT subcommand.

## SAVE Subcommand

The SAVE subcommand adds predicted, residual, leverage, or Cook's distance values to the working dataset.

- Specify one or more temporary variables, each followed by an optional new name in parentheses.
- The optional names must be unique, valid variable names.
- If new names are not specified, then GENLIN uses the default names. If the default names conflict with existing variable names, then a suffix is added to the default names to make them unique.

The following rules describe the functionality of the SAVE subcommand when the response variable—either the dependent variable or the events or trials variable—has an invalid value for a case.

- If all factors and covariates in the model have valid values for the case, then the procedure computes predicted values but not the residuals. (The MISSING subcommand setting is taken into account when defining valid/invalid values for a factor.)
- An additional restriction for factors is that only those values of the factor actually used in building the model are considered valid. For example, suppose factor A takes values 1, 2, and 3 when the procedure builds the model. Also suppose there is a case with an invalid dependent variable value, a value of 4 on factor A, and valid values on all other factors and covariates. For this case, no predicted value is saved because there is no model coefficient corresponding to factor A = 4.

### **XBPREP (varname | rootname:n)**

*Predicted value(s) of the linear predictor.* For all distributions except the multinomial, XBPREP creates one variable and the default variable name is *XBPredicted*. Specify a variable name in parentheses to override the default.

For the multinomial distribution, one variable is created for each dependent variable category except the last (see the dependent variable ORDER keyword in the section [Variable List](#)). XBPREP saves the predicted values of the linear predictor for the first 25 categories, up to but not including the last, by default. The default root name is *XBPredicted*, and the default variable names are *XBPredicted\_1*, *XBPredicted\_2*, and so on, corresponding to the order of the dependent variable categories. Specify a root name in parentheses to override the default. Specify a colon and a positive integer giving the number of categories to override the default 25. To specify a number without a root name, simply enter a colon before the number.

### **XBSTDERROR (varname | rootname:n)**

*Estimated standard error(s) of the predicted value of the linear predictor.* For all distributions except the multinomial, XBSTDERROR creates one variable and the default variable name is *XBStandardError*. Specify a variable name in parentheses to override the default.



For the multinomial distribution, one variable is created for each dependent variable category except the last (see the dependent variable `ORDER` keyword in the section [Variable List](#)). `XBSTDERROR` saves the estimated standard errors for the first 25 categories, up to but not including the last, by default. The default root name is `XBStandardError`, and the default variable names are `XBStandardError_1`, `XBStandardError_2`, and so on, corresponding to the order of the dependent variable categories. Specify a root name in parentheses to override the default. Specify a colon and a positive integer giving the number of categories to override the default 25. To specify a number without a root name, simply enter a colon before the number.

#### **MEANPRED (varname | rootname:n)**

*Predicted value(s) of the mean of the response.* For all distributions except the multinomial, `MEANPRED` creates one variable and the default variable name is `MeanPredicted`. Specify a variable name in parentheses to override the default.

If the binomial distribution is used and the dependent variable is in single variable format, then `MEANPRED` computes a predicted probability. Suppose the dependent variable has data values 0 and 1. If the default reference category is in effect, that is, `REFERENCE = LAST` on the `GENLIN` command line, then 1 is the reference category and `MEANPRED` computes the predicted probability that the dependent variable equals 0. To compute the predicted probability that the dependent variable equals 1 instead, specify `REFERENCE = FIRST` on the `GENLIN` command line.

If the binomial distribution is used and the dependent variable is in events/trials format, then `MEANPRED` computes the predicted number of events.

For the multinomial distribution, one variable is created for each dependent variable category except the last (see the dependent variable `ORDER` keyword in the section [Variable List](#)). `MEANPRED` saves the cumulative predicted probability for the first 25 categories, up to but not including the last, by default. The default root name is `CumMeanPredicted`, and the default variable names are `CumMeanPredicted_1`, `CumMeanPredicted_2`, and so on, corresponding to the order of the dependent variable categories. Specify a root name in parentheses to override the default. Specify a colon and a positive integer giving the number of categories to override the default 25. To specify a number without a root name, simply enter a colon before the number.

#### **CIMEANPREDL (varname | rootname:n)**

*Lower bound(s) of the confidence interval for the mean of the response.* For all distributions except the multinomial, `CIMEANPREDL` creates one variable and the default variable name is `CIMeanPredictedLower`. Specify a variable name in parentheses to override the default.

For the multinomial distribution, one variable is created for each dependent variable category except the last (see the dependent variable `ORDER` keyword in the section [Variable List](#)). `CIMEANPREDL` saves the lower bound of the cumulative predicted probability for the first 25 categories, up to but not including the last, by default. The default root name is `CICumMeanPredictedLower`, and the default variable names are `CICumMeanPredictedLower_1`, `CICumMeanPredictedLower_2`, and so on, corresponding to the order of the dependent variable categories. Specify a root name in parentheses to override the default. Specify a colon and a positive integer giving the number of categories to override the default 25. To specify a number without a root name, simply enter a colon before the number.

#### **CIMEANPREDU (varname | rootname:n)**

*Upper bound(s) of the confidence interval for the mean of the response.* For all distributions except the multinomial, `CIMEANPREDU` creates one variable and the default variable name is `CIMeanPredictedUpper`. Specify a variable name in parentheses to override the default.

For the multinomial distribution, one variable is created for each dependent variable category except the last (see the dependent variable `ORDER` keyword in the section [Variable List](#)). `CIMEANPREDU` saves the upper bound of the cumulative predicted probability for the first 25 categories, up to but not including the last, by default. The default root name is `CICumMeanPredictedUpper`, and the default variable names are `CICumMeanPredictedUpper_1`, `CICumMeanPredictedUpper_2`, and so on, corresponding to the order of the dependent variable categories. Specify a root name in parentheses to override the default. Specify a colon and a positive integer giving the number of categories to override the default 25. To specify a number without a root name, simply enter a colon before the number.

**PREDVAL (varname)**

*Predicted category value for binomial or multinomial distribution.* The class or value predicted by the model if the binomial or multinomial distribution is in effect. This keyword is honored only if the binomial distribution is used, that is, if `DISTRIBUTION = BINOMIAL` is specified or implied on the `MODEL` subcommand and the dependent variable is in single variable format, or the multinomial distribution is used (`DISTRIBUTION = MULTINOMIAL`). Otherwise, the `PREDVAL` keyword is ignored with a warning. The default variable name is *PredictedValue*.

**LEVERAGE (varname)**

*Leverage value.* The default variable name is *Leverage*. Leverage values are not available for the multinomial distribution or generalized estimating equations.

**RESID (varname)**

*Raw residual.* The default variable name is *Residual*. Raw residuals are not available for the multinomial distribution.

**PEARSONRESID (varname)**

*Pearson residual.* The default variable name is *PearsonResidual*. Pearson residuals are not available for the multinomial distribution.

**DEVIANCERESID (varname)**

*Deviance residual.* The default variable name is *DevianceResidual*. Deviance residuals are not available for the multinomial distribution or generalized estimating equations.

**STDPEARSONRESID (varname)**

*Standardized Pearson residual.* The default variable name is *StdPearsonResidual*. Standardized Pearson residuals are not available for the multinomial distribution or generalized estimating equations.

**STDDEVIANCERESID (varname)**

*Standardized deviance residual.* The default variable name is *StdDevianceResidual*. Standardized deviance residuals are not available for the multinomial distribution or generalized estimating equations.

**LIKELIHOODRESID (varname)**

*Likelihood residual.* The default variable name is *LikelihoodResidual*. Likelihood residuals are not available for the multinomial distribution or generalized estimating equations.

**COOK (varname)**

*Cook's distance.* The default variable name is *CooksDistance*. Cook's distances are not available for the multinomial distribution or generalized estimating equations.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand saves an external IBM® SPSS® Statistics data file or previously declared dataset containing the parameter correlation or covariance matrix with parameter estimates, standard errors, significance values, and degrees of freedom. It also saves the parameter estimates and the parameter covariance matrix in XML format.

- At least one keyword and a filename are required.
- The `COVB` and `CORB` keywords are mutually exclusive, as are the `MODEL` and `PARAMETER` keywords.
- The filename must be specified in full. `GENLIN` does not supply an extension.

**COVB = 'savfile' | 'dataset'**

*Writes the parameter covariance matrix and other statistics to an external SPSS Statistics data file or previously declared dataset.*

**CORB = 'savfile' | 'dataset'**

*Writes the parameter correlation matrix and other statistics to an external SPSS Statistics data file or previously declared dataset..*

**MODEL = 'file'**

*Writes the parameter estimates and the parameter covariance matrix to an XML file.*

**PARAMETER = 'file'**

*Writes the parameter estimates to an XML file.*

# GENLINMIXED

## Syntax Chart

GENLINMIXED is available in the Advanced Models option.

```
GENLINMIXED
  [/DATA_STRUCTURE]
    SUBJECTS = fieldName[*fieldName[*...]]
    [REPEATED_MEASURES = fieldName[*fieldName[*...]]]
    [GROUPING = fieldName[*fieldName[*...]]]
    [COVARIANCE_TYPE = DIAGONAL** | AR1 | ARMA11 | COMPOUND_SYMMETRY | IDENTITY
    TOEPLITZ | UNSTRUCTURED | VARIANCE_COMPONENTS]
  /FIELDS
    TARGET = fieldName
    [TRIALS = {NONE** }
    {VALUE(number) }
    {FIELD(fieldName)}]
    [ANALYSIS_WEIGHT = fieldName]
    [OFFSET = {NONE** }
    {VALUE(number) }
    {FIELD(fieldName)}]
  [/TARGET_OPTIONS]
    [REFERENCE = value]
    [DISTRIBUTION = NORMAL | BINOMIAL | MULTINOMIAL | GAMMA |
    INVERSE_GAUSSIAN | NEGATIVE_BINOMIAL | POISSON]
    [LINK = IDENTITY | CLOGLOG | LOG | LOGC | LOGIT | NLOGLOG | PROBIT | POWER]
    [LINK_PARAMETER = number]
  [/FIXED]
    [EFFECTS = effect [effect ...]]
    [USE_INTERCEPT = TRUE** | FALSE]
  [/RANDOM]
    [EFFECTS = effect [effect ...]]
    [USE_INTERCEPT = TRUE | FALSE**]
    [SUBJECTS = fieldName[*fieldName[*...]]]
    [GROUPING = fieldName[*fieldName[*...]]]
    [COVARIANCE_TYPE = DIAGONAL | AR1 | ARMA11 | COMPOUND_SYMMETRY | IDENTITY
    TOEPLITZ | UNSTRUCTURED | VARIANCE_COMPONENTS**]
  [/RANDOM ...]
  [/BUILD_OPTIONS]
    [TARGET_CATEGORY_ORDER = {ASCENDING**}
    {DESCENDING }
    {DATA }]
    [INPUTS_CATEGORY_ORDER = {ASCENDING**}
    {DESCENDING }
    {DATA }]
    [MAX_ITERATIONS = {100** }
    {number}]
    [CONFIDENCE_LEVEL = {95** }
    {number}]
    [DF_METHOD = {RESIDUAL** }
    {SATERTHWAITE}]
    [COVB = {MODEL**}
    {ROBUST }]
  [/EMMEANS]
    [TABLES = factor[*factor...]]
    [COMPARE = factor]
    [CONTRAST = {NONE** }
    {PAIRWISE }
    {DEVIATION}
    {SIMPLE }]
  [/EMMEANS...]
  [/EMMEANS_OPTIONS]
    [CONTROL = fieldName (number) [fieldName (number) ...]]
    [SCALE = {ORIGINAL** }
    {TRANSFORMED}]
    [PADJUST = {LSD** }
    {SEQBONFERRONI}
    {SEQSIDAK }]
```

```

[/OUTFILE MODEL = 'file']
[/SAVE [PREDICTED_VALUES(fieldName)]]
[PREDICTED_PROBABILITY(rootname)]
[MAX_CATEGORIES({25**  })]
           {integer}
[CONFIDENCE(fieldName)]
[CONFIDENCE_METHOD({PREDICTEDPROBABILITY**})]
           {PROBABILITYIMPROVEMENT}
[CONFIDENCE_INTERVALS(rootname)]
[PEARSON_RESIDUALS(fieldName)].

```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 19

- Command introduced.

### **Example**

```

GENLINMIXED
  /FIELDS TARGET=y1.

```

## **Overview**

The GENLINMIXED procedure fits generalized linear mixed models. Generalized linear mixed models extend the linear model so that:

- The target is linearly related to the factors and covariates via a specified link function
- The target can have a non-normal distribution
- The observations can be correlated.

Generalized linear mixed models cover a wide variety of models, from simple linear regression to complex multilevel models for non-normal longitudinal data.

### **Options**

**Probability Distribution of Target.** The probability distribution of the target can be specified as normal, binomial, gamma, inverse Gaussian, multinomial, negative binomial, or Poisson.

**Events/Trials Target Specification.** The typical target specification will be a single field, but for the binomial distribution the target can be specified using a number-of-events field and a number-of-trials field. Alternatively, if the number of trials is the same across all subjects, then trials can be specified using a fixed number instead of a field.

**Link Function.** The target is linearly related to the model effects via one of the following link functions: Identity, complementary log-log, log, log-complement, logit, negative log-log, power, and probit.

**Repeated Measurements.** You can specify the subject structure for repeated measurements and how the errors of the repeated measurements are correlated. Complete independence is assumed across subject blocks.

**Random Effects.** Random-effects are effects whose values in the data file can be considered a random sample from a larger population of values. They are useful for explaining excess variability in the target.

**Estimated Marginal Means.** Estimated marginal means can be computed for one or more crossed factors and can be based on either the response or the linear predictor.

**Output.** The procedure produces a model viewer object containing tables and charts for assessing the quality of the model; also you can save predicted values to the active dataset and the model to a file in PMML format.

### ***Basic Specification***

The basic specification is the `GENLINMIXED` command and the `FIELDS` subcommand with the `TARGET` keyword. This fits an intercept-only model that can be used as a basis for comparing more complicated models.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### ***Syntax Rules***

- The `FIELDS` subcommand is required. All other subcommands are optional.
- Subcommands can be specified in any order.
- There can be multiple `RANDOM` and `EMMEANS` subcommands; each `RANDOM` subcommand defines a separate random effects block, while each `EMMEANS` subcommand defines a separate set of estimated means. Only a single instance of any other subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not allowed.
- Any split field defined on the `SPLIT FILE` command cannot be used on this command.
- Any field specified on the `WEIGHT` command cannot be used on this command.

### ***Limitations***

- The `SPLIT FILE` setting is ignored with a warning by the procedure.

## Examples

### Common Distribution and Link Function Combinations

#### Linear model

```
GENLINMIXED
  /FIELDS TARGET=y
  /TARGET_OPTIONS DISTRIBUTION=NORMAL LINK=IDENTITY
  /FIXED EFFECTS=x1 x2 x3.
```

- The `FIELDS` subcommand specifies  $y$  as the target.
- The `TARGET_OPTIONS` subcommand that the target has a normal distribution and is linearly related to the model effects.
- The `FIXED` subcommand specifies a main effects model with fields  $x1$ ,  $x2$ , and  $x3$ . If they are continuous, they will be treated as covariates, if categorical, they will be treated as factors.

#### Gamma regression

```
GENLINMIXED
  /FIELDS TARGET=claim_amount ANALYSISWEIGHT=number_of_claims
  /TARGET_OPTIONS DISTRIBUTION=GAMMA LINK=POWER(-1)
  /FIXED EFFECTS=holder_age vehicle_group vehicle_age
  /BUILD_OPTIONS INPUTS_CATEGORY_ORDER=DESCENDING.
```

- The procedure fits a model for the target *claim\_amount*, using *holder\_age*, *vehicle\_group*, and *vehicle\_age* as main effects. In order to account for the varying number of claims used to compute the average claim amounts, you specify *number\_of\_claims* as the analysis weight.
- The `TARGET_OPTIONS` specification assumes that *claim\_amount* has a gamma distribution. A power link function with  $-1$  as the exponent relates the distribution of *claim\_amount* to a linear combination of the predictors, including an intercept term.
- The `BUILD_OPTIONS` subcommand specifies that the category order for factors is descending values of factor levels; thus, the first category of each categorical field is used as the reference category.

To compare this specification to the `GENLIN` procedure, see [Examples](#) on p. 767

#### (Poisson) Loglinear model

```
GENLINMIXED
  /FIELDS TARGET=damage_incidents OFFSET=log_months_service
  /TARGET_OPTIONS DISTRIBUTION=POISSON LINK=LOG
  /FIXED EFFECTS=type construction operation
  /BUILD_OPTIONS INPUTS_CATEGORY_ORDER=DESCENDING.
```

- The procedure fits a model for the target *damage\_incidents*, using *type*, *construction*, and *operation* as main effects.

- The `TARGET_OPTIONS` specification assumes that *damage\_incidents* has a Poisson distribution. A log link function relates the distribution of *damage\_incidents* to a linear combination of the predictors, including an intercept term, and an offset equal to the values of *log\_months\_service*.
- The `BUILD_OPTIONS` subcommand specifies that the category order for factors is descending values of factor levels; thus, the first category of each categorical field is used as the reference category.

To compare this specification to the `GENLIN` procedure, see [Examples](#) on p. 767

### **Negative binomial regression**

```
GENLINMIXED
  /FIELDS TARGET=damage_incidents OFFSET=log_months_service
  /TARGET_OPTIONS DISTRIBUTION=NEGATIVE_BINOMIAL LINK=LOG
  /FIXED EFFECTS=type construction operation.
```

- The negative binomial distribution is an alternative to the Poisson when the observations are overdispersed; that is, since the mean and variance of the Poisson distribution are the same, when the data show greater variability, the negative binomial distribution can provide a better fit.

### **Multinomial logistic regression**

```
GENLINMIXED
  /FIELDS TARGET=bfast
  /TARGET_OPTIONS DISTRIBUTION=MULTINOMIAL LINK=LOGIT
  /FIXED EFFECTS=agecat gender active.
```

- The procedure fits a model for *bfast* using *agecat*, *gender*, and *active* as main effects.
- The `TARGET_OPTIONS` specification assumes that *bfast* has a multinomial distribution. A logit link function relates the distribution of *bfast* to a linear combination of the predictors, including an intercept term.

To compare this specification to the `NOMREG` procedure, see [Examples](#) on p. 1342

### **Binary logistic regression**

```
GENLINMIXED
  /FIELDS TARGET=default
  /TARGET_OPTIONS DISTRIBUTION=BINOMIAL LINK=LOGIT
  /FIXED EFFECTS=age ed employ address income debtinc creddebt othdebt.
```

- The procedure fits a model for *default* using *age*, *ed*, *employ*, *address*, *income*, *debtinc*, *creddebt*, and *othdebt* as main effects.
- The `TARGET_OPTIONS` specification assumes that *default* has a binomial distribution. A logit link function relates the distribution of *default* to a linear combination of the predictors, including an intercept term.

To compare this specification to the `LOGISTIC REGRESSION` procedure, see [Examples](#) on p. 1031



**Binary probit model**

```

GENLINMIXED
  /FIELDS TARGET=response TRIALS=nsubj
  /TARGET_OPTIONS DISTRIBUTION=BINOMIAL LINK=PROBIT
  /FIXED EFFECTS=site value.

```

- The procedure fits a model for the number of responses *response* within trials *nsubj* using site and value as main effects.
- The TARGET\_OPTIONS specification assumes that *response* within *nsubj* has a binomial distribution. A probit link function relates the distribution of the target to a linear combination of the predictors, including an intercept term.

To compare this specification to the PROBIT procedure, see [Examples](#) on p. 1605

**Interval censored survival**

```

GENLINMIXED
  /FIELDS TARGET=result2
  /TARGET_OPTIONS DISTRIBUTION=BINOMIAL LINK=CLOGLOG
  /FIXED EFFECTS=duration treatment period age USE_INTERCEPT=FALSE
  /BUILD_OPTIONS TARGET_CATEGORY_ORDER=DESCENDING INPUTS_CATEGORY_ORDER=DESCENDING.

```

- The procedure fits a model for the target *result2*, using *duration*, *treatment*, *period*, and *age* as main effects.
- The BUILD\_OPTIONS subcommand specifies that the category order for the target and all factors is descending values of factor levels; thus, the first category of each categorical field is used as the reference category.
- The TARGET\_OPTIONS specification assumes that *result2* has a binomial distribution. A complementary log-log link function relates the probability of *result2* to a linear combination of the predictors, excluding an intercept term.

To compare this specification to the GENLIN procedure, see [Examples](#) on p. 767

**Linear mixed model**

\*Generalized Linear Mixed Models.

```

GENLINMIXED
  /DATA_STRUCTURE SUBJECTS=school*classroom*student_id
  /FIELDS TARGET=posttest TRIALS=NONE OFFSET=NONE
  /TARGET_OPTIONS DISTRIBUTION=NORMAL LINK=IDENTITY
  /FIXED EFFECTS=school_setting school_type teaching_method n_student gender lunch pretest
  USE_INTERCEPT=TRUE
  /RANDOM USE_INTERCEPT=TRUE SUBJECTS=school COVARIANCE_TYPE=VARIANCE_COMPONENTS
  /RANDOM USE_INTERCEPT=TRUE SUBJECTS=school*classroom COVARIANCE_TYPE=VARIANCE_COMPONENTS
  /BUILD_OPTIONS TARGET_CATEGORY_ORDER=ASCENDING INPUTS_CATEGORY_ORDER=ASCENDING MAX_ITERATIONS=
  CONFIDENCE_LEVEL=95 DF_METHOD=RESIDUAL COVB=MODEL
  /EMMEANS_OPTIONS SCALE=ORIGINAL PADJUST=LSD.

```

- The DATA\_STRUCTURE subcommand specifies that subjects are defined by *school*, *classroom*, and *student\_id*.
- The FIELDS and TARGET\_OPTIONS subcommands specify that the procedure creates a model for *posttest* using a normal distribution to fit the test score and an identity link to relate the target to a linear combination of the predictors.

- The FIXED subcommand specifies a model with *school\_setting*, *school\_type*, *teaching\_method*, *n\_student*, *gender*, *lunch*, and *pretest* as main effects.
- The first RANDOM subcommand specifies an intercept-only random effect block with *school* as the subject field. This should account for correlation between classrooms within the same school.
- The second RANDOM subcommand specifies an intercept-only random effect block with *school\*classroom* as the subject field. This should account for correlation between students within the same classroom.
- All other options are set to their default values.

### ***Poisson loglinear mixed model***

\*Generalized Linear Mixed Models.

```
GENLINMIXED
  /DATA_STRUCTURE SUBJECTS=patient_id
  /FIELDS TARGET=convulsions TRIALS=NONE OFFSET=NONE
  /TARGET_OPTIONS DISTRIBUTION=POISSON LINK=LOG
  /FIXED EFFECTS=after_t treatment*after_t USE_INTERCEPT=TRUE
  /RANDOM EFFECTS=after_t USE_INTERCEPT=TRUE SUBJECTS=patient_id COVARIANCE_TYPE=UNSTRUCTURED
  /BUILD_OPTIONS TARGET_CATEGORY_ORDER=ASCENDING INPUTS_CATEGORY_ORDER=ASCENDING MAX_ITERATIONS=100
  CONFIDENCE_LEVEL=95 DF_METHOD=RESIDUAL COVB=MODEL
  /EMMEANS_OPTIONS SCALE=ORIGINAL PADJUST=LSD.
```

- The DATA\_STRUCTURE subcommand specifies that subjects are defined by *patient\_id*.
- The FIELDS and TARGET\_OPTIONS subcommands specify that the procedure creates a model for *convulsions* using a Poisson distribution to fit the number of convulsions and a log link to relate the target to a linear combination of the predictors.
- The FIXED subcommand specifies a model with *after\_t* and *treatment\*after\_t* as effects.
- The RANDOM subcommand specifies *after\_t* and an intercept as effects in a random effect block with *patient\_id* as the subject field. This should account for correlation between repeated observations of the same patient.
- All other options are set to their default values.

### ***Multinomial logistic mixed model***

\*Generalized Linear Mixed Models.

```
GENLINMIXED
  /DATA_STRUCTURE SUBJECTS=customer_id
  /FIELDS TARGET=service_usage TRIALS=NONE OFFSET=NONE
  /TARGET_OPTIONS DISTRIBUTION=MULTINOMIAL LINK=LOGIT
  /FIXED EFFECTS=edcat inccat reside service_type USE_INTERCEPT=TRUE
  /RANDOM USE_INTERCEPT=TRUE SUBJECTS=customer_id COVARIANCE_TYPE=VARIANCE_COMPONENTS
  /BUILD_OPTIONS TARGET_CATEGORY_ORDER=ASCENDING INPUTS_CATEGORY_ORDER=ASCENDING MAX_ITERATIONS=100
  CONFIDENCE_LEVEL=95 DF_METHOD=RESIDUAL COVB=MODEL
  /EMMEANS_OPTIONS SCALE=ORIGINAL PADJUST=LSD.
```

- The DATA\_STRUCTURE subcommand specifies that subjects are defined by *customer\_id*.
- The FIELDS and TARGET\_OPTIONS subcommands specify that the procedure creates a model for *service\_usage* using a multinomial distribution and a logit link to relate the target to a linear combination of the predictors.

- The FIXED subcommand specifies a model with *edcat*, *inccat*, *reside*, and *service\_type* as main effects.
- The RANDOM subcommand specifies an intercept-only random effect block with *customer\_id* as the subject field. This should account for correlation between answers to the service usage questions across service types (tv, phone, internet) within a given survey responder's answers.
- All other options are set to their default values.

## ***DATA\_STRUCTURE Subcommand***

The `DATA_STRUCTURE` subcommand specifies the subject structure for repeated measurements and how the errors of the repeated measurements are correlated. If the `DATA_STRUCTURE` subcommand is not specified, then the model assumes that all error terms are independent.

- Records with missing values for any field on the `DATA_STRUCTURE` subcommand are not used in the analysis.

### ***SUBJECTS Keyword***

The `SUBJECTS` keyword identifies subjects in the active dataset. Complete independence is assumed *across* subjects, but responses *within* subjects are assumed to be correlated.

- Specify a single categorical field or a list of categorical fields connected by asterisks (\*) or the keyword `BY`.
- The number of subjects equals the number of distinct combinations of values of the fields.
- Any field specified on the `SUBJECTS` keyword cannot be used as a `REPEATED_MEASURES` field or on the `FIELDS` subcommand.
- The `SUBJECTS` keyword is required if the `DATA_STRUCTURE` subcommand is used.

### ***REPEATED\_MEASURES Keyword***

The `REPEATED_MEASURES` keyword gives the repeated (or within-subject) effect. This effect defines the ordering of repeated measurements within subjects. If some measurements do not appear in the data for some subjects, then the existing measurements are ordered and the omitted measurements are treated as missing values. If `REPEATED_MEASURES` is not specified, then no repeated measures model is built.

- Specify a single field or a list of fields connected by asterisks (\*) or the keyword `BY`.
- Each distinct combination of the values of the fields defines a separate repeated measure.
- Any field specified on the `REPEATED_MEASURES` keyword cannot be used as a `SUBJECTS` keyword or on the `FIELDS` subcommand.
- Each distinct repeated measures value must occur only once within a subject. Generally speaking, you should not use a `WEIGHT` field if `REPEATED_MEASURES` is specified.

**GROUPING Keyword**

The `GROUPING` keyword allows you to define independent sets of covariance parameters. All subjects have the same covariance type; subjects within the same covariance grouping will have the same values for the parameters.

- Specify a single categorical field or a list of categorical fields connected by asterisks (\*) or the keyword `BY`.
- Each distinct combination of the values of the fields defines a separate covariance grouping.
- A field specified on the `GROUPING` keyword can also be used as a `SUBJECTS` or `REPEATED_MEASURES` field, the `TARGET` on the `FIELDS` subcommand, or on the `EFFECTS` keyword or the `FIXED` or `RANDOM` subcommand.
- The `GROUPING` keyword is optional.
- If the `GROUPING` keyword is used, there must also be a `REPEATED_MEASURES` specification.

**COVARIANCE\_TYPE Keyword**

The `COVARIANCE_TYPE` keyword gives the covariance structure of the residual covariance matrix.

**AR1.** *First-order autoregressive.* The constraint  $|\rho| \leq 1$  is imposed for stationarity.

$$\sigma^2 \begin{bmatrix} 1 & \rho & \rho^2 & \rho^3 \\ \rho & 1 & \rho & \rho^2 \\ \rho^2 & \rho & 1 & \rho \\ \rho^3 & \rho^2 & \rho & 1 \end{bmatrix}$$

**ARMA11.** *Autoregressive moving average (1,1).* The constraints  $|\phi| \leq 1$  and  $|\rho| \leq 1$  are imposed for stationarity.

$$\sigma^2 \begin{bmatrix} 1 & \phi\rho & \phi\rho^2 & \phi\rho^3 \\ \phi\rho & 1 & \phi\rho & \phi\rho^2 \\ \phi\rho^2 & \phi\rho & 1 & \phi\rho \\ \phi\rho^3 & \phi\rho^2 & \phi\rho & 1 \end{bmatrix}$$

**COMPOUND\_SYMMETRY.** This structure has constant variance and constant covariance.

$$\begin{bmatrix} \sigma^2 + \sigma_1 & \sigma_1 & \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma^2 + \sigma_1 & \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma_1 & \sigma^2 + \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma_1 & \sigma_1 & \sigma^2 + \sigma_1 \end{bmatrix}$$

**DIAGONAL.** This is a diagonal structure with heterogenous variance. This is the default.

$$\begin{bmatrix} \sigma_1^2 & 0 & 0 & 0 \\ 0 & \sigma_2^2 & 0 & 0 \\ 0 & 0 & \sigma_3^2 & 0 \\ 0 & 0 & 0 & \sigma_4^2 \end{bmatrix}$$

**IDENTITY.** This is a scaled identity matrix.

$$\sigma^2 \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix}$$

**TOEPLITZ.** ( $|\rho_k| \leq 1$ ).

$$\sigma^2 \begin{bmatrix} 1 & \rho_1 & \rho_2 & \rho_3 \\ \rho_1 & 1 & \rho_1 & \rho_2 \\ \rho_2 & \rho_1 & 1 & \rho_1 \\ \rho_3 & \rho_2 & \rho_1 & 1 \end{bmatrix}$$

**UNSTRUCTURED.** This is a completely general covariance matrix.

$$\begin{bmatrix} \sigma_1^2 & \sigma_{21} & \sigma_{31} & \sigma_{41} \\ \sigma_{21} & \sigma_2^2 & \sigma_{32} & \sigma_{42} \\ \sigma_{31} & \sigma_{32} & \sigma_3^2 & \sigma_{43} \\ \sigma_{41} & \sigma_{42} & \sigma_{43} & \sigma_4^2 \end{bmatrix}$$

**VARIANCE\_COMPONENTS.** For repeated measures, the variance components structure is equivalent to the diagonal (**DIAGONAL**) structure.

## ***FIELDS Subcommand***

The **FIELDS** subcommand specifies the target, optional offset, and optional analysis (regression) weight.

- This subcommand, and the **TARGET** keyword, are required.
- A field can only be specified once on the **FIELDS** subcommand.
- A categorical **TARGET** can be specified as a **GROUPING** field on the **DATA\_STRUCTURE** or **RANDOM** subcommand, but not in any other capacity. Any other field specified on the **FIELDS** subcommand cannot be used on any other subcommand.

**TARGET.** The target field is what the model tries to predict. Specify a single field.

**TRIALS = NONE\*\* | VALUE(number) | FIELD(field).** If the model response can be expressed as the number of events occurring within a number of trials, then the **TARGET** keyword specifies the number of events and **TRIALS** specifies the number of trials. Use **VALUE** with the number of trials in parentheses if the number of trials is fixed for all subjects and **FIELD** with a field name in parentheses if the number of trials varies across subjects and there is a field containing the number of trials.

The procedure automatically computes the ratio of the events field over the trials field or number. Technically, the procedure treats the events field as the target in the sense that predicted values and residuals are based on the events field rather than the events/trials ratio.

**ANALYSIS\_WEIGHT.** The model uses analysis (regression) weights to account for differences in variability across levels of the target field. Specify a single continuous field.

**OFFSET = NONE\*\* | VALUE(value) | FIELD(field).** The offset is a constant added to the intercept. Specify a continuous field or number.

- Records with missing values on the `OFFSET` field are not used in the analysis.
- Specifying a number when `/FIXED USE_INTERCEPT = YES` is equivalent to adding a constant to the intercept.
- Specifying a number when `/FIXED USE_INTERCEPT = NO` is equivalent to fixing the intercept at the specified number.

## **TARGET\_OPTIONS Subcommand**

The `TARGET_OPTIONS` subcommand specifies the reference category, distribution, and link function for the target.

- This subcommand is optional. If the `TARGET_OPTIONS` subcommand is not specified, then `GENLINMIXED` assigns a default distribution and link function based on the measurement level of the target. A continuous target has `NORMAL` distribution and `IDENTITY` link. A categorical target has `MULTINOMIAL` distribution and `LOGIT` link.

**REFERENCE.** For categorical targets, this is the target value to use as the reference category for parameter estimation. No model parameters are assigned to the reference category. Specify a single value. Put the value inside a pair of quotes if it is formatted (such as date or time) or if the target is of string storage type. The `REFERENCE` value must be found in the dataset. If `REFERENCE` is not specified, the last category in ascending order is used as the reference.

### ***DISTRIBUTION Keyword***

The `DISTRIBUTION` keyword specifies the probability distribution of the target.

- The default probability distribution depends on the measurement level of the target. If the target is continuous, the default distribution is `NORMAL`. If the target is categorical, then the default is `MULTINOMIAL`. If an events/trials specification is used, then the default distribution is `BINOMIAL`.
- Caution must be exercised when the `LINK` but not the `DISTRIBUTION` keyword is used. The discussion of the `LINK` keyword below gives details about proper and improper combinations of `DISTRIBUTION` and `LINK` settings.

**BINOMIAL.** *Binomial probability distribution.* If the target is specified as a single field, then it can be numeric or string and it can have only two distinct valid data values.

If the events and trials options are specified, then the procedure automatically computes the ratio of the events field over the trials field or number. The events field—and the trials field if specified—must be numeric. Data values for the events field must be integers greater than or equal to zero. Data values for the trials field must be integers greater than zero. For each record, the trials value must be greater than or equal to the events value. If an events value is noninteger, less than zero, or missing, then the corresponding record is not used in the analysis. If a trials value is noninteger, less than or equal to zero, less than the events value, or missing, then the corresponding record is not used in the analysis.

If the trials option specifies a number, then it must be a positive integer, and it must be greater than or equal to the events value for each record. Records with invalid values are not used in the analysis.

This is the default probability distribution if the target is specified using events/trials format.

**GAMMA.** *Gamma probability distribution.* The target must be numeric, with data values greater than zero. If a data value is less than or equal to zero, or missing, then the corresponding record is not used in the analysis.

**INVERSE\_GAUSSIAN.** *Inverse Gaussian probability distribution.* The target must be numeric, with data values greater than zero. If a data value is less than or equal to zero, or missing, then the corresponding record is not used in the analysis.

**MULTINOMIAL.** *Multinomial probability distribution.* The target can be numeric or string, and it must have at least two distinct, valid data values. This is the default probability distribution if the target is categorical.

**NEGATIVE\_BINOMIAL.** *Negative binomial probability distribution.* The target must be numeric, with data values that are integers greater than or equal to zero. If a data value is noninteger, less than zero, or missing, then the corresponding record is not used in the analysis.

**NORMAL.** *Normal probability distribution.* The target must be numeric. This is the default probability distribution if the target is continuous.

**POISSON.** *Poisson probability distribution.* The target must be numeric, with data values that are integers greater than or equal to zero. If a data value is noninteger, less than zero, or missing, then the corresponding record is not used in the analysis.

### **LINK Keyword**

The LINK keyword specifies the link function. The following link functions are available.

**IDENTITY.** *Identity link function.*  $f(x)=x$

**CLOGLOG.** *Complementary log-log link function.*  $f(x)=\ln(-\ln(1-x))$

**LOG.** *Log link function.*  $f(x)=\ln(x)$

**LOGC.** *Log complement link function.*  $f(x)=\ln(1-x)$

**LOGIT.** *Logit link function.*  $f(x)=\ln(x / (1-x))$

**NLOGLOG.** *Negative log-log link function.*  $f(x)=-\ln(-\ln(x))$

**POWER.** *Power link function.*  $f(x)=x^\alpha$ , if  $\alpha \neq 0$ .  $f(x)=\ln(x)$ , if  $\alpha=0$ .  $\alpha$  is a required number specified on the LINKPARAMETER keyword and must be a real number. If  $|\alpha| < 2.2e-16$ ,  $\alpha$  is treated as 0.

**PROBIT.** *Probit link function.*  $f(x)=\Phi^{-1}(x)$ , where  $\Phi^{-1}$  is the inverse standard normal cumulative distribution function.

- The default setting for LINK depends on the DISTRIBUTION setting, whether DISTRIBUTION is set explicitly or is the default, as shown in the following table.

DISTRIBUTION Setting	Default LINK Setting
NORMAL	IDENTITY
BINOMIAL	LOGIT
GAMMA	POWER(-1)
INVERSE_GAUSSIAN	POWER(-2)
MULTINOMIAL	LOGIT
NEGATIVE_BINOMIAL	LOG
POISSON	LOG

- The procedure will fit a model if a permissible combination of LINK and DISTRIBUTION specifications is given. The table below indicates the permissible LINK and DISTRIBUTION combinations. Specifying an improper combination will yield an error message.
- Note that the default setting for DISTRIBUTION is irrespective of the LINK specification, and that not all LINK specifications are valid for the default DISTRIBUTION. Thus, if LINK = CLOGLOG is specified for a continuous target but DISTRIBUTION is not, then the default DISTRIBUTION = NORMAL yields an improper combination of DISTRIBUTION and LINK settings.

Table 95-1

Valid combinations of distribution and link function

Link	Distribution							
	NORMAL	BINOMIAL	GAMMA	INVERSE_GAUSSIAN	MULTINOMIAL	NEGATIVE_BINOMIAL	POISSON	BINOMIAL
IDENTITY	X	X	X	X		X		X
CLOGLOG		X						
LOG	X	X	X	X		X		X
LOGC		X						
LOGIT		X			X			
NLOGLOG		X						
PROBIT		X						
POWER	X	X	X	X		X		X

### **LINKPARAMETER Keyword**

The LINKPARAMETER keyword specifies the value of  $\alpha$  for the POWER link.

- Specify a number. There is no default.

## **FIXED Subcommand**

The FIXED subcommand is used to specify fixed model effects, including the intercept.

- This subcommand is optional. If not specified, then the procedure builds an intercept-only model.

**EFFECTS.** The effect list includes all effects to be included in the fixed model except for the intercept, which is specified using the USE\_INTERCEPT keyword.



- To include a term for the main effect of a factor (categorical predictor) or covariate (continuous predictor), enter its field name. Whether a field is treated as a factor or covariate depends upon its measurement level. Fields with categorical (flag, nominal, or ordinal) measurement level are treated as factors while fields with continuous (scale) measurement level are treated as covariates.
- To include a term for an interaction between factors, use the keyword `BY` or an asterisk (\*) to join the factors involved in the interaction. For example, `A*B` means a two-way interaction effect of `A` and `B`, where `A` and `B` are factors. `A*A` is not allowed because factors in an interaction effect must be distinct.
- To include a term for nesting one effect within another, use a pair of parentheses. For example, `A(B)` means that `A` is nested within `B`.
- Multiple nesting is allowed. For example, `A(B(C))` means that `B` is nested within `C`, and `A` is nested within `B(C)`. When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.
- Interactions between nested effects are not valid. For example, neither `A(C)*B(C)` nor `A(C)*B(D)` is valid.
- Covariates can be connected, but not nested, through the \* operator to form another covariate effect. Interactions among covariates such as `X1*X1` and `X1*X2` are valid, but `X1(X2)` is not.
- Factor and covariate effects can be connected only by the \* operator. Suppose `A` and `B` are factors, and `X1` and `X2` are covariates. Examples of valid factor-by-covariate interaction effects are `A*X1`, `A*B*X1`, `X1*A(B)`, `A*X1*X1`, and `B*X1*X2`.

`USE_INTERCEPT = TRUE** | FALSE`. This keyword controls whether an intercept term is included in the model. Specifying `USE_INTERCEPT=FALSE` when there are no effects in the model fits a “null” model. The default is `TRUE`.

## ***RANDOM Subcommand***

The `RANDOM` subcommand specifies the random effects in the mixed model.

- Depending on the covariance type specified, random effects specified in one `RANDOM` subcommand may be correlated.
- One block of the covariance `G` matrix will be constructed for each `RANDOM` subcommand. The dimension of the random effect covariance `G` matrix is equal to the sum of the levels of all random effects across all `RANDOM` subcommands.
- When the variance components (`VC`) structure is specified, a scaled identity (`ID`) structure will be assigned to each of the effects specified. This is the default covariance type for the `RANDOM` subcommand.
- Use a separate `RANDOM` subcommand when a different covariance structure is assumed for a list of random effects. If the same effect is listed on more than one `RANDOM` subcommand, it must be associated with a different `SUBJECT` combination.
- No random effects are included in the mixed model unless a `RANDOM` subcommand is specified correctly.

**EFFECTS.** The effect list includes all effects to be included in the random effects model block except for the intercept, which is specified using the `USE_INTERCEPT` keyword.

- To include a term for the main effect of a factor (categorical predictor) or covariate (continuous predictor), enter its field name. Whether a field is treated as a factor or covariate depends upon its measurement level.
- To include a term for an interaction between factors, use the keyword `BY` or an asterisk (\*) to join the factors involved in the interaction. For example, `A*B` means a two-way interaction effect of `A` and `B`, where `A` and `B` are factors. `A*A` is not allowed because factors in an interaction effect must be distinct.
- To include a term for nesting one effect within another, use a pair of parentheses. For example, `A(B)` means that `A` is nested within `B`.
- Multiple nesting is allowed. For example, `A(B(C))` means that `B` is nested within `C`, and `A` is nested within `B(C)`. When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.
- Interactions between nested effects are not valid. For example, neither `A(C)*B(C)` nor `A(C)*B(D)` is valid.
- Covariates can be connected, but not nested, through the \* operator to form another covariate effect. Interactions among covariates such as `X1*X1` and `X1*X2` are valid, but `X1(X2)` is not.
- Factor and covariate effects can be connected only by the \* operator. Suppose `A` and `B` are factors, and `X1` and `X2` are covariates. Examples of valid factor-by-covariate interaction effects are `A*X1`, `A*B*X1`, `X1*A(B)`, `A*X1*X1`, and `B*X1*X2`.

**USE\_INTERCEPT.** This keyword controls whether an intercept term is included in the model. It is invalid to specify `USE_INTERCEPT=FALSE` if there are no effects specified on the `EFFECTS` keyword. The default is `FALSE`.

### Examples

```
GENLINMIXED
  /FIXED EFFECTS=a b
  /RANDOM SUBJECTS=id.
```

```
GENLINMIXED
  /FIXED EFFECTS=a b
  /RANDOM USE_INTERCEPT=TRUE SUBJECTS=id.
```

```
GENLINMIXED
  /FIXED EFFECTS=a b
  /RANDOM EFFECTS=c d SUBJECTS=id.
```

- The first command fails because the `RANDOM` subcommand has no effects or intercept. The next two commands succeed.

### SUBJECTS Keyword

The `SUBJECTS` keyword identifies subjects in the active dataset. Complete independence is assumed *across* subjects, but responses *within* subjects are assumed to be correlated.

- Specify a single field or a list of fields connected by asterisks (\*) or the keyword BY.
- The number of subjects equals the number of distinct combinations of values of the fields.
- The fields in the RANDOMSUBJECTS list must be a subset of the fields on the DATA\_STRUCTURE SUBJECTS list.
- The SUBJECTS keyword is optional.

### **GROUPING Keyword**

The GROUPING keyword allows you to define independent sets of covariance parameters. All subjects have the same covariance type; subjects within the same covariance grouping will have the same values for the parameters.

- Specify a single field or a list of fields connected by asterisks (\*) or the keyword BY.
- Each distinct combination of the values of the fields defines a separate covariance grouping.
- A field specified on the GROUPING keyword can also be used as a SUBJECTS or REPEATED\_MEASURES field, the TARGET on the FIELDS subcommand, or on the EFFECTS keyword or the FIXED or RANDOM subcommand.
- The GROUPING keyword is optional.

### **COVARIANCE\_TYPE Keyword**

The COVARIANCE\_TYPE keyword gives the covariance structure of the random effect covariance matrix.

- Random effects are considered independent of each other, and a separate covariance matrix is computed for each effect.

**AR1.** *First-order autoregressive.* The constraint  $|\rho| \leq 1$  is imposed for stationarity.

$$\sigma^2 \begin{bmatrix} 1 & \rho & \rho^2 & \rho^3 \\ \rho & 1 & \rho & \rho^2 \\ \rho^2 & \rho & 1 & \rho \\ \rho^3 & \rho^2 & \rho & 1 \end{bmatrix}$$

**ARMA11.** *Autoregressive moving average (1,1).* The constraints  $|\varphi| \leq 1$  and  $|\rho| \leq 1$  are imposed for stationarity.

$$\sigma^2 \begin{bmatrix} 1 & \phi\rho & \phi\rho^2 & \phi\rho^3 \\ \phi\rho & 1 & \phi\rho & \phi\rho^2 \\ \phi\rho^2 & \phi\rho & 1 & \phi\rho \\ \phi\rho^3 & \phi\rho^2 & \phi\rho & 1 \end{bmatrix}$$

**COMPOUND\_SYMMETRY.** This structure has constant variance and constant covariance.

$$\begin{bmatrix} \sigma^2 + \sigma_1 & \sigma_1 & \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma^2 + \sigma_1 & \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma_1 & \sigma^2 + \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma_1 & \sigma_1 & \sigma^2 + \sigma_1 \end{bmatrix}$$

**DIAGONAL.** This is a diagonal structure with heterogenous variance.

$$\begin{bmatrix} \sigma_1^2 & 0 & 0 & 0 \\ 0 & \sigma_2^2 & 0 & 0 \\ 0 & 0 & \sigma_3^2 & 0 \\ 0 & 0 & 0 & \sigma_4^2 \end{bmatrix}$$

**IDENTITY.** This is a scaled identity matrix.

$$\sigma^2 \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix}$$

**TOEPLITZ.** ( $|\rho_k| \leq 1$ ).

$$\sigma^2 \begin{bmatrix} 1 & \rho_1 & \rho_2 & \rho_3 \\ \rho_1 & 1 & \rho_1 & \rho_2 \\ \rho_2 & \rho_1 & 1 & \rho_1 \\ \rho_3 & \rho_2 & \rho_1 & 1 \end{bmatrix}$$

**UNSTRUCTURED.** This is a completely general covariance matrix.

$$\begin{bmatrix} \sigma_1^2 & \sigma_{21} & \sigma_{31} & \sigma_{41} \\ \sigma_{21} & \sigma_2^2 & \sigma_{32} & \sigma_{42} \\ \sigma_{31} & \sigma_{32} & \sigma_3^2 & \sigma_{43} \\ \sigma_{41} & \sigma_{42} & \sigma_{43} & \sigma_4^2 \end{bmatrix}$$

**VARIANCE\_COMPONENTS.** *Variance components.* This is the default covariance structure for random effects. The variance components structure for random effects is a scaled identity (ID) structure assigned to each of the effects specified on the subcommand.

## ***BUILD\_OPTIONS Subcommand***

The `BUILD_OPTIONS` subcommand specifies the criteria used to build the model.

**TARGET\_CATEGORY\_ORDER.** This determines the order of the target's categories for purposes of determining the "last" category, which is used as the reference category if the `REFERENCE` keyword is not used. This keyword is ignored if the target is not categorical.

- **ASCENDING.** *Target categories are sorted in ascending order, from the lowest value to the highest value.* The highest value is treated as the "last" category. This is the default.
- **DESCENDING.** *Target categories are sorted in descending order, from the highest value to the lowest value.* The lowest value is treated as the "last" category.
- **DATA.** *Target values are not sorted.* The first value encountered in the data defines the first category, the last value encountered defines the last category. This option is technically not valid if splits are defined on the `SPLIT FILE` command, but the procedure will run while it does not honor splits.

**INPUTS\_CATEGORY\_ORDER.** This determines the order of the categories for factors (categorical inputs) for purposes of determining the "last" category, which is used as the reference category. This keyword is ignored if there are no factors.

- **ASCENDING.** *Factor values are sorted in ascending order, from the lowest value to the highest value.* The highest value is treated as the “last” category. This is the default.
- **DESCENDING.** *Factor values are sorted in descending order, from the highest value to the lowest value.* The lowest value is treated as the “last” category.
- **DATA.** *Factor values are not sorted.* The first value encountered in the data defines the first category, the last value encountered defines the last category. This option is technically not valid if splits are defined on the `SPLIT FILE` command, but the procedure will run while it does not honor splits.

**MAX\_ITERATIONS = number.** The model building algorithm stops after a certain number of iterations. By default, this is 100. Alternatively, specify a non-negative integer maximum number of iterations.

**CONFIDENCE\_LEVEL.** This is the level of confidence used to compute interval estimates of the model coefficients. Specify a value greater than 0 and less than 100. The default is 95.

**DF\_METHOD.** This specifies how degrees of freedom are computed for significance tests.

- **RESIDUAL.** The residual method has a fixed degrees of freedom for all tests. It is useful if your sample size is sufficiently large, or the data are balanced, or the model uses a simpler covariance type; for example, scaled identity or diagonal. This is the default.
- **SATTERTHWAITE.** The Satterthwaite method has a field degrees of freedom across tests. It is useful if your sample size is small, or the data are unbalanced, or the model uses a complicated covariance type; for example, unstructured.

**COVB.** This is the method for computing the parameter estimates covariance matrix.

- **MODEL.** The model-based estimate assumes that the model assumptions are correct. This is the default.
- **ROBUST.** The robust estimate is useful if you are concerned that the model assumptions are violated.

## ***EMMEANS Subcommand***

The `EMMEANS` subcommand displays estimated marginal means of the target for all level combinations of a set of factors. Note that these are predicted, not observed, means. Estimated marginal means can be computed based on the original scale of the target or the based on the link function transformation.

- Multiple `EMMEANS` subcommands are allowed. Each is treated independently.
- The `EMMEANS` subcommand can be specified with no additional keywords. The output for an empty `EMMEANS` subcommand is the overall estimated marginal mean of the response, collapsing over any factors.
- Estimated marginal means are not available if the multinomial distribution is used. If `DISTRIBUTION = MULTINOMIAL` on the `MODEL` subcommand and the `EMMEANS` subcommand is specified, then `EMMEANS` is ignored and a warning is issued.

**TABLES Keyword**

The TABLES keyword specifies the cells for which estimated marginal means are displayed.

- Valid options are all-factor effects appearing on the EFFECTS keyword of the FIXED subcommand, including main effects factors such as  $a$   $b$ , all-factor interactions such as  $a*b$ , and all-factor nested effects such as  $a(b)$ , if  $a$  and  $b$  are factors.
- If the TABLES keyword is specified, then the procedure collapses over any other factors specified on the command but not on the TABLES keyword before computing the estimated marginal means for the target.
- If the TABLES keyword is not specified, then the overall estimated marginal mean of the target, collapsing over any factors, is computed.

**Examples**

```
GENLINMIXED
  /FIXED EFFECTS=a b
  /EMMEANS TABLES=a*b.
```

```
GENLINMIXED
  /FIXED EFFECTS=a*b
  /EMMEANS TABLES=a
  /EMMEANS TABLES=b.
```

```
GENLINMIXED
  /FIXED EFFECTS=a b a*b
  /EMMEANS TABLES=a
  /EMMEANS TABLES=b
  /EMMEANS TABLES=a*b.
```

- The first two commands fail because the effects listed on the TABLES keywords are not listed on the EFFECTS keyword. The third command succeeds.

**COMPARE Keyword**

The COMPARE keyword specifies a factor, the levels of which are compared using the contrast type specified on the CONTRAST keyword.

- Valid options are factors appearing on the TABLES keyword.
- The COMPARE keyword is valid only if the TABLES and CONTRAST keyword is also specified.
- By default, the procedure sorts levels of the factors in ascending order and defines the highest level as the last level. (If the factor is a string variable, then the value of the highest level is locale-dependent.) However, the sort order can be modified using the INPUTS\_SORT\_ORDER keyword on the BUILD\_OPTIONS subcommand.

**CONTRAST Keyword**

The CONTRAST keyword specifies the type of contrast to use for the levels of the factor on the COMPARE keyword. The CONTRAST keyword creates an **L** matrix (that is, a coefficient matrix) such that the columns corresponding to the factor match the contrast given. The other columns are adjusted so that the **L** matrix is estimable.

- The `CONTRAST` keyword is valid only if the `COMPARE` keyword is also specified.
- If the `COMPARE` keyword is specified without `CONTRAST`, then pairwise comparisons are performed for the factor on `COMPARE`.
- `SIMPLE` contrasts are defined with respect to the last level as determined by the `INPUTS_CATEGORY_ORDER` specification on the `BUILD_OPTIONS` subcommand.

The following contrast types are available.

**NONE.** *No comparisons.* This is the default.

**PAIRWISE.** *Pairwise comparisons are computed for all levels of the specified or factor.* Pairwise contrasts are not orthogonal.

**DEVIATION.** *Each level of the factor is compared to the grand mean.* Deviation contrasts are not orthogonal.

**SIMPLE.** *Each level of the factor except the last is compared to the last level.* Simple contrasts are not orthogonal.

## ***EMMEANS\_OPTIONS Subcommand***

The `EMMEANS_OPTIONS` subcommand contains specifications that are used across all `EMMEANS` subcommands.

### ***CONTROL Keyword***

The `CONTROL` keyword specifies the covariate values to use when computing the estimated marginal means.

- Specify one or more covariates appearing on the `FIXED` subcommand `EFFECTS` list, each of which must be followed by a numeric value.
- If a covariate is not specified on the `CONTROL` option, then its overall mean will be used in estimated marginal means calculations.
- Any covariate can occur only once on the `CONTROL` keyword.

### ***SCALE Keyword***

The `SCALE` keyword specifies whether to compute estimated marginal means based on the original scale of the target or based on the link function transformation.

**ORIGINAL.** *Estimated marginal means are based on the original scale of the target.* Estimated marginal means are computed for the target. This is the default. Note that when the target is specified using the `events/trials` option, `ORIGINAL` gives the estimated marginal means for the `events/trials` proportion rather than for the number of events.

**TRANSFORMED.** *Estimated marginal means are based on the link function transformation.* Estimated marginal means are computed for the linear predictor.

**PADJUST Keyword**

The PADJUST keyword indicates the method of adjusting the significance level.

**LSD.** *Least significant difference.* This method does not control the overall probability of rejecting the hypotheses that some linear contrasts are different from the null hypothesis value(s). This is the default.

**SEQBONFERRONI.** *Sequential Bonferroni.* This is a sequentially step-down rejective Bonferroni procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.

**SEQSIDAK.** *Sequential Sidak.* This is a sequentially step-down rejective Sidak procedure that is much less conservative in terms of rejecting individual hypotheses but maintains the same overall significance level.

**OUTFILE Subcommand**

The OUTFILE subcommand saves a file containing the scoring model.

**MODEL = 'file'.** Writes the fixed effects portion of the model to an external .zip file. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#) Specify a unique, valid filename. If the file specification refers to an existing file, then the file is overwritten.

**SAVE Subcommand**

The SAVE subcommand writes optional temporary fields to the active dataset.

**PREDICTED\_VALUES(varname).** *Predicted value of the target.* The default field name is *PredictedValue*.

**PREDICTED\_PROBABILITY(rootname).** *Predicted probabilities of the target.* If the target is categorical, this keyword saves the predicted probabilities of the first n categories. The value of n is specified by the MAX\_CATEGORIES keyword.

- Specify a unique, valid variable name. The default root name is *PredictedProbability*.
- If the dependent variable is continuous, this keyword is ignored with a warning.

**MAX\_CATEGORIES.** *The maximum number of categories for which predicted probabilities and confidence intervals (for a categorical target) should be saved.*

- Specify a positive integer. The default is 25.
- This keyword is ignored if predicted probabilities or confidence intervals (for a categorical target) are not saved.



**CONFIDENCE(varname).** *Confidence in predicted value for categorical target.* The computed confidence can be based on the probability of the predicted value (the highest predicted probability) or the difference between the highest predicted probability and the second highest predicted probability. The default field name is *Confidence*.

**CONFIDENCE\_METHOD.** *Method for computing confidence.* The computed confidence can be based on the probability of the predicted value (the highest predicted probability) or the difference between the highest predicted probability and the second highest predicted probability.

**CONFIDENCE\_INTERVALS(rootname).** *Upper and lower confidence intervals.* For all distributions except the multinomial, this creates two variables and the default root name is *CI*, with *\_Lower* and *\_Upper* as the suffixes. Specify a root name in parentheses to override the default.

For the multinomial distribution, one field is created for each dependent variable category except the last (see the `TARGET_SORT_ORDER` keyword). `CONFIDENCE_INTERVALS` saves the lower and upper bounds of the cumulative predicted probability for the first 25 categories, up to but not including the last, by default. The default root name is *CI*, and the default field names are *CI\_Lower\_1*, *CI\_Upper\_1*, *CI\_Lower\_2*, *CI\_Upper\_2*, and so on, corresponding to the order of the target categories. Specify a root name in parentheses to override the default. The value of *n* is specified by the `MAX_CATEGORIES` keyword.

**PEARSON\_RESIDUALS(varname).** *Pearson residuals.* The default field name is *PearsonResidual*.

# GENLOG

GENLOG is available in the Advanced Statistics option.

```
GENLOG varlist[BY] varlist [WITH covariate varlist]
  [/CSTRUCTURE=varname]
  [/GRESID=varlist]
  [/GLOR=varlist]
  [/MODEL={POISSON** }
          {MULTINOMIAL}]
  [/CRITERIA=[CONVERGE({0.001**})] [ITERATE({20**})] [DELTA({0.5**})]
             {n }           {n }           {n }
             [CIN({95**})] [EPS({1E-8**})]
             {n }           {n }
             [DEFAULT]]
  [/PRINT={FREQ**] [RESID**] [ADJRESID**] [DEV**]
           [ZRESID] [ITERATION] [COV] [DESIGN] [ESTIM] [CORR]
           [ALL] [NONE]
           [DEFAULT]]
  [/PLOT={DEFAULT**           }
         {RESID([ADJRESID] [DEV]) }
         {NORMPROB([ADJRESID] [DEV]) }
         {NONE }
         ]
  [/SAVE=tempvar (newvar) [tempvar (newvar)...]]
  [/MISSING={(EXCLUDE**)}}
           {INCLUDE }
           ]
  [/DESIGN=effect[(n)] effect[(n)]... effect {BY} effect...
           {* }
           ]
```

**\*\***Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
GENLOG DPREF RACE CAMP.
```

## Overview

GENLOG is a general procedure for model fitting, hypothesis testing, and parameter estimation for any model that has categorical variables as its major components. As such, GENLOG subsumes a variety of related techniques, including general models of multiway contingency tables, logit models, logistic regression on categorical variables, and quasi-independence models.

GENLOG, following the regression approach, uses dummy coding to construct a design matrix for estimation and produces maximum likelihood estimates of parameters by means of the Newton-Raphson algorithm. Since the regression approach uses the original parameter spaces, the parameter estimates correspond to the original levels of the categories and are therefore easier to interpret.

HIOGLOGLINEAR, which uses an iterative proportional-fitting algorithm, is more efficient for hierarchical models and useful in model building, but it cannot produce parameter estimates for unsaturated models, does not permit specification of contrasts for parameters, and does not display a correlation matrix of the parameter estimates.

The General Loglinear Analysis and Logit Loglinear Analysis dialog boxes are both associated with the GENLOG command. In previous releases, these dialog boxes were associated with the LOGLINEAR command. The LOGLINEAR command is now available only as a syntax command. The differences are described in the discussion of the LOGLINEAR command.

### **Options**

**Cell Weights.** You can specify cell weights (such as structural zero indicators) for the model with the CSTRUCTURE subcommand.

**Linear Combinations.** You can compute linear combinations of observed cell frequencies, expected cell frequencies, and adjusted residuals using the GRESID subcommand.

**Generalized Log-Odds Ratios.** You can specify contrast variables on the GLOR subcommand and test whether the generalized log-odds ratio equals 0.

**Model Assumption.** You can specify POISSON or MULTINOMIAL on the MODEL subcommand to request the Poisson loglinear model or the product multinomial loglinear model.

**Tuning the Algorithm.** You can control the values of algorithm-tuning parameters with the CRITERIA subcommand.

**Output Display.** You can control the output display with the PRINT subcommand.

**Optional Plots.** You can request plots of adjusted or deviance residuals against observed and expected counts, or normal plots and detrended normal plots of adjusted or deviance residuals using the PLOT subcommand.

### **Basic Specification**

The basic specification is one or more factor variables that define the tabulation. By default, GENLOG assumes a Poisson distribution and estimates the saturated model. Default output includes the factors or effects, their levels, and any labels; observed and expected frequencies and percentages for each factor and code; and residuals, adjusted residuals, and deviance residuals.

### **Limitations**

- Maximum 10 factor variables (dependent *and* independent).
- Maximum 200 covariates.

**Subcommand Order**

- The variable specification must come first.
- Subcommands can be specified in any order.
- When multiple subcommands are specified, only the last specification takes effect.

**Examples**

```
GENLOG DPREF RACE CAMP.
```

- *DPREF*, *RACE*, and *CAMP* are categorical variables.
- This is a general loglinear model because no `BY` keyword appears.
- The design defaults to a saturated model that includes all main effects and two-way and three-way interaction effects.

**Example: Specifying a Custom Model**

```
GENLOG GSLEVEL EDUC SEX
  /DESIGN=GSLEVEL EDUC SEX.
```

- *GSLEVEL*, *EDUC*, and *SEX* are categorical variables.
- `DESIGN` specifies a model with main effects only.

**Poisson Regression**

```
WEIGHT
  BY accid .
GENLOG
  agecat gender /CSTRUCTURE = pop
  /MODEL = POISSON
  /PRINT = FREQ RESID DEV ADJRESID ESTIM CORR COV
  /CRITERIA = CIN(95) ITERATE(20) CONVERGE(.001) DELTA(.5)
  /DESIGN .
```

- The procedure examines the relationship between *agecat* and *gender*. Cases are weighted by *accid* to give the raw cell counts, and the `CSTRUCTURE` subcommand specifies that *pop* should be used to weight the cells to account for cell-to-cell variation in the population at risk.
- The empty `DESIGN` subcommand means that a saturated model will be fit. No plots are produced for a saturated model.
- The `PRINT` subcommand requests, in addition to the default output, the parameter estimates and the correlation and covariance matrix of the estimates.

**Paired Data : Symmetry, Quasi-Symmetry, and Quasi-Independence Models**

```
WEIGHT
  BY count .
GENLOG
  pref1 pref2 WITH s12 s13 s14 s23 s24 s34 /CSTRUCTURE = offdiag
  /MODEL = POISSON
  /PRINT = FREQ RESID DEV ADJRESID
  /PLOT = NONE
  /CRITERIA = CIN(95) ITERATE(20) CONVERGE(.001) DELTA(.5)
  /DESIGN s12 s13 s14 s23 s24 s34 .
```

- The procedure examines the relationship between *pref1* and *pref2*, using *s12*, *s13*, *s14*, *s23*, *s24*, and *s34* as cell covariates. Cases are weighted by *count* to give the raw cell counts, and the `CSTRUCTURE` subcommand specifies that *offdiag* should be used to indicate the off-diagonal cells of interest that correspond to cases that switched their preferred candidate.
- The `DESIGN` subcommand specifies that a symmetry model will be fit; the symmetry model for paired data supposes that for each supporter a candidate gained from another contender, they lost a supporter to that same contender.
- The `PLOT` subcommand turns off the default plots.
- All other options are set to their default values.

```
GENLOG
pref1 pref2 WITH s12 s13 s14 s23 s24 s34 /CSTRUCTURE = offdiag
/MODEL = POISSON
/PRINT = FREQ RESID DEV ADJRESID
/PLOT = NONE
/CRITERIA = CIN(95) ITERATE(20) CONVERGE(.001) DELTA(.5)
/DESIGN pref1 pref2 s12 s13 s14 s23 s24 s34 .
```

- The `DESIGN` subcommand now specifies a quasi-symmetry model will be fit; the likelihood ratio statistic from symmetry and quasi-symmetry models can be compared to determine whether there is a statistically significant change in voter preference after the debate.

```
GENLOG
pref1 pref2 /CSTRUCTURE = offdiag
/MODEL = POISSON
/PRINT = FREQ RESID DEV ADJRESID
/PLOT = NONE
/CRITERIA = CIN(95) ITERATE(20) CONVERGE(.001) DELTA(.5)
/DESIGN pref1 pref2 .
```

- The procedure now removes the cell covariates to specify the quasi-independence model, which is a main-effect model of *pref1* and *pref2* with *offdiag* as the cell structure variable. When marginal homogeneity does not hold, the quasi-independence model tests the null hypothesis that respondents who switched from a given candidate did not tend to move to particular opponents.

### **Logit Loglinear Analysis**

```
GENLOG
bfast BY active gender agecat
/MODEL = MULTINOMIAL
/PRINT = FREQ RESID DEV ADJRESID ESTIM CORR COV
/PLOT = NONE
/CRITERIA = CIN(95) ITERATE(20) CONVERGE(.001) DELTA(.5)
/DESIGN bfast bfast*active bfast*agecat .
```

- The keyword `BY` on the variable list specifies a logit model in which *bfast* is the dependent variable and *active*, *gender*, and *agecat* are the independent variables.
- A logit model is multinomial.
- The `DESIGN` subcommand specifies a model that tests for the absence of the joint effect of *active* and *agecat* on *bfast*.
- The `PLOT` subcommand turns off the default plots.

- The `PRINT` subcommand requests, in addition to the default output, the parameter estimates and the correlation and covariance matrix of the estimates.
- All other options are set to their default values.

## Variable List

The variable list specifies the variables to be included in the model. `GENLOG` analyzes two classes of variables—categorical and continuous. Categorical variables are used to define the cells of the table. Continuous variables are used as cell covariates.

- The list of categorical variables must be specified first. Categorical variables must be numeric.
- Continuous variables can be specified only after the `WITH` keyword following the list of categorical variables.
- To specify a logit model, use the keyword `BY` (see [Logit Model on p. 824](#)). A variable list without the keyword `BY` generates a general loglinear model.
- A variable can be specified only once in the variable list—as a dependent variable immediately following `GENLOG`, as an independent variable following the keyword `BY`, or as a covariate following the keyword `WITH`.
- No range needs to be specified for categorical variables.

## Logit Model

The logit model examines the relationships between dependent and independent factor variables.

- To separate the independent variables from the dependent variables in a logit model, use the keyword `BY`. The categorical variables preceding `BY` are the dependent variables; the categorical variables following `BY` are the independent variables.
- Up to 10 variables can be specified, including both dependent and independent variables.
- For the logit model, you must specify `MULTINOMIAL` on the `MODEL` subcommand.
- `GENLOG` displays an analysis of dispersion and two measures of association—entropy and concentration. These measures are discussed elsewhere (Haberman, 1982) and can be used to quantify the magnitude of association among the variables. Both are proportional-reduction-in-error measures. The entropy statistic is analogous to Theil's entropy measure, while the concentration statistic is analogous to Goodman and Kruskal's tau-*b*. Both statistics measure the strength of association between the dependent variable and the independent variable set.

### Example

```
GENLOG  GSLEVEL BY EDUC SEX
      /MODEL=MULTINOMIAL
      /DESIGN=GSLEVEL, GSLEVEL BY EDUC, GSLEVEL BY SEX.
```

- The keyword `BY` on the variable list specifies a logit model in which `GSLEVEL` is the dependent variable and `EDUC` and `SEX` are the independent variables.
- A logit model is multinomial.

- `DESIGN` specifies a model that can test for the absence of the joint effect of *SEX* and *EDUC* on *GSLEVEL*.

### **Cell Covariates**

- Continuous variables can be used as covariates. When used, the covariates must be specified after the `WITH` keyword following the list of categorical variables.
- A variable cannot be named as both a categorical variable and a cell covariate.
- To enter cell covariates into a model, the covariates must be specified on the `DESIGN` subcommand.
- Cell covariates are not applied on a case-by-case basis. The weighted covariate mean for a cell is applied to that cell.

#### **Example**

```
GENLOG DPREF RACE CAMP WITH X
  /DESIGN=DPREF RACE CAMP X.
```

- The variable *X* is a continuous variable specified as a cell covariate. Cell covariates must be specified after the keyword `WITH` following the variable list. No range is defined for cell covariates.
- To include the cell covariate in the model, the variable *X* is specified on `DESIGN`.

### **CSTRUCTURE Subcommand**

`CSTRUCTURE` specifies the variable that contains values for computing cell weights, such as structural zero indicators. By default, cell weights are equal to 1.

- The specification must be a numeric variable.
- Variables specified as dependent or independent variables in the variable list cannot be specified on `CSTRUCTURE`.
- Cell weights are not applied on a case-by-case basis. The weighted mean for a cell is applied to that cell.
- `CSTRUCTURE` can be used to impose structural, or *a priori*, zeros on the model. This feature is useful in specifying a quasi-symmetry model and in excluding cells from entering into estimation.
- If multiple `CSTRUCTURE` subcommands are specified, the last specification takes effect.

#### **Example**

```
COMPUTE CWT=(HUSED NE WIFED) .
GENLOG HUSED WIFED WITH DISTANCE
  /CSTRUCTURE=CWT
  /DESIGN=HUSED WIFED DISTANCE.
```

- The Boolean expression assigns *CWT* the value of 1 when *HUSED* is not equal to *WIFED*, and the value of 0 otherwise.
- `CSTRUCTURE` imposes structural zeros on the diagonal of the symmetric crosstabulation.

## ***GRESID Subcommand***

GRESID (Generalized Residual) calculates linear combinations of observed and expected cell frequencies as well as simple, standardized, and adjusted residuals.

- The variables specified must be numeric, and they must contain coefficients of the desired linear combinations.
- Variables specified as dependent or independent variables in the variable list cannot be specified on GRESID.
- The generalized residual coefficient is not applied on a case-by-case basis. The weighted coefficient mean of the value for all cases in a cell is applied to that cell.
- Each variable specified on the GRESID subcommand contains a single linear combination.
- If multiple GRESID subcommands are specified, the last specification takes effect.

### ***Example***

```
COMPUTE GR_1=(MONTH LE 6) .
COMPUTE GR_2=(MONTH GE 7) .
GENLOG MONTH WITH Z
  /GRESID=GR_1 GR_2
  /DESIGN=Z .
```

- The first variable, *GR\_1*, combines the first six months into a single effect; the second variable, *GR\_2*, combines the rest of the months.
- For each effect, GENLOG displays the observed and expected counts as well as the simple, standardized, and adjusted residuals.

## ***GLOR Subcommand***

GLOR (Generalized Log-Odds Ratio) specifies the population contrast variable(s). For each variable specified, GENLOG tests the null hypothesis that the generalized log-odds ratio equals 0 and displays the Wald statistic and the confidence interval. You can specify the level of the confidence interval using the CIN significance-level keyword on CRITERIA. By default, the confidence level is 95%.

- The variable sum is 0 for the loglinear model and for each combined level of independent variables for the logit model.
- Variables specified as dependent or independent variables in the variable list cannot be specified on GLOR.
- The coefficient is not applied on a case-by-case basis. The weighted mean for a cell is applied to that cell.
- If multiple GLOR subcommands are specified, the last specification takes effect.

### ***Example***

```
GENLOG A B
  /GLOR=COEFF
  /DESIGN=A B .
```



- The variable *COEFF* contains the coefficients of two dichotomous factors *A* and *B*.
- If the weighted cell mean for *COEFF* is 1 when *A* equals *B* and  $-1$  otherwise, this example tests whether the log-odds ratio equals 0, or in this case, whether variables *A* and *B* are independent.

## MODEL Subcommand

MODEL specifies the assumed distribution of your data.

- You can specify only one keyword on MODEL. The default is POISSON.
- If more than one MODEL subcommand is specified, the last specification takes effect.

<b>POISSON</b>	<i>The Poisson distribution.</i> This is the default.
<b>MULTINOMIAL</b>	<i>The multinomial distribution.</i> For the logit model, you must specify MULTINOMIAL.

## CRITERIA Subcommand

CRITERIA specifies the values used in tuning the parameters for the Newton-Raphson algorithm.

- If multiple CRITERIA subcommands are specified, the last specification takes effect.

<b>CONVERGE(n)</b>	<i>Convergence criterion.</i> Specify a positive value for the convergence criterion. The default is 0.001.
<b>ITERATE(n)</b>	<i>Maximum number of iterations.</i> Specify an integer. The default number is 20.
<b>DELTA(n)</b>	<i>Cell delta value.</i> Specify a non-negative value to add to each cell frequency for the first iteration. (For the saturated model, the delta value is added for all iterations.) The default is 0.5. The delta value is used to solve mathematical problems created by 0 observations; if all of your observations are greater than 0, we recommend that you set DELTA to 0.
<b>CIN(n)</b>	<i>Level of confidence interval.</i> Specify the percentage interval used in the test of generalized log-odds ratios and parameter estimates. The value must be between 50 and 99.99, inclusive. The default is 95.
<b>EPS(n)</b>	<i>Epsilon value used for redundancy checking in design matrix.</i> Specify a positive value. The default is 0.00000001.
<b>DEFAULT</b>	<i>Default values are used.</i> DEFAULT can be used to reset all criteria to default values.

### Example

```
GENLOG DPREF BY RACE ORIGIN CAMP
  /MODEL=MULTINOMIAL
  /CRITERIA=ITERATION(50) CONVERGE(.0001).
```

- ITERATION increases the maximum number of iterations to 50.
- CONVERGE lowers the convergence criterion to 0.0001.

## PRINT Subcommand

PRINT controls the display of statistics.

- By default, GENLOG displays the frequency table and simple, adjusted, and deviance residuals.

- When PRINT is specified with one or more keywords, only the statistics requested by these keywords are displayed.
- When multiple PRINT subcommands are specified, the last specification takes effect.

The following keywords can be used on PRINT:

<b>FREQ</b>	<i>Observed and expected cell frequencies and percentages.</i> This is displayed by default.
<b>RESID</b>	<i>Simple residuals.</i> This is displayed by default.
<b>ZRESID</b>	<i>Standardized residuals.</i>
<b>ADJRESID</b>	<i>Adjusted residuals.</i> This is displayed by default.
<b>DEV</b>	<i>Deviance residuals.</i> This is displayed by default.
<b>DESIGN</b>	<i>The design matrix of the model.</i> The design matrix corresponding to the specified model is displayed.
<b>ESTIM</b>	<i>The parameter estimates of the model.</i> The parameter estimates refer to the original categories.
<b>CORR</b>	<i>The correlation matrix of the parameter estimates.</i>
<b>COV</b>	<i>The covariance matrix of the parameter estimates.</i>
<b>ALL</b>	<i>All available output.</i>
<b>DEFAULT</b>	<i>FREQ, RESID, ADJRESID, and DEV.</i> This keyword can be used to reset PRINT to its default setting.
<b>NONE</b>	<i>The design and model information with goodness-of-fit statistics only.</i> This option overrides all other specifications on the PRINT subcommand.

### Example

```
GENLOG A B
  /PRINT=ALL
  /DESIGN=A B.
```

- The DESIGN subcommand specifies a main-effects model, which tests the hypothesis of no interaction. The PRINT subcommand displays all available output for this model.

## PLOT Subcommand

PLOT specifies which plots you want to display. Plots of adjusted residuals against observed and expected counts, and normal and detrended normal plots of the adjusted residuals are displayed if PLOT is not specified or is specified without a keyword. When multiple PLOT subcommands are specified, only the last specification is executed.

<b>DEFAULT</b>	<i>RESID (ADJRESID) and NORMPROB (ADJRESID).</i> This is the default if PLOT is not specified or is specified with no keyword.
<b>RESID (type)</b>	<i>Plots of residuals against observed and expected counts.</i> You can specify the type of residuals to plot. ADJRESID plots adjusted residuals; DEV plots deviance residuals. ADJRESID is the default if you do not specify a type.
<b>NORMPROB (type)</b>	<i>Normal and detrended normal plots of the residuals.</i> You can specify the type of residuals to plot. ADJRESID plots adjusted residuals; DEV plots deviance residuals. ADJRESID is the default if you do not specify a type.
<b>NONE</b>	<i>No plots.</i>

**Example**

```
GENLOG RESPONSE BY SEASON
  /MODEL=MULTINOMIAL
  /PLOT=RESID(ADJRESID,DEV)
  /DESIGN=RESPONSE SEASON(1) BY RESPONSE.
```

- This example requests plots of adjusted and deviance residuals against observed and expected counts.
- Note that if you specify `/PLOT=RESID(ADJRESID) RESID(DEV)`, only the deviance residuals are plotted. The first keyword specification, `RESID(ADJRESID)`, is ignored.

**MISSING Subcommand**

MISSING controls missing values. By default, GENLOG excludes all cases with system- or user-missing values for any variable. You can specify INCLUDE to include user-missing values.

- |                |   |
|----------------|---|
| <b>EXCLUDE</b> | <i>Delete cases with user-missing values.</i> This is the default if the subcommand is omitted. You can also specify the keyword <code>DEFAULT</code> . |
| <b>INCLUDE</b> | <i>Include cases with user-missing values.</i> Only cases with system-missing values are deleted.   |

**Example**

```
MISSING VALUES A(0).
GENLOG A B
  /MISSING=INCLUDE
  /DESIGN=B.
```

- Even though 0 was specified as missing, it is treated as a nonmissing category of *A* in this analysis.

**SAVE Subcommand**

SAVE saves specified temporary variables into the active dataset. You can assign a new name to each temporary variable saved.

- The temporary variables you can save include *RESID* (raw residual), *ZRESID* (standardized residual), *ADJRESID* (adjusted residual), *DEV* (deviance residual), and *PRED* (predicted cell frequency). An explanatory label is assigned to each saved variable.
- A temporary variable can be saved only once on a SAVE subcommand.
- To assign a name to a saved temporary variable, specify the new name in parentheses following that temporary variable. The new name must conform to IBM® SPSS® Statistics naming conventions and must be unique in the active dataset. The names cannot begin with # or \$.
- If you do not specify a variable name in parentheses, GENLOG assigns default names to the saved temporary variables. A default name starts with the first three characters of the name of the saved temporary variable, followed by an underscore and a unique number. For example, *RESID* will be saved as *RES\_n*, where *n* is a number incremented each time a default name is assigned to a saved *RESID*.

- The saved variables are pertinent to cells in the contingency table, not to individual observations. In the Data Editor, all cases that define one cell receive the same value. To make sense of these values, you need to aggregate the data to obtain cell counts.

### Example

```
GENLOG A B
  /SAVE PRED (PREDA_B)
  /DESIGN = A, B.
```

- `SAVE` saves the predicted values for two independent variables *A* and *B*.
- The saved variable is renamed *PREDA\_B* and added to the active dataset.

## DESIGN Subcommand

`DESIGN` specifies the model to be fit. If `DESIGN` is omitted or used with no specifications, the saturated model is produced. The saturated model fits all main effects and all interaction effects.

- Only one design can be specified on the subcommand.
- To obtain main-effects models, name all of the variables listed on the variables specification.
- To obtain interactions, use the keyword `BY` or an asterisk (`*`) to specify each interaction, for example, `A BY B` or `C*D`. To obtain the single-degree-of-freedom partition of a specified factor, specify the partition in parentheses following the factor (see the example below).
- To include cell covariates in the model, first identify them on the variable list by naming them after the keyword `WITH`, and then specify the variable names on `DESIGN`.
- Effects that involve only independent variables result in redundancy. `GENLOG` removes these effects from the model.
- If your variable list includes a cell covariate (identified by the keyword `WITH`), you cannot imply the saturated model by omitting `DESIGN` or specifying it alone. You need to request the model explicitly by specifying all main effects and interactions on `DESIGN`.

### Example

```
COMPUTE X=MONTH.
GENLOG MONTH WITH X
  /DESIGN X.
```

- This example tests the linear effect of the dependent variable.
- The variable specification identifies *MONTH* as a categorical variable. The keyword `WITH` identifies *X* as a covariate.
- `DESIGN` tests the linear effect of *MONTH*.

### Example

```
GENLOG A B
  /DESIGN=A.
```

```
GENLOG A B
  /DESIGN=A, B.
```

- Both designs specify main-effects models.
- The first design tests the homogeneity of category probabilities for  $B$ ; it fits the marginal frequencies on  $A$  but assumes that membership in any of the categories of  $B$  is equiprobable.
- The second design tests the independence of  $A$  and  $B$ . It fits the marginals on both  $A$  and  $B$ .

### **Example**

```
GENLOG A B C
  /DESIGN=A,B,C, A BY B.
```

- This design consists of the  $A$  main effect, the  $B$  main effect, the  $C$  main effect, and the interaction of  $A$  and  $B$ .

### **Example**

```
GENLOG A BY B
  /MODEL=MULTINOMIAL
  /DESIGN=A, A BY B(1).
```

- This example specifies single-degree-of-freedom partitions.
- The value 1 following  $B$  to the first category of  $B$ .

### **Example**

```
GENLOG HUSED WIFED WITH DISTANCE
  /DESIGN=HUSED WIFED DISTANCE.
```

- The continuous variable *DISTANCE* is identified as a cell covariate by the keyword *WITH*. The cell covariate is then included in the model by naming it on *DESIGN*.

### **Example**

```
COMPUTE X=1.
GENLOG MONTH WITH X
  /DESIGN=X.
```

- This example specifies an equiprobability model.
- The design tests whether the frequencies in the table are equal by using a constant of 1 as a cell covariate.

## **References**

Haberman, S. J. 1982. Analysis of dispersion of multinomial responses. *Journal of the American Statistical Association*, 77, 568–580.

# GET

```
GET FILE='file' [LOCK={YES**}]
                    {NO  }
[/KEEP={ALL**  }] [/DROP=varlist]
                    {varlist}
[/RENAME=(old varnames=new varnames)...]
[/MAP]
```

\*\*Default if the subcommand is omitted.

## Release History

Release 17.0

- LOCK keyword introduced on the FILE subcommand.

## Example

```
GET FILE='/data/empl.sav'.
```

## Overview

GET reads IBM® SPSS® Statistics data files. It also reads SPSS/PC+ data files, but you should not read SPSS/PC+ data files in Unicode mode (see Operations below).

GET is used only for reading SPSS Statistics data files. See DATA LIST for information on reading and defining data in a text data file. See MATRIX DATA for information on defining matrix materials in a text data file. For information on defining complex data files that cannot be defined with DATA LIST alone, see FILE TYPE and REPEATING DATA.

The program can also read data files created for other software applications. See IMPORT for information on reading *portable files* created with EXPORT. See the relevant commands, such as GET DATA, GET STATA, and GET SAS, for information on reading files created by other software programs.

## Options

**Variable Subsets and Order.** You can read a subset of variables and reorder the variables that are copied into the active dataset using the DROP and KEEP subcommands.

**Variable Names.** You can rename variables as they are copied into the active dataset with the RENAME subcommand.

**Variable Map.** To confirm the names and order of variables in the active dataset, use the MAP subcommand.

**Basic Specification**

- The basic specification is the `FILE` subcommand, which specifies the SPSS Statistics data file to be read.

**Subcommand Order**

- `FILE` must be specified first.
- The remaining subcommands can be specified in any order.

**Syntax Rules**

- `FILE` is required and can be specified only once.
- `KEEP`, `DROP`, `RENAME`, and `MAP` can be used as many times as needed.
- `GET` cannot be used inside a `DO IF-END IF` or `LOOP-END LOOP` structure.

**Operations**

- If `KEEP` is not specified, variables in the active dataset are in the same order as the original data file.
- A file saved with weighting in effect maintains weighting the next time the file is accessed. For a discussion of turning off weights, see `WEIGHT`.
- In Unicode mode, for code page data files and data files created in releases prior to 16.0, the defined width of string variables is tripled. You can use `ALTER TYPE` to automatically adjust the width of all string variables. See `SET command`, `UNICODE subcommand` for more information.
- In Unicode mode, SPSS/PC+ data files may not be read correctly.
- SPSS/PC+ data files can be read only on Windows operating systems.

## **FILE Subcommand**

`FILE` specifies the IBM® SPSS® Statistics data file to be read. `FILE` is required and can be specified only once. It must be the first specification on `GET`.

**LOCK Keyword**

The `LOCK` keyword only applies when opening a file from a IBM® SPSS® Collaboration and Deployment Services Repository. It specifies whether the file is locked (in the repository) when opened. Locking a file prevents other users from modifying it while you are working with it. Other users can open locked files but they cannot save them back to the repository. The lock is released when the file is closed or when the current session ends.

- By default, files are opened as locked.
- The `LOCK` keyword is ignored if the specified file does not reside in a IBM SPSS Collaboration and Deployment Services Repository.

For more information, see the topic [File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects](#) on p. 2230.

## ***DROP and KEEP Subcommands***

`DROP` and `KEEP` are used to copy a subset of variables into the active dataset. `DROP` specifies variables that should not be copied into the active dataset. `KEEP` specifies variables that should be copied. Variables not specified on `KEEP` are dropped.

- Variables can be specified in any order. The order of variables on `KEEP` determines the order of variables in the active dataset. The order of variables on `DROP` does not affect the order of variables in the active dataset.
- The keyword `ALL` on `KEEP` refers to all remaining variables not previously specified on `KEEP`. `ALL` must be the last specification on `KEEP`.
- If a variable is specified twice on the same subcommand, only the first mention is recognized.
- Multiple `DROP` and `KEEP` subcommands are allowed. However, specifying a variable named on a previous `DROP` or not named on a previous `KEEP` results in an error, and the `GET` command is not executed.
- The keyword `TO` can be used to specify a group of consecutive variables in the data file.

### ***Example***

```
GET FILE='/data/hubtemp.sav' /DROP=DEPT79 TO DEPT84 SALARY79.
```

- All variables between and including `DEPT79` and `DEPT84`, as well as `SALARY79`, are excluded from the active dataset. All other variables are included in the active dataset.
- Variables in the active dataset are in the same order as the variables in the original file.

### ***Example***

```
GET FILE='/data/prsnl.sav' /DROP=GRADE STORE
/KEEP=LNAME NAME TENURE JTENURE ALL.
```

- The variables `GRADE` and `STORE` are dropped when the file `prsnl.sav` is copied into the active dataset.
- `KEEP` specifies that `LNAME`, `NAME`, `TENURE`, and `JTENURE` are the first four variables in the active dataset, followed by all remaining variables (except those dropped by the previous `DROP` subcommand). These remaining variables are copied into the active dataset in the same sequence in which they appear in the `prsnl.sav` file.

## ***RENAME Subcommand***

`RENAME` changes the names of variables as they are copied into the active dataset.

- The specification on `RENAME` is a list of old variable names followed by an equals sign and a list of new variable names. The same number of variables must be specified on both lists. The keyword `TO` can be used on the first list to refer to consecutive variables in the data file and on the second list to generate new variable names. The entire specification must be enclosed in parentheses.



- Alternatively, you can specify each old variable name individually, followed by an equals sign and the new variable name. Multiple sets of variable specifications are allowed. The parentheses around each set of specifications are optional.
- Old variable names do not need to be specified according to their order in the data file.
- Name changes take place in one operation. Therefore, variable names can be exchanged between two variables.
- Variables cannot be renamed to scratch variables.
- Multiple RENAME subcommands are allowed.
- On a subsequent DROP or KEEP subcommand, variables are referred to by their new names.

### Example

```
GET FILE='/data/empl88.sav'  
  /RENAME AGE=AGE88 JOBCAT=JOBCAT88.
```

- RENAME specifies two name changes for the active dataset. *AGE* is renamed to *AGE88* and *JOBCAT* is renamed to *JOBCAT88*.

### Example

```
GET FILE='/data/empl88.sav'  
  /RENAME (AGE JOBCAT=AGE88 JOBCAT88).
```

- The name changes are identical to those in the previous example. *AGE* is renamed to *AGE88* and *JOBCAT* is renamed to *JOBCAT88*. The parentheses are required with this method.

## MAP Subcommand

MAP displays a list of the variables in the active dataset and their corresponding names in the original data file.

- The only specification is the keyword MAP. There are no additional specifications.
- Multiple MAP subcommands are allowed. Each MAP subcommand maps the results of subcommands that precede it; results of subcommands that follow it are not mapped.

### Example

```
GET FILE='/data/empl88.sav'  
  /RENAME=(AGE=AGE88) (JOBCAT=JOBCAT88)  
  /KEEP=LNAME NAME JOBCAT88 ALL /MAP.
```

- MAP is specified to confirm the new names for the variables *AGE* and *JOBCAT* and the order of variables in the active dataset (*LNAME*, *NAME*, and *JOBCAT88*, followed by all remaining variables in the data file).

# GET CAPTURE

GET CAPTURE is supported for compatibility purposes. GET DATA is the preferred command for reading databases. For more information, see the topic GET DATA on p. 839.

```
GET CAPTURE {ODBC }*

[/CONNECT='connection string']
[/UNENCRYPTED]
[/LOGIN=login] [/PASSWORD=password]
[/SERVER=host] [/DATABASE=database name]†

/SQL 'select statement'
    ['continuation of select statement']
```

\* You can import data from any database for which you have an ODBC driver installed.

† Optional subcommands are database-specific. For more information, see [Overview](#) below.

Release 15.0

- UNENCRYPTED subcommand introduced.

## Example

```
GET CAPTURE ODBC
  /CONNECT='DSN=sales.mdb;DBQ=/data/saledata.mdb;DriverId=281;FIL=MS'+
  ' Access;MaxBufferSize=2048;PageTimeout=5;'
  /SQL = 'SELECT T0.ID AS ID`, `T0.JOBCAT AS JOBCAT, `
  '`T0`.`REGION` AS `REGION`, `T0`.`DIVISION` AS `DIVISION`, `T0`.`TRAVEL`
  '` AS `TRAVEL`, `T0`.`SALES` AS `SALES`, `T0`.`VOLUME96` AS `VOLUME96`,
  '`T1`.`REGION` AS `REGION1`, `T1`.`AVGINC` AS `AVGINC`, `T1`.`AVGAGE` AS
  '`AVGAGE`, `T1`.`POPULAT` AS `POPULAT` FROM { oj `Regions` `T1` LEFT
  `OUTER JOIN `EmployeeSales` `T0` ON `T1`.`REGION` = `T0`.`REGION` } `.
```

## Overview

GET CAPTURE retrieves data from a database and converts them to a format that can be used by program procedures. GET CAPTURE retrieves data and data information and builds an active dataset for the current session.

*Note:* Although GET CAPTURE is still supported, equivalent functionality and additional features are provided in the newer GET DATA command.

### Basic Specification

The basic specification is one of the subcommands specifying the database type followed by the SQL subcommand and any select statement in quotation marks or apostrophes. Each line of the select statement should be enclosed in quotation marks or apostrophes, and no quoted string should exceed 255 characters.

### ***Subcommand Order***

The subcommand specifying the type of database must be the first specification. The `SQL` subcommand must be the last.

### ***Syntax Rules***

- Only one subcommand specifying the database type can be used.
- The `CONNECT` subcommand must be specified if you use the Microsoft ODBC (Open Database Connectivity) driver.

### ***Operations***

- `GET CAPTURE` retrieves the data specified on `SQL`.
- The variables are in the same order in which they are specified on the `SQL` subcommand.
- The data definition information captured from the database is stored in the active dataset dictionary.

### ***Limitations***

- A maximum of 3,800 characters (approximately) can be specified on the `SQL` subcommand. This translates to 76 lines of 50 characters. Characters beyond the limit are ignored.

## ***CONNECT Subcommand***

`CONNECT` is required to access any database that has an installed Microsoft ODBC driver.

- You cannot specify the connection string directly in the syntax window, but you can paste it with the rest of the command from the Results dialog box, which is the last of the series of dialog boxes opened with the Database Wizard.

## ***UNENCRYPTED Subcommands***

`UNENCRYPTED` allows unencrypted passwords to be used in the `CONNECT` subcommand. By default, passwords are assumed to be encrypted.

## ***SQL Subcommand***

`SQL` specifies any SQL select statement accepted by the database that you access. With ODBC, you can now select columns from more than one related table in an ODBC data source using either the inner join or the outer join.

## ***Data Conversion***

`GET CAPTURE` converts variable names, labels, missing values, and data types, wherever necessary, to a format that conforms to IBM® SPSS® Statistics conventions.

### **Variable Names and Labels**

Database columns are read as variables.

- A column name is converted to a variable name if it conforms to variable naming conventions and is different from all other names created for the active dataset. If not, `GET CAPTURE` gives the column a name formed from the first few letters of the column and its column number. If this is not possible, the letters `COL` followed by the column number are used. For example, the seventh column specified in the select statement could be `COL7`.
- `GET CAPTURE` labels each variable with its full column name specified in the original database.
- You can display a table of variable names with their original database column names using the `DISPLAY LABELS` command.

### **Missing Values**

Null values in the database are transformed into the system-missing value in numeric variables or into blanks in string variables.

# GET DATA

```
GET DATA
  /TYPE = {ODBC }
         {OLEDB}
         {XLS  }
         {XLSX }
         {XLSM }
         {TXT  }

  /FILE = 'filename'

Subcommands for TYPE = ODBC and OLEDB
/CONNECT='connection string'
[{/ENCRYPTED  }]
{/UNENCRYPTED}
/SQL 'select statement'
    ['select statement continued']

Subcommands for TYPE=ODBC, TYPE=OLEDB, XLS, XLSX, and XLSM
[/ASSUMEDSTRWIDTH={255**}]
    {n      }

Subcommands for TYPE = XLS, XLSX, and XLSM*
[/SHEET = {INDEX**} {sheet number}]
    {NAME  } {'sheet name'}
[/CELLRANGE = {RANGE } {'start point:end point' }]
    {FULL**}
[/READNAMES = {on** } ]
    {off  }

Subcommands for TYPE = TXT
[/ARRANGEMENT = {FIXED  }]
    {DELIMITED**}
[/FIRSTCASE = {n}]
[/DELCASE = {LINE**  } ]1
    {VARIABLES n}

[/FIXCASE = n]
[/IMPORTCASE = {ALL**  }]
    {FIRST n }
    {PERCENT n}
[/DELIMITERS = {"delimiters"}]
[/QUALIFIER = "qualifier"]

VARIABLES subcommand for ARRANGEMENT = DELIMITED
/VARIABLES = varname {format}

VARIABLES subcommand for ARRANGEMENT = FIXED
/VARIABLES varname {startcol - endcol} {format}
    {/rec#} varname {startcol - endcol} {format}
```

\*For Excel 4.0 or earlier files, use GET TRANSLATE.

\*\*Default if the subcommand is omitted.

**Release history**

## Release 13.0

- ASSUMEDSTRWIDTH subcommand introduced for TYPE=ODBC.

## Release 14.0

- ASSUMEDSTRWIDTH subcommand extended to TYPE=XLS.
- TYPE=OLEDB introduced.

## Release 15.0

- ASSUMEDSTRWIDTH subcommand extended to TYPE=OLEDB.

## Release 16.0

- TYPE=XLSX and TYPE=XLSM introduced.

## Release 17.0

- ENCRYPTED subcommand introduced.

**Example**

```
GET DATA
  /TYPE=XLS
  /FILE='/PlanningDocs/files10.xls'
  /SHEET=name 'First Quarter'
  /CELLRANGE=full
  /READNAMES=on.
```

**Overview**

GET DATA reads data from ODBC OLE DB data sources (databases), Excel files (release 5 or later), and text data files. It contains functionality and syntax similar to GET CAPTURE, GET TRANSLATE, and DATA LIST.

- GET DATA /TYPE=ODBC is almost identical to GET CAPTURE ODBC in both syntax and functionality.
- GET DATA /TYPE=XLS reads Excel 95 through Excel 2003 files; GET DATA /TYPE=XLSX and GET DATA /TYPE=XLSM read Excel 2007 or later files. GET TRANSLATE reads Excel 4 or earlier, Lotus, and dBASE files.
- GET DATA /TYPE=TXT is similar to DATA LIST but does not create a temporary copy of the data file, significantly reducing temporary file space requirements for large data files.

**Limitations****TYPE Subcommand**

The `TYPE` subcommand is required and must be the first subcommand specified.

<b>ODBC</b>	<i>Data sources accessed with ODBC drivers.</i>
<b>OLEDB</b>	<i>Data sources accessed with Microsoft OLEDB technology. Available only on Windows platforms and requires .NET framework and IBM® SPSS® Data Collection Survey Reporter Developer Kit. For information on obtaining a compatible version of SPSS Survey Reporter Developer Kit, go to <a href="http://support.spss.com">support.spss.com</a> (<a href="http://support.spss.com">http://support.spss.com</a>).</i>
<b>XLS</b>	<i>Excel 95 through Excel 2003 files. For earlier versions of Excel files, Lotus 1-2-3 files, and dBASE files, see the <code>GET TRANSLATE</code> command.</i>
<b>XLSX and XLSM</b>	<i>Excel 2007 files. Macros in XLSM files are ignored. XLSB (binary) format files are not supported.</i>
<b>TXT</b>	<i>Simple (ASCII) text data files. The encoding used to read text data files is UTF-8 in Unicode mode or the code page determined by the current locale in code page mode. For more information, see the topic <code>UNICODE Subcommand</code> on p. 1878.</i>

**FILE subcommand**

The `FILE` subcommand is required for `TYPE=XLS`, `TYPE=XLSX`, `TYPE=XLSM`, and `TYPE=TXT` and must immediately follow the `TYPE` subcommand. It specifies the file to read. File specifications should be enclosed in quotes.

**Subcommands for TYPE=ODBC and TYPE=OLEDB**

The `CONNECT` and `SQL` subcommands are both required, and `SQL` must be the last subcommand.

**Example**

```
GET DATA /TYPE=ODBC
  /CONNECT=
  'DSN=MS Access Database;DBQ=/examples/data/dm_demo.mdb;' +
  'DriverId=25;FIL=MS Access;MaxBufferSize=2048;PageTimeout=5;'
  /SQL = 'SELECT * FROM CombinedTable'.
```

Note: If you are running the Windows 64-bit version of IBM® SPSS® Statistics, you cannot read Excel, Access, or dBASE database sources with `GET DATA /TYPE=ODBC`. The 32-bit ODBC drivers for these products are not compatible. You can, however, read Excel files with `GET DATA /TYPE=XLS`.

**CONNECT subcommand**

The `CONNECT` subcommand identifies the database source. The recommended method for generating a valid `CONNECT` specification is to initially use the Database Wizard and paste the resulting syntax to a syntax window in the last step of the wizard.

- The entire connect string must be enclosed in quotation marks.
- For long connect strings, you can use multiple quoted strings on separate lines, using a plus sign (+) to combine the quoted strings.

**ENCRYPTED and UNENCRYPTED subcommands**

`UNENCRYPTED` allows unencrypted passwords to be used in the `CONNECT` subcommand. The default is `ENCRYPTED`.

**SQL subcommand**

`SQL` specifies any SQL select statement accepted by the database that you access.

- You can select columns from more than one related table in a data source using either the inner join or the outer join.
- Each line of SQL must be enclosed in quotation marks and cannot exceed 255 characters.
- When the command is processed, all of the lines of the SQL statement are merged together in a very literal fashion; so each line should either begin or end with a blank space where spaces should occur between specifications.
- For `TYPE=OLEDB` (available only on Windows operating systems), table joins are not supported; you can specify fields only from a single table.

**Example**

```
GET DATA /TYPE=ODBC /CONNECT=
  'DSN=Microsoft Access;DBQ=/data/demo.mdb;DriverId=25;'+
  'FIL=MS Access;MaxBufferSize=2048;PageTimeout=5;'
/SQL =
'SELECT SurveyResponses.ID, SurveyResponses.Internet, '
' [Value Labels].[Internet Label]'
' FROM SurveyResponses LEFT OUTER JOIN [Value Labels]'
' ON SurveyResponses.Internet'
' = [Value Labels].[Internet Value]'.
```

If the SQL contains `WHERE` clauses with expressions for case selection, dates and times in expressions need to be specified in a special manner (including the curly braces shown in the examples):

- Date literals should be specified using the general form {d 'YYYY-mm-dd'}.
- Time literals should be specified using the general form {t 'hh:mm:ss'}.



- Date/time literals (timestamps) should be specified using the general form {ts 'yyyy-mm-dd hh:mm:ss'}.
- The entire date and/or time value must be enclosed in single quotes. Years must be expressed in four-digit form, and dates and times must contain two digits for each portion of the value. For example January 1, 2005, 1:05 AM would be expressed as:

```
{ts '2005-01-01 01:05:00'}
```

For functions used in expressions, a list of standard functions is available at [http://msdn.microsoft.com/en-us/library/ms711813\(VS.85\).aspx](http://msdn.microsoft.com/en-us/library/ms711813(VS.85).aspx).

## ***ASSUMEDSTRWIDTH subcommand***

For TYPE=ODBC, TYPE=OLEDB, and TYPE=XLS, this controls the width of variable-width string values. By default, the width is 255 bytes, and only the first 255 bytes will be read. The width can be up to 32,767 bytes. Although you probably don't want to truncate string values, you also don't want to specify an unnecessarily large value, since this will be used as the display width for those string values.

## ***Subcommands for TYPE=XLS, XLSX, and XLSM***

For Excel 95 or later files, you can specify a spreadsheet within the workbook, a range of cells to read, and the contents of the first row of the spreadsheet (variable names or data). For files from earlier versions of Excel, use GET TRANSLATE .

### ***Example***

```
GET DATA /TYPE=XLS
  /FILE='/data/sales.xls'
  /SHEET=name 'June Sales'
  /CELLRANGE=range 'A1:C3'
  /READNAMES=on.
```

## ***SHEET subcommand***

The SHEET subcommand indicates the worksheet in the Excel file that will be read. Only one sheet can be specified. If no sheet is specified, the first sheet will be read.

- |                    |  |
|--------------------|--|
| <b>INDEX n</b>     | <i>Read the specified sheet number.</i> The number represents the sequential order of the sheet within the workbook. |
| <b>NAME 'name'</b> | <i>Read the specified sheet name.</i> If the name contains spaces, it must be enclosed in quotes.                    |

**CELLRANGE subcommand**

The CELLRANGE subcommand specifies a range of cells to read within the specified worksheet. By default, the entire worksheet is read.

<b>FULL</b>	<i>Read the entire worksheet.</i> This is the default.
<b>RANGE 'start:end'</b>	<i>Read the specified range of cells.</i> Specify the beginning column letter and row number, a colon, and the ending column letter and row number, as in A1:K14. The cell range must be enclosed in quotes.

**READNAMES subcommand**

<b>ON</b>	<i>Read the first row of the sheet or specified range as variable names.</i> This is the default. Values that contain invalid characters or do not meet other criteria for variable names are converted to valid variable names. <a href="#">For more information, see the topic Variable Names on p. 49.</a>
<b>OFF</b>	<i>Read the first row of the sheet or specified range as data.</i> Default variable names are assigned, and all rows are read as data.

**Subcommands for TYPE=TXT**

The VARIABLES subcommand is required and must be the last GET DATA subcommand.

**Example**

```
GET DATA /TYPE = TXT
  /FILE = '/data/textdata.dat'
  /DELCASE = LINE
  /DELIMITERS = "\t ,"
  /ARRANGEMENT = DELIMITED
  /FIRSTCASE = 2
  /IMPORTCASE = FIRST 200
  /VARIABLES = id F3.0 gender A1 bdate DATE10 educ F2.0
  jobcat F1.0 salary DOLLAR8 salbegin DOLLAR8
  jobtime F4.2 prevexp F4.2 minority F3.0.
```

**ARRANGEMENT subcommand**

The ARRANGEMENT subcommand specifies the data format.

<b>DELIMITED</b>	<i>Spaces, commas, tabs, or other characters are used to separate variables.</i> The variables are recorded in the same order for each case but not necessarily in the same column locations. This is the default.
<b>FIXED</b>	<i>Each variable is recorded in the same column location for every case.</i>

**FIRSTCASE subcommand**

FIRSTCASE specifies the first line (row) to read for the first case of data. This allows you to bypass information in the first *n* lines of the file that either don't contain data or contain data that you don't want to read. This subcommand applies to both fixed and delimited file formats.

The only specification for this subcommand is an integer greater than zero that indicates the number of lines to skip. The default is 1.

### ***DELCASE subcommand***

The DELCASE subcommand applies to delimited data (ARRANGEMENT=DELIMITED) only.

<b>LINE</b>	<i>Each case is contained on a single line (row).</i> This is the default.
<b>VARIABLES n</b>	<i>Each case contains n variables.</i> Multiple cases can be contained on the same line, and data for one case can span more than one line. A case is defined by the number of variables.

### ***FIXCASE subcommand***

The FIXCASE subcommand applies to fixed data (ARRANGEMENT=FIXED) only. It specifies the number of lines (records) to read for each case.

The only specification for this subcommand is an integer greater than zero that indicates the number of lines (records) per case. The default is 1.

### ***IMPORTCASES subcommand***

The IMPORTCASES subcommand allows you to specify the number of cases to read.

<b>ALL</b>	<i>Read all cases in the file.</i> This is the default.
<b>FIRST n</b>	<i>Read the first n cases.</i> The value of <i>n</i> must be a positive integer.
<b>PERCENT n</b>	<i>Read approximately the first n percent of cases.</i> The value of <i>n</i> must be a positive integer less than 100. The percentage of cases actually selected only approximates the specified percentage. The more cases there are in the data file, the closer the percentage of cases selected is to the specified percentage.

### ***DELIMITERS subcommand***

The DELIMITERS subcommand applies to delimited data (ARRANGEMENT=DELIMITED) only. It specifies the characters to read as delimiters between data values.

- Each delimiter can be only a single character, except for the specification of a tab or a backslash as a delimiter (see below).
- The list of delimiters must be enclosed in quotes.
- There should be no spaces or other delimiters between delimiter specifications, except for a space that indicates a space as a delimiter.
- To specify a tab as a delimiter use "\t". This must be the first delimiter specified.
- To specify a backslash as a delimiter, use two backslashes ("\\"). This must be the first delimiter specified unless you also specify a tab as a delimiter, in which case the backslash specification should come second—immediately after the tab specification.

**Missing data with delimited data.** Multiple consecutive spaces in a data file are treated as a single space and cannot be used to indicate missing data. For any other delimiter, multiple delimiters without any intervening data indicate missing data.

**Example**

```
DELIMITERS "\t\ , ;"
```

In this example, tabs, backslashes, spaces, commas, and semicolons will be read as delimiters between data values.

### **QUALIFIER subcommand**

The `QUALIFIERS` subcommand applies to delimited data (`ARRANGEMENT=DELIMITED`) only. It specifies the character used to enclose values that contain delimiter characters. For example, if a comma is the delimiter, values that contain commas will be read incorrectly unless there is a text qualifier enclosing the value, preventing the commas in the value from being interpreted as delimiters between values. CSV-format data files exported from Excel use a double quote (") as a text qualifier.

- The text qualifier appears at both the beginning and end of the value, enclosing the entire value.
- The qualifier value must be enclosed in single or double quotes. If the qualifier is a single quote, the value should be enclosed in double quotes. If the qualifier value is a double quote, the value should be enclosed in single quotes.

**Example**

```
/QUALIFIER = ` `
```

### **VARIABLES subcommand for ARRANGEMENT = DELIMITED**

For delimited files, the `VARIABLES` subcommand specifies the variable names and variable formats.

- Variable names must conform to variable naming rules. [For more information, see the topic Variable Names on p. 49.](#)
- Each variable name must be followed by a format specification. [For more information, see the topic Variable Format Specifications for TYPE = TXT on p. 847.](#)

### **VARIABLES subcommand for ARRANGEMENT = FIXED**

For fixed-format files, the `VARIABLES` subcommand specifies variable names, start and end column locations, and variable formats.

- Variable names must conform to variable naming rules. [For more information, see the topic Variable Names on p. 49.](#)
- Each variable name must be followed by column specifications. Start and end columns must be separated by a dash, as in 0-10.

- Column specifications must include both the start and end column positions, even if the width is only one column, as in 32–32.
- Each column specification must be followed by a format specification.
- Column numbering starts with 0, not 1 (in contrast to `DATA LIST`).

**Multiple records.** If each case spans more than one record (as specified with the `FIXCASE` subcommand), delimit variable specifications for each record with a slash (/) followed by the record number, as in:

```
VARIABLES = /1 var1 0-10 F var2 11-20 DATE
           /2 var3 0-5 A var4 6-10 F
           /3 var5 0-20 A var6 21-30 DOLLAR
```

### **Variable Format Specifications for `TYPE = TXT`**

For both fixed and delimited files, available formats include (but are not limited to):

<b>F<math>n</math>.<math>d</math></b>	<i>Numeric.</i> Specification of the total number of characters ( $n$ ) and decimals ( $d$ ) is optional.
<b>A<math>n</math></b>	<i>String (alphanumeric).</i> Specification of the maximum string length ( $n$ ) is optional.
<b>DATE<math>n</math></b>	<i>nDates of the general format dd-mmm-yyyy.</i> Specification of the maximum length ( $n$ ) is optional but must be eight or greater if specified.
<b>ADATE<math>n</math></b>	<i>Dates of the general format mm/dd/yyyy.</i> Specification of the maximum length ( $n$ ) is optional but must be eight or greater if specified.
<b>DOLLAR<math>n</math>.<math>d</math></b>	<i>Currency with or without a leading dollar sign (\$).</i> Input values can include a leading dollar sign, but it is not required. Specification of the total number of characters ( $n$ ) and decimals ( $d$ ) is optional.

For a complete list of variable formats, see [Variable Types and Formats on p. 54](#).

*Note:* For default numeric (F) format and scientific notation (E) format, the decimal indicator of the input data must match the IBM® SPSS® Statistics locale decimal indicator (period or comma). Use `SHOW DECIMAL` to display the current decimal indicator and `SET DECIMAL` to set the decimal indicator. (Comma and Dollar formats recognize only the period as the decimal indicator, and Dot format recognizes only the comma as the decimal indicator.)

# GET SAS

```
GET SAS DATA='file' [DSET(dataset)]  
  
  [/ENCODING={"Locale" **}  
             {"System"  }  
             {"UTF-8"   }  
             {"value"   }  
  
  [/FORMATS=file]
```

\*\*Default if subcommand omitted.

## **Release History**

Release 19

- ENCODING subcommand introduced.

## **Example**

```
GET SAS DATA='/data/elect.sd7'.
```

## **Overview**

GET SAS builds an active dataset from a SAS dataset or a SAS transport file. A SAS transport file is a sequential file written in SAS transport format and can be created by the SAS export engine available in SAS Release 6.06 or higher or by the EXPORT option on the COPY or XCOPY procedure in earlier versions. GET SAS reads SAS version 6-9 files.

### **Options**

**Retrieving User-Defined Value Labels.** For native SAS datasets, you can specify a file on the FORMATS subcommand to retrieve user-defined value labels associated with the data being read. This file must be created by the SAS PROC FORMAT statement and can be used only for native SAS datasets. For SAS transport files, the FORMATS subcommand is ignored.

**Specifying the Dataset.** You can name a dataset contained in a specified SAS file, using DSET on the DATA subcommand. GET SAS reads the specified dataset from the SAS file.

### **Basic Specification**

The basic specification is the DATA subcommand followed by the name of the SAS file to read. By default, the first SAS dataset is copied into the active dataset and any necessary data conversions are made.

### **Syntax Rules**

- The subcommand DATA and the SAS filename are required and must be specified first.

- The subcommand `FORMATS` is optional. This subcommand is ignored for SAS transport files.
- The subcommand `ENCODING` is optional.
- `GET SAS` does not allow `KEEP`, `DROP`, `RENAME`, and `MAP` subcommands. To use a subset of the variables, rename them, or display the file content, you can specify the appropriate commands after the active dataset is created.

### **Operations**

- `GET SAS` reads data from the specified or default dataset contained in the SAS file named on the `DATA` subcommand.
- Value labels retrieved from a SAS user-defined format are used for variables associated with that format, becoming part of the IBM® SPSS® Statistics dictionary.
- All variables from the SAS dataset are included in the active dataset, and they are in the same order as in the SAS dataset.

## **DATA Subcommand**

`DATA` specifies the file that contains the SAS dataset to be read.

- `DATA` is required and must be the first specification on `GET SAS`.
- The file specification varies from operating system to operating system. File specifications should be enclosed in quotes.
- The optional `DSET` keyword on `DATA` determines which dataset within the specified SAS file is to be read. The default is the first dataset.

**DSET (dataset)**      *Dataset to be read.* Specify the name of the dataset in parentheses. If the specified dataset does not exist in the SAS file, `GET SAS` displays a message informing you that the dataset was not found.

### **Example**

```
GET SAS DATA='/data/elect.sd7' DSET(Y1948) .
```

- The SAS file *elect.sd7* is opened and the dataset named *Y1948* is used to build the active dataset for the session.

## **ENCODING Subcommand**

The `ENCODING` subcommand specifies the character encoding of the SAS data file.

- The encoding must be correctly identified or the file cannot be read.
- If there is a `formats` file, the encoding of the `formats` file must be the same as the encoding of the data file. [For more information, see the topic FORMATS Subcommand on p. 851.](#)
- The subcommand is followed by an optional equals sign and a quoted encoding value.

- The quoted value can be any of the values in the Encoding column in the [Character Encoding](#) table or “UTF-8” for UTF-8 Unicode encoding.
- The default encoding is “Locale”, which is the encoding of the current IBM® SPSS® Statistics locale. [For more information, see the topic LOCALE Subcommand on p. 1876.](#)

### Example

```
GET SAS DATA='/data/elect.sas7bdat'
  /ENCODING='Windows-1252'.
```

Table 100-1  
Character Encoding

Character Set	Encoding
SPSS Statistics Locale	Locale
Operating System Locale	System
Western European	
Western	ISO-8859-1
Western	ISO-8859-15
Western	IBM-850
Western	Windows-1252
Celtic	ISO-8859-14
Greek	ISO-8859-7
Greek	Windows-1253
Nordic	ISO-8859-10
Eastern European	
Baltic	Windows-1257
Central European	IBM-852
Central European	ISO-8859-2
Cyrillic	IBM-855
Cyrillic	ISO-8859-5
Cyrillic	Windows-1251
Cyrillic/Russian	CP-866
Eastern Asian	
Chinese Simplified	GBK
Chinese Simplified	ISO-2022-CN
Chinese Traditional	Big5
Chinese Traditional	EUC-TW
Japanese	EUC-JP
Japanese	ISO-2022-JP
Japanese	Shift-JIS
Korean	EUC-KR
Southeastern and Southwestern Asian	
Thai	Windows-874
Turkish	IBM-857
Turkish	ISO-8859-9
Middle Eastern	
Arabic	Windows-1256



Arabic	IBM-864
Hebrew	ISO-8859-8
Hebrew	Windows-1255
Hebrew	IBM-862

## FORMATS Subcommand

FORMATS specifies the file containing user-defined value labels to be applied to the retrieved data.

- File specifications should be enclosed in quotation marks.
- If FORMATS is omitted, no value labels are available.
- Value labels are applied only to numeric integer values. They are not applied to non-integer numeric values or string variables.
- The file specified on the FORMATS subcommand must be created with the SAS PROC FORMAT statement.
- For SAS transport files, the FORMATS subcommand is ignored.

### Example

```
GET SAS /DATA='/data/elect.sas7bdat' DSET(Y1948)
  /FORMATS='data/formats.sas7bcat'.
```

- Value labels read from the SAS catalog file *formats.sas7bcat* are converted to conform to IBM® SPSS® Statistics conventions.

## Creating a Formats File with PROC FORMAT

The following is an example of how to create a SAS format catalog file:

```
libname mylib 'c:\mydir\' ;

proc format library = mylib ;
value YesNo
  0='No'
  1='Yes' ;
value HighLow
  1 = 'Low'
  2 = 'Medium'
  3 = 'High' ;

options  fmtsearch=(mylib);

proc datasets library = mylib ;
modify mydata;
  format  var1 var2 var3 YesNo.;
  format  var4 var5 var6 HighLow.;

quit;
```

- libname defines a “library,” which is a directory path.
- proc format defines two formats: *YesNo* and *HighLow*. Each format defines a set of value labels associated with data values.

- `proc datasets` identifies the data file—*mydata*—and the variables to which each of the defined formats should be applied. So the *YesNo* format is applied to variables *var1*, *var2*, and *var3*, and the *HighLow* format is applied to the variables *var4*, *var5*, and *var6*.
- This creates the SAS catalog file *c:\mydir\formats.sas7bcat*.

## **SAS Data Conversion**

Although SAS and IBM® SPSS® Statistics data files have similar attributes, they are not identical. The following conversions are made to force SAS datasets to comply with SPSS Statistics conventions.

### **Variable Names**

SAS variable names that do not conform to IBM® SPSS® Statistics variable name rules are converted to valid variable names.

### **Variable Labels**

SAS variable labels specified on the `LABEL` statement in the `DATA` step are used as variable labels.

### **Value Labels**

SAS value formats that assign value labels are read from the dataset specified on the `FORMATS` subcommand. The SAS value labels are then converted to IBM® SPSS® Statistics value labels in the following manner:

- Labels assigned to single values are retained.
- Labels assigned to a range of values are ignored.
- Labels assigned to the SAS keywords `LOW`, `HIGH`, and `OTHER` are ignored.
- Labels assigned to string variables and non-integer numeric values are ignored.

### **Missing Values**

Since SAS has no user-defined missing values, all SAS missing codes are converted to IBM® SPSS® Statistics system-missing values.

### **Variable Types**

- Both SAS and IBM® SPSS® Statistics allow two basic types of variables: numeric and character string. During conversion, SAS numeric variables become SPSS Statistics numeric variables, and SAS string variables become SPSS Statistics string variables of the same length.
- Date, time, and date/time SAS variables are converted to equivalent SPSS Statistics date, time, and date/time variables. All other numeric formats are converted to the default numeric format.

# GET STATA

```
GET STATA FILE='file'  
    [/ENCODING={"Locale" **}]  
        {"System"  }  
        {"value"   }
```

\*\*Default if subcommand omitted.

## **Release History**

Release 14.0

- Command introduced.

Release 19

- ENCODING subcommand introduced.

## **Example**

```
GET STATA FILE=' /data/empl.dta' .
```

## **Overview**

GET STATA reads Stata-format data files created by Stata versions 4–8.

### **Basic Specification**

- The basic specification is the FILE keyword, which specifies the Stata data file to be read.

### **Operations**

- **Variable names.** Stata variable names are converted to IBM® SPSS® Statistics variable names in case-sensitive form. Stata variable names that are identical except for case are converted to valid variable names by appending an underscore and a sequential letter (*\_A*, *\_B*, *\_C*, ..., *\_Z*, *\_AA*, *\_AB*, ..., and so forth).
- **Variable labels.** Stata variable labels are converted to SPSS Statistics variable labels.
- **Value labels.** Stata value labels are converted to SPSS Statistics value labels, except for Stata value labels assigned to “extended” missing values.
- **Missing values.** Stata “extended” missing values are converted to system-missing values.
- **Date conversion.** Stata date format values are converted to SPSS Statistics DATE format (d-m-y) values. Stata “time-series” date format values (weeks, months, quarters, and so on) are converted to simple numeric (F) format, preserving the original, internal integer value, which is the number of weeks, months, quarters, and so on, since the start of 1960.

## FILE Keyword

FILE specifies the Stata data file to be read. FILE is the only specification; it is required and can be specified only once. The keyword name is followed by an equals sign and a quoted file specification (or quoted file handle) that specifies the Stata data file to read.

## ENCODING Subcommand

The ENCODING subcommand specifies the character encoding of the Stata data file.

- The encoding must be correctly identified or the file cannot be read.
- The subcommand is followed by an optional equals sign and a quoted encoding value.
- The quoted value can be any of the values in the Encoding column in the [Character Encoding table](#).
- The default encoding is “Locale”, which is the encoding of the current IBM® SPSS® Statistics locale. [For more information, see the topic LOCALE Subcommand on p. 1876.](#)

### Example

```
GET STATA FILE=' /data/empl.dta'
  /ENCODING='Windows-1252'.
```

Table 101-1  
Character Encoding

Character Set	Encoding
SPSS Statistics Locale	Locale
Operating System Locale	System
Western European	
Western	ISO-8859-1
Western	ISO-8859-15
Western	IBM-850
Western	Windows-1252
Celtic	ISO-8859-14
Greek	ISO-8859-7
Greek	Windows-1253
Nordic	ISO-8859-10
Eastern European	
Baltic	Windows-1257
Central European	IBM-852
Central European	ISO-8859-2
Cyrillic	IBM-855
Cyrillic	ISO-8859-5
Cyrillic	Windows-1251
Cyrillic/Russian	CP-866
Eastern Asian	
Chinese Simplified	GBK
Chinese Simplified	ISO-2022-CN

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Chinese Traditional	Big5
Chinese Traditional	EUC-TW
Japanese	EUC-JP
Japanese	ISO-2022-JP
Japanese	Shift-JIS
Korean	EUC-KR
Southeastern and Southwestern Asian	
Thai	Windows-874
Turkish	IBM-857
Turkish	ISO-8859-9
Middle Eastern	
Arabic	Windows-1256
Arabic	IBM-864
Hebrew	ISO-8859-8
Hebrew	Windows-1255
Hebrew	IBM-862

# GET TRANSLATE

```
GET TRANSLATE FILE=file

[/TYPE={WK }]
    {WK1}
    {WKS}
    {SYM}
    {SLK}
    {XLS}
    {DBF}
    {TAB}
    {SYS}

[/FIELDNAMES]*

[/RANGE={range name }]*
    {start..stop}
    {start:stop }

[/KEEP={ALL** }] [/DROP=varlist]
    {varlist}

[/MAP]
```

\*Available only for spreadsheet and tab-delimited ASCII files.

\*\*Default if the subcommand is omitted.

<b>Keyword</b>	<b>Type of file</b>
WK	Any Lotus 1-2-3 or Symphony file
WK1	1-2-3 Release 2.0
WKS	1-2-3 Release 1A
WR1	Symphony Release 2.0
WRK	Symphony Release 1.0
SLK	Microsoft Excel and Multiplan in SYLK (symbolic link) format
XLS	Microsoft Excel (for Excel 5 or later, use GET DATA)
DBF	All dBASE files
TAB	Tab-delimited ASCII file
SYS	Systat data file

## **Example**

```
GET TRANSLATE FILE='PROJECT.WKS'
  /FIELDNAMES
  /RANGE=D3..J279.
```

## **Overview**

GET TRANSLATE creates an active dataset from files produced by other software applications. Supported formats are 1-2-3, Symphony, Multiplan, Excel, dBASE II, dBASE III, dBASE IV, and tab-delimited ASCII files.

### **Options**

**Variable Subsets.** You can use the `DROP` and `KEEP` subcommands to specify variables to omit or retain in the resulting active dataset.

**Variable Names.** You can rename variables as they are translated using the `RENAME` subcommand.

**Variable Map.** To confirm the names and order of the variables in the active dataset, use the `MAP` subcommand. `MAP` displays the variables in the active dataset and their corresponding names in the other application.

**Spreadsheet Files.** You can use the `RANGE` subcommand to translate a subset of cells from a spreadsheet file. You can use the `FIELDNAMES` subcommand to translate field names in the spreadsheet file to variable names.

### **Basic Specification**

- The basic specification is `FILE` with a file specification enclosed in apostrophes.
- If the file's extension is not the default for the type of file you are reading, `TYPE` must also be specified.

### **Subcommand Order**

Subcommands can be named in any order.

### **Limitations**

The maximum number of variables that can be translated into the active dataset depends on the maximum number of variables that the other software application can handle:

<b>Application</b>	<b>Maximum variables</b>
1-2-3	256
Symphony	256
Multiplan	255
Excel	256
dBASE IV	255
dBASE III	128
dBASE II	32

## **Operations**

`GET TRANSLATE` replaces an existing active dataset.

### **Spreadsheets**

A spreadsheet file suitable for this program should be arranged so that each row represents a case and each column, a variable.

- By default, the new active dataset contains all rows and up to 256 columns from Lotus 1-2-3, Symphony, or Excel, or up to 255 columns from Multiplan.
- By default, GET TRANSLATE uses the column letters as variable names in the active dataset.
- The first row of a spreadsheet or specified range may contain field labels immediately followed by rows of data. These names can be transferred as IBM® SPSS® Statistics variable names. [For more information, see the topic FIELDNAMES Subcommand on p. 861.](#)
- The current value of a formula is translated to the active dataset.
- Blank, ERR, and NA values in 1-2-3 and Symphony and error values such as #N/A in Excel are translated as system-missing values in the active dataset.
- Hidden columns and cells in 1-2-3 Release 2 and Symphony files are translated and copied into the active dataset.
- Column width and format type are transferred to the dictionary of the active dataset.
- The format type is assigned from values in the first data row. By default, the first data row is row 1. If RANGE is specified, the first data row is the first row in the range. If FIELDNAMES is specified, the first data row follows immediately after the single row containing field names.
- If a cell in the first data row is empty, the variable is assigned the global default format from the spreadsheet.

The formats from 1-2-3, Symphony, Excel, and Multiplan are translated as follows:

1-2-3/Symphony	Excel	SYLK	SPSS Statistics
Fixed	0.00; #,##0.00 0; #,##0	Fixed Integer	F F
Scientific	0.00E+00	Exponent	E
Currency , (comma)	\$#,##0_);...	\$ (dollar)	DOLLAR COMMA
General +/-	General	General * (bargraph)	F F
Percent	0%; 0.00%	Percent	PCT
Date	m/d/yy;d-mmm-yy...		DATE
Time	h:mm; h:mm:ss...		TIME
Text/Literal			F
Label		Alpha	String

- If a string is encountered in a column with numeric format, it is converted to the system-missing value in the active dataset.
- If a numeric value is encountered in a column with string format, it is converted to a blank in the active dataset.
- Blank lines are translated as cases containing the system-missing value for numeric variables and blanks for string variables.
- 1-2-3 and Symphony date and time indicators (shown at the bottom of the screen) are not transferred from *WKS*, *WK1*, or *SYM* files.



## Databases

Database tables are logically very similar to IBM® SPSS® Statistics data files.

- By default, all fields and records from dBASE II, dBASE III, or dBASE IV files are included in the active dataset.
- Field names are automatically translated into variable names. If the `FIELDNAMES` subcommand is used with database files, it is ignored.
- Field names are converted to valid SPSS Statistics variable names.
- Colons used in dBASE II field names are translated to underscores.
- Records in dBASE II, dBASE III, or dBASE IV that have been marked for deletion but that have not actually been purged are included in the active dataset. To differentiate these cases, `GET TRANSLATE` creates a new string variable, `D_R`, which contains an asterisk for cases marked for deletion. Other cases contain a blank for `D_R`.
- Character, floating, and numeric fields are transferred directly to variables. Logical fields are converted into string variables. Memo fields are ignored.

dBASE formats are translated as follows:

<b>dBASE</b>	<b>SPSS Statistics</b>
Character	String
Logical	String
Date	Date
Numeric	Number
Floating	Number
Memo	Ignored

## Tab-Delimited ASCII Files

Tab-delimited ASCII files are simple spreadsheets produced by a text editor, with the columns delimited by tabs and rows, by carriage returns. The first row is usually occupied by column headings.

- By default all columns of all rows are treated as data. Default variable names `VARI`, `VAR2`, and so on are assigned to each column. The data type (numeric or string) for each variable is determined by the first data value in the column.
- If `FIELDNAMES` is specified, the program reads in the first row as variable names and determines the data type by the values read in from the second row.
- Any value that contains non-numeric characters is considered a string value. Dollar and date formats are not recognized and are treated as strings. When string values are encountered for a numeric variable, they are converted to the system-missing value.
- For numeric variables, the assigned format is `F8.2` or the format of the first data value in the column, whichever is wider. Values that exceed the defined width are rounded for display, but the entire value is stored internally.

- For string variables, the assigned format is A8 or the format of the first data value in the column, whichever is wider. Values that exceed the defined width are truncated.
- ASCII data files delimited by space (instead of tabs) or in fixed format should be read by DATA LIST.

## **FILE Subcommand**

FILE names the file to read. The only specification is the name of the file.

- File specifications should be enclosed in quotation marks or apostrophes.

### **Example**

```
GET TRANSLATE FILE='PROJECT.WKS'.
```

- GET TRANSLATE creates an active dataset from the 1-2-3 Release 1.0 spreadsheet with the name *PROJECT.WKS*.
- The active dataset contains all rows and columns and uses the column letters as variable names.
- The format for each variable is determined by the format of the value in the first row of each column.

## **TYPE Subcommand**

TYPE indicates the format of the file.

- TYPE can be omitted if the file extension named on FILE is the default for the type of file that you are reading.
- The TYPE subcommand takes precedence over the file extension.
- You can create a Lotus format file in Multiplan and translate it to an active dataset by specifying WKS on TYPE.

<b>WK</b>	<i>Any Lotus 1-2-3 or Symphony file.</i>
<b>WK1</b>	<i>1-2-3 Release 2.0.</i>
<b>WKS</b>	<i>1-2-3 Release 1A.</i>
<b>SYM</b>	<i>Symphony Release 2.0 or Symphony Release 1.0.</i>
<b>SLK</b>	<i>Microsoft Excel and Multiplan saved in SYLK (symbolic link) format.</i>
<b>XLS</b>	<i>Microsoft Excel. For Excel 5 or later, use GET DATA.</i>
<b>DBF</b>	<i>All dBASE files.</i>
<b>TAB</b>	<i>Tab-delimited ASCII data file.</i>

### **Example**

```
GET TRANSLATE FILE='PROJECT.OCT' /TYPE=SLK.
```

- GET TRANSLATE creates an active dataset from the Multiplan file *PROJECT.OCT*.

## **FIELDNAMES Subcommand**

FIELDNAMES translates spreadsheet field names into variable names.

- FIELDNAMES can be used with spreadsheet and tab-delimited ASCII files only. FIELDNAMES is ignored when used with database files.
- Each cell in the first row of the spreadsheet file (or the specified range) must contain a field name. If a column does not contain a name, the column is dropped.
- Field names are converted to valid IBM® SPSS® Statistics variable names.
- If two or more columns in the spreadsheet have the same field name, digits are appended to all field names after the first, making them unique.
- Illegal characters in field names are changed to underscores in this program.
- If the spreadsheet file uses reserved words (ALL, AND, BY, EQ, GE, GT, LE, LT, NE, NOT, OR, TO, or WITH) as field names, GET TRANSLATE appends a dollar sign (\$) to the variable name. For example, columns named *GE*, *GT*, *EQ*, and *BY* will be renamed *GE\$*, *GT\$*, *EQ\$*, and *BY\$* in the active dataset.

### **Example**

```
GET TRANSLATE FILE='MONTHLY.SYM' /FIELDNAMES.
```

- GET TRANSLATE creates a active dataset from a Symphony 1.0 spreadsheet. The first row in the spreadsheet contains field names that are used as variable names in the active dataset.

## **RANGE Subcommand**

RANGE translates a specified set of cells from a spreadsheet file.

- RANGE cannot be used for translating database files.
- For 1-2-3 or Symphony, specify the beginning of the range with a column letter and row number followed by two periods and the end of the range with a column letter and row number, as in *A1..K14*.
- For Multiplan spreadsheets, specify the beginning and ending cells of the range separated by a colon, as in *R1C1:R14C11*.
- For Excel files, specify the beginning column letter and row number, a colon, and the ending column letter and row number, as in *A1:K14*.
- You can also specify the range using range names supplied in Symphony, 1-2-3, or Multiplan.
- If you specify FIELDNAMES with RANGE, the first row of the range must contain field names.

### **Example**

```
GET TRANSLATE FILE='PROJECT.WKS' /FIELDNAMES /RANGE=D3..J279.
```

- GET TRANSLATE creates an active dataset from the 1-2-3 Release 1A file *PROJECT.WKS*.
- The field names in the first row of the range (row 3) are used as variable names.
- Data from cells D4 through J279 are transferred to the active dataset.

## ***DROP and KEEP Subcommands***

DROP and KEEP are used to copy a subset of variables into the active dataset. DROP specifies the variables not to copy into the active dataset. KEEP specifies the variables to copy. Variables not specified on KEEP are dropped.

- DROP and KEEP cannot precede the FILE or TYPE subcommands.
- DROP and KEEP specifications use variable names. By default, this program uses the column letters from spreadsheets and the field names from databases as variable names.
- If FIELDNAMES is specified when translating from a spreadsheet, the DROP and KEEP subcommands must refer to the field names, not the default column letters.
- Variables can be specified in any order. Neither DROP nor KEEP affects the order of variables in the resulting file. Variables are kept in their original order.
- If a variable is referred to twice on the same subcommand, only the first mention of the variable is recognized.
- Multiple DROP and KEEP subcommands are allowed; the effect is cumulative. Specifying a variable named on a previous DROP or not named on a previous KEEP results in an error and the command is not executed.
- If you specify both RANGE and KEEP, the resulting file contains only variables that are both within the range and specified on KEEP.
- If you specify both RANGE and DROP, the resulting file contains only variables within the range and excludes those mentioned on DROP, even if they are within the range.

### ***Example***

```
GET TRANSLATE FILE='ADDRESS.DBF' /DROP=PHONENO, ENTRY.
```

- GET TRANSLATE creates an active dataset from the dBASE file *ADDRESS.DBF*, omitting the fields named *PHONENO* and *ENTRY*.

### ***Example***

```
GET TRANSLATE FILE='PROJECT.OCT' /TYPE=WK1 /FIELDNAMES
/KEEP=NETINC, REP, QUANTITY, REGION, MONTH, DAY, YEAR.
```

- GET TRANSLATE creates a active dataset from the 1-2-3 Release 2.0 file called *PROJECT.OCT*.
- The subcommand FIELDNAMES indicates that the first row of the spreadsheet contains field names, which will be translated into variable names in the active dataset.
- The subcommand KEEP translates columns with the field names *NETINC*, *REP*, *QUANTITY*, *REGION*, *MONTH*, *DAY*, and *YEAR* to the active dataset.

## ***MAP Subcommand***

MAP displays a list of the variables in the active dataset and their corresponding names in the other application.

- The only specification is the keyword MAP. There are no additional specifications.

- Multiple `MAP` subcommands are allowed. Each `MAP` subcommand maps the results of subcommands that precede it; results of subcommands that follow it are not mapped.

**Example**

```
GET TRANSLATE FILE='ADDRESS.DBF' /DROP=PHONENO, ENTRY /MAP.
```

- `MAP` is specified to confirm that the variables *PHONENO* and *ENTRY* have been dropped.

# GGRAPH

*Note:* Square brackets used in the GGRAPH syntax chart are required parts of the syntax and are not used to indicate optional elements. Any equals signs (=) displayed in the syntax chart are required. The GRAPHSPEC subcommand is required.

```
GGRAPH
  /GRAPHDATASET NAME="name"
    DATASET=datasetname
    VARIABLES=variablespec
    TRANSFORM={NO**
               {VARSTOCASES (SUMMARY="varname" INDEX="varname") }}
    MISSING={LISTWISE** } REPORTMISSING={NO**
        {VARIABLEWISE } {YES }}
    CASELIMIT={100000**
               {value }}
  /GRAPHSPEC SOURCE={INLINE
                   {GPLFILE("filespec")}
                   {VIZMLFILE("filespec")}
                   {VIZTEMPLATE(NAME={"template name"}[LOCATION={LOCAL**
                   {"repository folderspec"}
                   {"filespec"} ]
                   {FILE
                   MAPPING(templatemapping))
                   }}}
    EDITABLE={YES**
              {NO }}
    LABEL="string"
    DEFAULTTEMPLATE={YES**} TEMPLATE=["filespec" ...]
                    {NO }
    VIZSTYLESHEET="stylesheet name" [LOCATION={LOCAL**
                                         {"repository folderspec"}
                                         }]
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

### Release 14.0

- Command introduced.

### Release 15.0

- RENAME syntax qualifier deprecated.
- COUNTCI, MEDIANCI, MEANCI, MEANSD, and MEANSE functions introduced.

### Release 17.0

- Added SOURCE=VIZTEMPLATE to support visualization templates.
- Added VIZSTYLESHEET keyword to support visualization stylesheets.

#### Release 19.0

- Added LOCATION=FILE to support visualization templates stored in an arbitrary location on the file system.

### Examples

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl").
```

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: count=col(source(s), name("COUNT"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Count"))
  ELEMENT: interval(position(jobcat*count))
END GPL.
```

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat
  /GRAPHSPEC SOURCE=VIZTEMPLATE(NAME="Bar"
                                MAPPING("Categories"="jobcat" "Summary"="Count"))
  DEFAULTTEMPLATE=NO.
```

## Overview

GGRAPH generates a graph by computing statistics from variables in a data source and constructing the graph according to the graph specification, which may be written in the Graphics Productions Language (GPL) or ViZml.

### Basic Specification

The basic specification is the GRAPHSPEC subcommand.

### Syntax Rules

- Subcommands and keywords can appear in any order.
- Subcommand names and keywords must be spelled out in full.
- The GRAPHDATASET and GRAPHSPEC subcommands are repeatable.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.

- Strings in the GPL are enclosed in quotation marks. You cannot use single quotes (apostrophes).
- With the IBM® SPSS® Statistics Batch Facility (available only with SPSS Statistics Server), use the `-i` switch when submitting command files that contain `BEGIN GPL-END GPL` blocks.

## **GRAPHDATASET Subcommand**

`GRAPHDATASET` creates graph datasets based on open IBM® SPSS® Statistics data files. The subcommand is repeatable, allowing you to create multiple graph datasets that can be referenced in a graph specification. Furthermore, multiple graph specifications (the ViZml or GPL code that defines a graph) can reference the same graph dataset.

Graph datasets contain the data that accompany a graph. The actual variables and statistics in the graph dataset are specified by the `VARIABLES` keyword.

### **Example**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl").
```

## **NAME Keyword**

The `NAME` keyword specifies the name that identifies the graph dataset when it is referenced in a graph specification. There is no default name, so you must specify one. You can choose any name that honors variable naming rules. (For more information about naming rules, see [Variable Names on p. 49.](#)) When the same graph dataset name is used in multiple `GRAPHDATASET` subcommands, the name in the last `GRAPHDATASET` subcommand is honored.

## **DATASET Keyword**

The `DATASET` keyword specifies the dataset name of an open data file to use for the graph dataset. If the keyword is omitted, `GGRAPH` uses the active dataset. You can also use an asterisk (\*) to refer to the active dataset.

The following are honored only for the active dataset (which cannot be named except with an asterisk):

- `FILTER`
- `USE`
- `SPLIT FILE`
- Weight filtering (exclusion of cases with non-positive weights)
- Temporary transformations
- Pending transformations

### **Example**

```
GGRAPH
```



```
/GRAPHDATASET NAME="graphdataset" DATASET=DataSet2 VARIABLES=jobcat COUNT()
/GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl").
```

## VARIABLES Keyword

The `VARIABLES` keyword identifies the variables, statistics, and utility function results that are included in the graph dataset. These are collectively identified as a **variable specification**. The minimum variable specification is a variable. An aggregation or summary function is required when the variable specification includes a multiple-response set. The order of the variables and functions in the variable specification does not matter. Multiple aggregation or summary functions are allowed so that you can graph more than one statistic. You can also use the `ALL` and `TO` keywords to include multiple variables without explicitly listing them. For information about the `ALL` keyword, see [Keyword ALL on p. 51](#). For information about the `TO` keyword, see [Keyword TO on p. 50](#).

When the variable specification includes an aggregation function and does *not* include the `CASEVALUE` function, the graph dataset is aggregated. Any stand-alone variables in the variable specification act as categorical break variables for the aggregation (including scale variables that are not parameters of a summary function). The function is evaluated for each unique value in each break variable. When the variable specification includes only variables or includes the `CASEVALUE` function, the graph dataset is unaggregated. The built-in variable `$CASENUM` is included in the unaggregated dataset. `$CASENUM` cannot be specified or renamed in the variable specification, but you can refer to it in the graph specification.

An unaggregated graph dataset includes a case for every case in the IBM® SPSS® Statistics dataset. An aggregated dataset includes a case for every combination of unique break variable values. For example, assume that there are two categorical variables that act as break variables. If there are three categories in one variable and two in the other, there are six cases in the aggregated graph dataset, as long as there are values for each category.

*Note:* If the dataset is aggregated, be sure to include all of the break variables in the graph specification (the ViZml or GPL). For example, if the variable specification includes two categorical variables and a summary function of a scale variable, the graph specification should use one of the categorical variables as the *x*-axis variable and one as a grouping or panel variable. Otherwise, the resulting graph will not be correct because it does not contain all of the information used for the aggregation.

### Example

```
GGRAPH
/GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat[NAME="empcat" LEVEL=NOMINAL] COUNT()
/GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl").
```

- The `NAME` qualifier renames a variable. For more information, see the topic [Variable and Function Names on p. 868](#).
- The `LEVEL` qualifier specifies a temporary measurement level for a variable. For more information, see the topic [Measurement Level on p. 869](#).

### Variable and Function Names

The variable name that you use in the variable specification is the same as the name defined in the data dictionary. This also the default name for referencing the variable in the graph specification. To use a different name in the graph specification, rename the variable by appending the qualifier `[NAME="name"]` to the name in the variable specification. You might do this to avoid name conflicts across datasets, to shorten the name, or to reuse the same graph specification even if the datasets have different variable names. For example:

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat[NAME="catvar"] COUNT()
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: catvar=col(source(s), name("catvar"), unit.category())
  DATA: count=col(source(s), name("COUNT"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Count"))
  ELEMENT: interval(position(catvar*count))
END GPL.
```

The default name for a summary function is the function name in uppercase letters followed by the parameters separated by underscores. For example, if the function is `MEAN(salary)`, the default name for referencing this function in the graph specification is `MEAN_salary`. For `GPTILE(salary, 90)`, the default name is `GPTILE_salary_90`. You can also change the default function name using the qualifier `[NAME="name"]`, just as you do with variables. For example:

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset"
  VARIABLES=jobcat MEDIAN(salary) MEAN(salary)[NAME="meansal"]
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: medsal=col(source(s), name("MEDIAN_salary"))
  DATA: meansal=col(source(s), name("meansal"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Salary"))
  ELEMENT: line(position(jobcat*medsal), color("Median"))
  ELEMENT: line(position(jobcat*meansal), color("Mean"))
END GPL.
```

Error interval functions produce three values (a summary value, an upper bound, and a lower bound), so there are three default names for these functions. The default name for the summary value follows the same rule as the default name for a summary function: the function name in uppercase letters followed by the parameters separated by underscores. The other two values are this name with `_HIGH` appended to the name for the upper bound and `_LOW` appended to the name for the lower bound. For example, if the function is `MEANCI(salary, 95)`, the default names for referencing the results of this function in the graph specification are `MEANCI_salary_95`,

MEANCI\_salary\_95\_HIGH, and MEANCI\_salary\_95\_LOW. You can change the names of the values using the qualifiers [NAME="name" HIGH="name" LOW="name"]. For example:

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset"
    VARIABLES=jobcat COUNTCI(95)[NAME="stat" HIGH="high" LOW="low"]
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: stat=col(source(s), name("stat"))
  DATA: high=col(source(s), name("high"))
  DATA: low=col(source(s), name("low"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Count with 95% CI"))
  ELEMENT: point(position(jobcat*stat))
  ELEMENT: interval(position(region.spread.range(jobcat*(low+high))), shape(shape.ibeam))
END GPL.
```

## Measurement Level

You can change a variable's measurement level temporarily by appending the qualifier [LEVEL=measurement level] to the name in the variable specification. (The variable's measurement level in the dictionary is unaffected.) Valid values for the measurement level are SCALE, NOMINAL, and ORDINAL. Currently, the measurement level qualifier is used to influence the behavior of the REPORTMISSING keyword. If the measurement level is set to SCALE, missing values are not reported for that variable, even if the value of the REPORTMISSING keyword is YES. If you are using the NAME qualifier for the same variable, both qualifiers are enclosed in the same pair of square brackets. For example:

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat [NAME="empcat" LEVEL=NOMINAL] COUNT()
  /GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl").
```

## Functions

*Utility functions:*

**CASE-VALUE(var)**      *Yields the value of the specified variable for each case.* CASEVALUE always produces one value for each case and always results in GGRAPH creating an unaggregated graph dataset. Use this function when you are creating graphs of individual cases and want to use the values of the specified variable as the axis tick labels for each case. This function cannot be used with multiple response sets or aggregation functions.

*Aggregation functions:*

Three groups of aggregation functions are available: count functions, summary functions, and error interval functions.

*Count functions:*

*Note:* Percent and cumulative statistic functions are not available in the variable specification. Use the summary percent and cumulative statistic functions that are available in the Graphics Production Language (GPL) itself.

<b>COUNT()</b>	<i>Frequency of cases in each category.</i>
<b>RESPONSES()</b>	<i>Number of responses for a multiple dichotomy set.</i>
<b>RESPONSES(DUP / NODUP)</b>	<i>Number of responses for a multiple category set. The argument (DUP or NODUP) specifies whether the function counts duplicates. The argument is optional, and the default is not to count duplicates. This function cannot be used with a multiple dichotomy set.</i>

- Count functions yield the count of valid cases within categories determined by the other variables in the variable specification (including other scale variables that are not parameters of a summary function).
- Count functions do not use variables as parameters.

*Summary functions:*

<b>MINIMUM(var)</b>	<i>Minimum value of the variable.</i>
<b>MAXIMUM(var)</b>	<i>Maximum value of the variable.</i>
<b>VALIDN(var)</b>	<i>Number of cases for which the variable has a nonmissing value.</i>
<b>SUM(var)</b>	<i>Sum of the values of the variable.</i>
<b>MEAN(var)</b>	<i>Mean of the variable.</i>
<b>STDDEV(var)</b>	<i>Standard deviation of the variable.</i>
<b>VARIANCE(var)</b>	<i>Variance of the variable.</i>
<b>MEDIAN(var)</b>	<i>Median of the variable.</i>
<b>GMEDIAN(var)</b>	<i>Group median of the variable.</i>
<b>MODE(var)</b>	<i>Mode of the variable.</i>
<b>PTILE(var,x)</b>	<i>Xth percentile value of the variable. X must be greater than 0 and less than 100.</i>
<b>GPTILE(var,x)</b>	<i>Xth percentile value of the variable, where the percentile is calculated as if the values were uniformly distributed over the whole interval. X must be greater than 0 and less than 100.</i>
<b>PLT(var,x)</b>	<i>Percentage of cases for which the value of the variable is less than x.</i>
<b>PGT(var,x)</b>	<i>Percentage of cases for which the value of the variable is greater than x.</i>
<b>NLT(var,x)</b>	<i>Number of cases for which the value of the variable is less than x.</i>
<b>NGT(var,x)</b>	<i>Number of cases for which the value of the variable is greater than x.</i>
<b>PIN(var,x1,x2)</b>	<i>Percentage of cases for which the value of the variable is greater than or equal to x1 and less than or equal to x2. x1 cannot exceed x2.</i>
<b>NIN(var,x1,x2)</b>	<i>Number of cases for which the value of the variable is greater than or equal to x1 and less than or equal to x2. x1 cannot exceed x2.</i>
<b>NLE(var,x)</b>	<i>Number of cases for which the value of the variable is less than or equal to x.</i>
<b>PLE(var,x)</b>	<i>Percentage of cases for which the value of the variable is less than or equal to x.</i>
<b>NEQ(var,x)</b>	<i>Number of cases for which the value of the variable is equal to x.</i>
<b>PEQ(var,x)</b>	<i>Percentage of cases for which the value of the variable is equal to x.</i>

<b>NGE(var,x)</b>	<i>Number of cases for which the value of the variable is greater than or equal to x.</i>
<b>PGE(var,x)</b>	<i>Percentage of cases for which the value of the variable is greater than or equal to x.</i>

- Summary functions yield a single value.
- Summary functions operate on summary variables (variables that record continuous values, such as age or expenses). To use a summary function, specify the name of one or more variables as the first parameter of the function and then specify other required parameters as shown. The variable used as a parameter cannot contain string data.

*Error interval functions:*

<b>COUNTCI(alpha)</b>	<i>Confidence intervals for the count with a confidence level of alpha. alpha must be greater than or equal to 50 and less than 100.</i>
<b>MEDIANCI(var,alpha)</b>	<i>Confidence intervals for median of the variable with a confidence level of alpha. alpha must be greater than or equal to 50 and less than 100.</i>
<b>MEANCI(var,alpha)</b>	<i>Confidence intervals for mean of the variable with a confidence level of alpha. alpha must be greater than or equal to 50 and less than 100.</i>
<b>MEANSD(var,multiplier)</b>	<i>Standard deviations for mean of the variable with a multiplier. multiplier must be an integer greater than 0.</i>
<b>MEANSE(var,multiplier)</b>	<i>Standard errors for mean of the variable with a multiplier. multiplier must be an integer greater than 0.</i>

- Error functions yield three values: a summary value, a lower bound value, and an upper bound value.
- Error functions may or may not operate on summary variables (variables that record continuous values, such as age or expenses). To use a summary function that operates on a variable, specify the name of the variable as the first parameter of the function and then specify other required parameters as shown. The variable used as a parameter cannot contain string data.

## ***TRANSFORM Keyword***

The TRANSFORM keyword applies a transformation to the graph dataset.

**NO** *Do not transform the graph dataset.*

**VARSTOCASES(SUMMARY="varname" INDEX="varname")**

*Transform the summary function results to cases in the graph dataset. Use this when you are creating graphs of separate variables. The results of each summary function becomes a case in the graph dataset, and the data elements drawn for each case act like categories in a categorical graph. Each case is identified by an index variable whose value is a unique sequential number. The result of the summary function is stored in the summary variable. The upper and lower bound of error interval functions are also stored in two other variables. By default, the names of the variables are #INDEX for the index variable, #SUMMARY for the summary variable, #HIGH for the upper bound variable, and #LOW for the lower bound variable. You can change these names by using the SUMMARY, INDEX, HIGH, and LOW qualifiers. Furthermore, break variables in the variable*

specification are treated as fixed variables and are not transposed. Note that this transformation is similar to the VARSTOCASES command (see [VARSTOCASES](#) on p. 2136).

### Examples

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=MEAN(salbegin) MEAN(salary)
  TRANSFORM=VARSTOCASES(SUMMARY="meansal" INDEX="variables")
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: variables=col(source(s), name("variables"), unit.category())
  DATA: meansal=col(source(s), name("meansal"))
  GUIDE: axis(dim(2), label("Mean"))
  ELEMENT: interval(position(variables*meansal))
END GPL.
```

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=MEANCI(salbegin, 95) MEANCI(salary, 95)
  TRANSFORM=VARSTOCASES(SUMMARY="meansal" INDEX="variables" LOW="low" HIGH="high")
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: variables=col(source(s), name("variables"), unit.category())
  DATA: meansal=col(source(s), name("meansal"))
  DATA: low=col(source(s), name("low"))
  DATA: high=col(source(s), name("high"))
  GUIDE: axis(dim(2), label("Mean with 95% CI"))
  ELEMENT: point(position(variables*meansal))
  ELEMENT: interval(position(region.spread.range(variables*(low+high))),
                    shape(shape.ibeam))
END GPL.
```

## MISSING Keyword

The `MISSING` keyword specifies how missing values are handled when the variable specification includes an aggregation function. When the variable specification includes only variables or includes the `CASEVALUE` function, this keyword does not affect the treatment of missing values. The graph dataset is unaggregated, so cases with system- and user-missing values are always included in the graph dataset.

- |                     |  |
|---------------------|--|
| <b>LISTWISE</b>     | <i>Exclude the whole case if any one of the variables in the variable specification has a missing value. This is the default.</i>  |
| <b>VARIABLEWISE</b> | <i>Exclude a case from the aggregation function if the value is missing for a particular variable being analyzed. This means that a case is excluded if that case has a missing value for a variable that is a summary function parameter.</i> |

## **REPORTMISSING Keyword**

The REPORTMISSING keyword specifies whether to create a category for each unique user-missing value.

- NO** *Do not create a category for each unique user-missing value. User-missing values are treated like system-missing values. This is the default.*
- YES** *Create a category for each unique user-missing value. User-missing values are treated as valid categories, are included as break variables for aggregation functions, and are drawn in the graph. Note that this does not affect variables identified as SCALE by the LEVEL qualifier in the VARIABLES keyword.*

## **CASELIMIT Keyword**

The CASELIMIT keyword specifies a limit to the number of cases that are included in the graph dataset. The limit does not apply to the number of cases use for analysis in any functions specified by the VARIABLES keyword. It only limits the number of cases in the graph dataset, which may or may not affect the number cases drawn in the resulting chart. You may want to limit the number of cases for very large datasets that are not summarized by a function. A scatterplot is an example. Limiting cases may improve performance.

- value** *Limit the number of cases in the graph dataset to the specified value. The default value is 1000000.*

## **GRAPHSPEC Subcommand**

GRAPHSPEC defines a graph specification. A graph specification identifies the source used to create the graph, in addition to other features like templates. GRAPHSPEC is repeatable, allowing you to define multiple graph specifications to create multiple graphs with one GGRAPH command.

### **Example**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl").
```

## **SOURCE Keyword**

The SOURCE keyword specifies the source of the graph specification.

- INLINE** *An inline graph specification follows the GGRAPH command. Currently, the BEGIN GPL/END GPL block is used for the inline graph specification. This block must follow the GGRAPH command, and there must be as many blocks as there are GRAPHSPEC subcommands with SOURCE=INLINE. For more information, see the topic [BEGIN GPL-END GPL on p. 229](#). See [Overview on p. 229](#) for limitations.*
- GPLFILE("filespec")**

Use the specified GPL file as the graph specification. See the *GPL Reference Guide* on the manuals CD for more details about GPL. The examples in the GPL documentation may look different compared to the syntax pasted from the Chart Builder. The main difference is in when aggregation occurs. See [Working with the GPL](#) below for information about the differences. See [GPL Examples on p. 879](#) for examples with GPL that is similar to the pasted syntax.

**VIZMLFILE("filespec")**

Use the specified ViZml file as the graph specification. You can save ViZml from the Chart Editor.

**VIZTEMPLATE(NAME="template name"[LOCATION=LOCAL]  
MAPPING(templatemapping))**

**VIZTEMPLATE(NAME="template name"[LOCATION="repository folderspec"]  
MAPPING(templatemapping))**

**VIZTEMPLATE(NAME="filespec"[LOCATION=FILE]  
MAPPING(templatemapping))**

Use the specified visualization template and mapping as the graph specification. The LOCATION qualifier specifies the location in which to find the named template. You specify whether the template is a local imported template, is obtained from a specific IBM® SPSS® Collaboration and Deployment Services Repository, or is stored in an arbitrary location on the file system. Note that a visualization template is not the same as a template for use with the TEMPLATE keyword.

- If you received a template and want to reference it in syntax, you must first manually import it with the Graphboard Template Chooser to make it a local template.
- For built-in or imported templates, set the location to LOCAL.
- For templates stored in an IBM SPSS Collaboration and Deployment Services Repository, use an IBM SPSS Collaboration and Deployment Services Repository file specification. For more information about IBM SPSS Collaboration and Deployment Services Repository file specifications, see [File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects](#) on p. 2230.
- For templates stored on the file system, set the location to FILE and specify the full path to the visualization template file with NAME.

The template mapping is a series of key-value pairs separated by spaces. An example is `MAPPING("Categories"="jobcat" "Summary"="Count")`.

- The key is a slot defined by the template, and the value is a variable or string assigned to the slot.
- If a variable is renamed in the VARIABLES keyword using the [NAME="name"] qualifier, the value refers to the rename, not the original name.
- In the example, the keys are Categories and Summary, and the respective values are jobcat and Count.
- Individual keys and values are enclosed in quotes.
- Each value can also be followed by the qualifier [DATASET="dataset name"] to indicate a variable's dataset.
- At this time, only one dataset can be used with each template.
- To determine valid slot names, open the Graphboard Template Chooser, add variables to slots of interest, and paste the syntax. You can access the Template Chooser by selecting Graphboard Template Chooser from the Graphs menu. Visualization templates on the file system are not available in the Graphboard Template Chooser. You will have to import them first, or you will have to open them in a text editor to determine the slot names.

New templates can be created with the separate IBM® SPSS® Visualization Designer product. The GGRAPH command can use these templates if they are stored in the IBM SPSS Collaboration and Deployment Services Repository or if they are imported using the Graphboard Template Chooser.



*Note:* When using a visualization template, it is recommended that you include `DEFAULTTEMPLATE=NO`. This setting will prevent the default template from overriding the styles in the visualization template.

### Examples

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: count=col(source(s), name("COUNT"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Count"))
  ELEMENT: interval(position(jobcat*count))
END GPL.

GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl").

GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=VIZMLFILE("simplebarchart.xml").

GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat
  /GRAPHSPEC SOURCE=VIZTEMPLATE(NAME="Bar"
    MAPPING("Categories"="jobcat" "Summary"="Count"))
  DEFAULTTEMPLATE=NO.

GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=salbegin salary gender[NAME="sex"]
  /GRAPHSPEC SOURCE=VIZTEMPLATE(NAME="Binned Scatter" [LOCATION="SPSSCR:///viz/templates"]
    MAPPING("X"="salbegin" [DATASET="graphdataset"]
      "Y"="salary" [DATASET="graphdataset"]
      "Panel across"="sex" [DATASET="graphdataset"]))
  DEFAULTTEMPLATE=NO.

GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=salbegin salary gender[NAME="sex"]
  /GRAPHSPEC SOURCE=VIZTEMPLATE(NAME="/vizfiles/MyBinnedScatter.viztemplate" [LOCATION=FILE]
    MAPPING("X"="salbegin" [DATASET="graphdataset"]
      "Y"="salary" [DATASET="graphdataset"]
      "Panel across"="sex" [DATASET="graphdataset"]))
  DEFAULTTEMPLATE=NO.
```

### Working with the GPL

The Chart Builder allows you to paste GGRAPH syntax. This syntax contains inline GPL

You may want to edit the GPL to create a chart or add a feature that isn't available from the Chart Builder. You can use the GPL documentation to help you. However, the GPL documentation always uses unaggregated data and includes GPL statistics in the examples to aggregate the data. The pasted syntax, on the other hand, may use data aggregated by a GGRAPH summary function. Also, the pasted syntax includes defaults that you may have to change when you edit

the syntax. Therefore, it may be confusing how you can use the pasted syntax to create the examples. Following are some tips.

- Variables must be specified in two places: in the `VARIABLES` keyword in the `GGRAPH` command and in the `DATA` statements in the `GPL`. So, if you add a variable, make sure a reference to it appears in both places.
- Pasted syntax often uses the `VARIABLES` keyword to specify summary statistics. Like other variables, the summary function name is specified in the `GPL DATA` statement. You do not need to use `GGRAPH` summary functions. Instead, you can use the equivalent `GPL` statistic for aggregation. However, for very large data sets, you may find that pre-aggregating the data with `GGRAPH` is faster than using the aggregation in the `GPL` itself. Try both approaches and stick with the one that feels comfortable to you. In the examples that follow, you can compare the different approaches.
- Make sure that you understand how the functions are being used in the `GPL`. You may need to modify one or more of them when you add a variable to pasted syntax. For example, if you change the dimension on which a categorical variable appears, you may need to change references to the dimension in the `GUIDE` and `SCALE` statements. If you are unsure about whether you need a particular function, try removing it and see if you get the results you expect.

Here's an example from the `GPL` documentation:

**Figure 103-1**

*Example from GPL documentation*

```
SOURCE: s=usersource(id("Employeeedata"))
DATA: jobcat = col(source(s), name("jobcat"), unit.category())
DATA: gender = col(source(s), name("gender"), unit.category())
DATA: salary = col(source(s), name("salary"))
SCALE: linear(dim(2), include(0))
GUIDE: axis(dim(3), label("Gender"))
GUIDE: axis(dim(2), label("Mean Salary"))
GUIDE: axis(dim(1), label("Job Category"))
ELEMENT: interval(position(summary.mean(jobcat*salary*gender)))
```

The simplest way to use the example is to use unaggregated data and `VARIABLES=ALL` like this:

**Figure 103-2**

*Modified example with unaggregated data*

```
GGRAPH
  /GRAPHDATASET NAME="Employeeedata" VARIABLES=ALL
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
SOURCE: s=usersource(id("Employeeedata"))
DATA: jobcat = col(source(s), name("jobcat"), unit.category())
DATA: gender = col(source(s), name("gender"), unit.category())
DATA: salary = col(source(s), name("salary"))
SCALE: linear(dim(2), include(0))
GUIDE: axis(dim(3), label("Gender"))
GUIDE: axis(dim(2), label("Mean Salary"))
GUIDE: axis(dim(1), label("Job Category"))
ELEMENT: interval(position(summary.mean(jobcat*salary*gender)))END GPL
```

Note that specifying `VARIABLES=ALL` includes all the data in the graph. You can improve performance by using only those variables that you need. In this example, `VARIABLES=jobcat gender salary` would have been sufficient.

You can also use aggregated data like the following, which is more similar to the pasted syntax:

**Figure 103-3**

*Modified example with aggregated data*

```
GGRAPH
  /GRAPHDATASET NAME="Employeeedata" VARIABLES=jobcat gender MEAN(salary)
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("Employeeedata"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: gender=col(source(s), name("gender"), unit.category())
  DATA: MEAN_salary=col(source(s), name("MEAN_salary"))
  SCALE: linear(dim(2), include(0))
  GUIDE: axis(dim(3), label("Gender"))
  GUIDE: axis(dim(2), label("Mean Salary"))
  GUIDE: axis(dim(1), label("Job Category"))
  ELEMENT: interval(position(jobcat*MEAN_salary*gender))
END GPL.
```

## **EDITABLE Keyword**

The EDITABLE keyword specifies that the resulting graph can be edited. If you are creating a complicated graph with the graph specification, it may be useful to prevent editing because not all of the graph's features may be supported in the editor.

**YES**            *The graph can be edited.* This is the default.  
**NO**             *The graph cannot be edited.*

## **LABEL Keyword**

The LABEL keyword specifies the output label. This label appears in the Output Viewer. It is also used in Output XML (OXML) as a chartTitle element, which is not the same as the title in the graph itself.

**string**            *Use the specified string as the label.*

## **DEFAULTTEMPLATE Keyword**

The DEFAULTTEMPLATE keyword specifies whether GGRAPH applies the default styles to the graph. Most default styles are defined in the Options dialog box, which you can access by choosing Options from the Edit menu. Then click the Charts tab. Some SET commands also define default aesthetics. Finally, other default styles are set to improve the presentation of graphs. These are controlled by the *chart\_style.sgt* template file located in the installation directory.

**YES**            *Apply default styles to the graph.* This is the default.  
**NO**             *Do not apply default styles to the graph.* This option is useful when you are using a custom ViZml or GPL file that defines styles that you do not want to be overridden by the default styles.

## TEMPLATE Keyword

The `TEMPLATE` keyword identifies an existing template file or files and applies them to the graph requested by the current `GGRAPH` command. The template overrides the default settings that are used to create any graph, and the specifications on the current `GGRAPH` command override the template. Templates do *not* override visualization stylesheets specified with the `VIZSTYLESHEET` keyword.

Templates are created in the Chart Editor by saving an existing chart as a template. Note that these templates are not the same as visualization templates. Visualization templates are created in a separate IBM® SPSS® Visualization Designer product and are referenced by the `SOURCE` keyword.

The `TEMPLATE` keyword is followed by an equals sign (=) and square brackets ( [ ] ) that contain one or more file specifications. Each file specification is enclosed in quotation marks. The square brackets are optional if there is only one file, but the file must be enclosed in quotation marks. Note that the order in which the template files are specified is the order in which `GGRAPH` applies the templates. Therefore, template files that appear after other template files can override the templates that were applied earlier.

**filespec**                    *Apply the specified template file or files to the graph being created.*

### Example

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=GPLFILE("simplebarchart.gpl")
  TEMPLATE=["mytemplate.sgt"
           "/myothertemplate.sgt"].
```

## VIZSTYLESHEET Keyword

The `VIZSTYLESHEET` keyword identifies an existing visualization stylesheet and applies it to the graph requested by the current `GGRAPH` command. Styles stored in the original visualization template, GPL specification, ViZml specification, or user-specified templates are preserved. Because the default template may set styles that cannot be replaced by the visualization stylesheet, it is recommended that you use `DEFAULTTEMPLATE=NO` in combination with the `VIZSTYLESHEET` keyword.

The keyword is followed by an equals sign (=) and the name of the stylesheet in quotation marks. By default, `GGRAPH` looks for a built-in or imported template that is stored on the local machine. This default behavior is equivalent to appending the qualifier `[LOCATION=LOCAL]` to the quoted stylesheet name. To use a stylesheet in an IBM® SPSS® Collaboration and Deployment Services Repository, append the qualifier `[LOCATION="repository folderspec"]` to the quoted stylesheet name. For more information about IBM SPSS Collaboration and Deployment Services Repository file specifications, see [File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects](#) on p. 2230. The IBM SPSS Collaboration and Deployment Services Repository file specification is enclosed in quotation marks.

New stylesheets can be created with the separate IBM® SPSS® Visualization Designer product. The `GGRAPH` command can reference these stylesheets if they are stored in the IBM SPSS Collaboration and Deployment Services Repository or if they are imported using the

Graphboard Template Chooser. You can access the Template Chooser by selecting Graphboard Template Chooser from the Graphs menu.

**“stylesheet name”**

*Apply the specified stylesheet to the graph being created.*

**Example**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat
  /GRAPHSPEC SOURCE=VIZTEMPLATE(NAME="Binned Scatter"[LOCATION=LOCAL]
                                MAPPING("X"="salbegin" "Y"="salary"
                                          "Panel across"="gender"))
  VIZSTYLE SHEET="Corporate" [LOCATION="SPSSCR://viz/stylesheet"]
  DEFAULTTEMPLATE=NO.
```

## GPL Examples

Following are some graph examples using GPL. Pictures are not included to encourage you to run the examples. Except when noted, all examples use *Employee data.sav*, which is located in the product installation directory. See the *GPL Reference Guide* on the manuals CD for more examples.

**Simple Bar Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat MEAN(salary)
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: meansal=col(source(s), name("MEAN_salary"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Mean Current Salary"))
  ELEMENT: interval(position(jobcat*meansal))
END GPL.
```

**Simple Bar Chart Using a Multiple-Response Set**

*Note:* This example uses *1991 U.S. General Social Survey.sav*, which is located in the product installation directory.

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=$prob RESPONSES() [NAME="RESPONSES"]
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: prob=col(source(s), name("$prob"), unit.category())
  DATA: responses=col(source(s), name("RESPONSES"))
  GUIDE: axis(dim(1), label("Most Important Problems in Last 12 Months"))
  GUIDE: axis(dim(2), label("Responses"))
  ELEMENT: interval(position(prob*responses))
END GPL.
```

**Stacked Bar Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat gender COUNT()
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: gender=col(source(s), name("gender"), unit.category())
  DATA: count=col(source(s), name("COUNT"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Count"))
  ELEMENT: interval.stack(position(jobcat*count), color(gender))
END GPL.
```

**Clustered Bar Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat gender MEAN(salary)
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: gender=col(source(s), name("gender"), unit.category())
  DATA: meansal=col(source(s), name("MEAN_salary"))
  COORD: rect(dim(1,2), cluster(3,0))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Mean Current Salary"))
  ELEMENT: interval(position(gender*meansal*jobcat), color(gender))
END GPL.
```

**Paneled Bar Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat gender MEAN(salary)
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: gender=col(source(s), name("gender"), unit.category())
  DATA: meansal=col(source(s), name("MEAN_salary"))
  GUIDE: axis(dim(1), label("Gender"))
  GUIDE: axis(dim(2), label("Mean Current Salary"))
  GUIDE: axis(dim(3), label("Employment Category"))
  ELEMENT: interval(position(gender*meansal*jobcat))
END GPL.
```

**3-D Bar Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat gender MEAN(salary)
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: gender=col(source(s), name("gender"), unit.category())
  DATA: meansal=col(source(s), name("MEAN_salary"))
  GUIDE: axis(dim(1), label("Gender"))
  GUIDE: axis(dim(2), label("Employment Category"))
  GUIDE: axis(dim(3), label("Mean Current Salary"))
  COORD: rect(dim(1,2,3))
  ELEMENT: interval(position(gender*jobcat*meansal))
END GPL.
```

**Simple Scatterplot**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=salbegin salary
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: salbegin=col(source(s), name("salbegin"))
  DATA: salary=col(source(s), name("salary"))
  GUIDE: axis(dim(1), label("Beginning Salary"))
  GUIDE: axis(dim(2), label("Current Salary"))
  ELEMENT: point(position(salbegin*salary))
END GPL.
```

**Simple Scatterplot with Fit Line**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=salbegin salary
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: salbegin=col(source(s), name("salbegin"))
  DATA: salary=col(source(s), name("salary"))
  GUIDE: axis(dim(1), label("Beginning Salary"))
  GUIDE: axis(dim(2), label("Current Salary"))
  ELEMENT: point(position(salbegin*salary))
  ELEMENT: line(position(smooth.linear(salbegin*salary)))
END GPL.
```

**Pie Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat COUNT()
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE:s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: count=col(source(s), name("COUNT"))
  GUIDE: axis(dim(1), null())
  COORD: polar.theta()
  SCALE: linear(dim(1), dataMinimum(), dataMaximum())
  ELEMENT: interval.stack(position(summary.percent(count)),color(jobcat))
END GPL.
```

**Area Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat MEAN(salary)
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: meansal=col(source(s), name("MEAN_salary"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Mean Current Salary"))
  ELEMENT: area(position(jobcat*meansal))
END GPL.
```

**Grouped Line Chart**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=jobcat gender MEAN(salary)
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: gender=col(source(s), name("gender"), unit.category())
  DATA: meansal=col(source(s), name("MEAN_salary"))
  GUIDE: axis(dim(1), label("Employment Category"))
  GUIDE: axis(dim(2), label("Mean Current Salary"))
  ELEMENT: line(position(jobcat*meansal), color(gender), missing.wings())
END GPL.
```

**Bar Chart of Separate Variables**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=MEAN(salary) MEAN(salbegin)
  TRANSFORM=VARSTOCASES(SUMMARY="meansal" INDEX="variables")
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: variables=col(source(s), name("variables"), unit.category())
  DATA: meansal=col(source(s), name("meansal"))
  GUIDE: axis(dim(2), label("Mean"))
  ELEMENT: interval(position(variables*meansal))
END GPL.
```

**Bar Chart Clustered by Separate Variables**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=MEAN(salary) MEAN(salbegin) jobcat
  TRANSFORM=VARSTOCASES(SUMMARY="meansal" INDEX="variables")
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: variables=col(source(s), name("variables"), unit.category())
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: meansal=col(source(s), name("meansal"))
  COORD: rect(dim(1,2), cluster(3,0))
  GUIDE: axis(dim(2), label("Mean"))
  GUIDE: legend(aesthetic(aesthetic.color), label("Variables"))
  ELEMENT: interval(position(variables*meansal*jobcat), color(variables))
END GPL.
```

**Bar Chart of Separate Variables Clustered by Categorical Variable**

```
GGRAPH
  /GRAPHDATASET NAME="graphdataset" VARIABLES=MEAN(salary) MEAN(salbegin) jobcat
  TRANSFORM=VARSTOCASES(SUMMARY="meansal" INDEX="variables")
  /GRAPHSPEC SOURCE=INLINE.
BEGIN GPL
  SOURCE: s=userSource(id("graphdataset"))
  DATA: variables=col(source(s), name("variables"), unit.category())
  DATA: jobcat=col(source(s), name("jobcat"), unit.category())
  DATA: meansal=col(source(s), name("meansal"))
  COORD: rect(dim(1,2), cluster(3,0))
  GUIDE: axis(dim(2), label("Mean"))
  GUIDE: legend(aesthetic(aesthetic.color), label("Employment Category"))
  ELEMENT: interval(position(jobcat*meansal*variables), color(jobcat))
END GPL.
```



# GLM

GLM is available in the Advanced Statistics option.

GLM dependent varlist [BY factor list [WITH covariate list]]

```
[/WSFACTOR=name levels [{DEVIATION [(refcat)]          }] name...
                        {SIMPLE [(refcat)]             }
                        {DIFFERENCE                    }
                        {HELMERT                       }
                        {REPEATED                      }
                        {POLYNOMIAL [({1,2,3...})]**}
                        {          {metric }           }
                        {SPECIAL (matrix)              }

[/MEASURE=newname newname...]

[/WSDESIGN=effect effect...]+

[/RANDOM=factor factor...]

[/REGWGT=varname]

[/METHOD=SSTYPE({1  })]
                {2  }
                {3**}
                {4  }

[/INTERCEPT=[INCLUDE**] [EXCLUDE]]

[/MISSING=[INCLUDE] [EXCLUDE**]]

[/CRITERIA=[EPS({1E-8**})] [ALPHA({0.05**})]
           {a    }         {a    }

[/PRINT  = [DESCRIPTIVE] [HOMOGENEITY] [PARAMETER] [ETASQ] [RSSCP]
           [GEF] [LOF] [OPOWER] [TEST [({SSCP} [LMATRIX] [MMATRIX])]]]

[/PLOT=[SPREADLEVEL] [RESIDUALS]
       [PROFILE (factor factor*factor factor*factor*factor ...)
        [WITH(covariate={value} [...])]]
       {MEAN }

[/TEST=effect VS {linear combination [DF(df)]}]
           {value DF (df)              }

[/LMATRIX={{["label"] effect list effect list ...;...}}
           {["label"] effect list effect list ...   }
           {["label"] ALL list; ALL...              }
           {["label"] ALL list                      }

[/CONTRAST (factor name)={DEVIATION[(refcat)]** #   }]
                        {SIMPLE [(refcat)]         }
                        {DIFFERENCE                 }
                        {HELMERT                    }
                        {REPEATED                   }
                        {POLYNOMIAL [({1,2,3...})]}
                        {          {metric }        }
                        {SPECIAL (matrix)           }

[/MMATRIX= {["label"] depvar value depvar value ...;["label"]...}
           {["label"] depvar value depvar value ...   }
           {["label"] ALL list; ["label"] ...         }
           {["label"] ALL list                        }

[/KMATRIX= {list of numbers   }]
           {list of numbers;...}

[/POSTHOC = effect [effect...]
```

```

([SNK] [TUKEY] [BTUKEY] [DUNCAN]
 [SCHEFFE] [DUNNETT(refcat)] [DUNNETTL(refcat)]
 [DUNNETTR(refcat)] [BONFERRONI] [LSD] [SIDAK]
 [GT2] [GABRIEL] [FREGW] [QREGW] [T2] [T3] [GH] [C]
 [WALLER ({100** }))]
      {kratio}
[VS effect]

[/EMMEANS=TABLES({OVERALL          }) [WITH(covariate={value} [...])]
      {factor          }                {MEAN }
      {factor*factor... }
      {wsfactor        }
      {wsfactor*wsfactor ... }
      {factor*...wsfactor*...}

      [COMPARE ADJ(LSD) (BONFERRONI) (SIDAK)]

[/SAVE={tempvar [(list of names)] [tempvar [(list of names)]...]}
      [DESIGN]

[/OUTFILE={ {COVB('savfile' | 'dataset')} }
      {CORB('savfile' | 'dataset')} }
      [EFFECT('savfile' | 'dataset')] [DESIGN('savfile' | 'dataset')]

[/DESIGN={ {INTERCEPT... } }
      {[effect effect...]}

```

† WSDSIGN uses the same specification as DESIGN, with only within-subjects factors.

‡ DEVIATION is the default for between-subjects factors, while POLYNOMIAL is the default for within-subjects factors.

\*\* Default if the subcommand or keyword is omitted.

*Temporary variables (tempvar) are:*

PRED, WPRED, RESID, WRESID, DRESID, ZRESID, SRESID, SEPREP, COOK, LEVER

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 17.0

- POSTHOC subcommand: T2, T3, GH, and C keywords are not valid when multiple factors in the model.
- PLOT subcommand: new WITH keyword allows you to fix covariate values for profile plots.

## **Overview**

GLM (general linear model) is a general procedure for analysis of variance and covariance, as well as regression. GLM is the most versatile of the analysis-of-variance procedures and can be used for both univariate and multivariate designs. GLM allows you to:

- Include interaction and nested effects in your design model. Multiple nesting is allowed; for example, *A* within *B* within *C* is specified as *A(B(C))*.

- Include covariates in your design model. GLM also allows covariate-by-covariate and covariate-by-factor interactions, such as  $X$  by  $X$  (or  $X^*X$ ),  $X$  by  $A$  (or  $X^*A$ ), and  $X$  by  $A$  within  $B$  (or  $X^*A(B)$ ). Thus, polynomial regression or a test of the homogeneity of regressions can be performed.
- Select appropriate sums-of-squares hypothesis tests for effects in balanced design models, unbalanced all-cells-filled design models, and some-cells-empty design models. The estimable functions that correspond to the hypothesis test for each effect in the model can also be displayed.
- Display the general form of estimable functions.
- Display expected mean squares, automatically detecting and using the appropriate error term for testing each effect in mixed-effects and random-effects models.
- Select commonly used contrasts or specify custom contrasts to perform hypothesis tests.
- Customize hypothesis testing, based on the null hypothesis  $\mathbf{LBM} = \mathbf{K}$ , where  $\mathbf{B}$  is the parameter vector or matrix.
- Display a variety of post hoc tests for multiple comparisons.
- Display estimates of population marginal cell means for both between-subjects factors and within-subjects factors, adjusted for covariates.
- Perform multivariate analysis of variance and covariance.
- Estimate parameters by using the method of weighted least squares and a generalized inverse technique.
- Graphically compare the levels in a model by displaying plots of estimated marginal cell means for each level of a factor, with separate lines for each level of another factor in the model.
- Display a variety of estimates and measures that are useful for diagnostic checking. All of these estimates and measures can be saved in a data file for use by another procedure.
- Perform repeated measures analysis of variance.
- Display homogeneity tests for testing underlying assumptions in multivariate and univariate analyses.

## ***General Linear Model (GLM) and MANOVA***

MANOVA, the other generalized procedure for analysis of variance and covariance, is available only in syntax. The major distinction between GLM and MANOVA in terms of statistical design and functionality is that GLM uses a non-full-rank, or overparameterized, indicator variable approach to parameterization of linear models instead of the full-rank reparameterization approach that is used in MANOVA. GLM employs a generalized inverse approach and employs aliasing of redundant parameters to 0. These processes employed by GLM allow greater flexibility in handling a variety of data situations, particularly situations involving empty cells. GLM offers the following features that are unavailable in MANOVA:

- Identification of the general forms of estimable functions.
- Identification of forms of estimable functions that are specific to four types of sums of squares (Types I–IV).

- Tests that use the four types of sums of squares, including Type IV, specifically designed for situations involving empty cells.
- Flexible specification of general comparisons among parameters, using the syntax subcommands `LMATRIX`, `MMATRIX`, and `KMATRIX`; sets of contrasts can be specified that involve any number of orthogonal or nonorthogonal linear combinations.
- Nonorthogonal contrasts for within-subjects factors (using the syntax subcommand `WSFACTORS`).
- Tests against nonzero null hypotheses, using the syntax subcommand `KMATRIX`.
- Feature where estimated marginal means (`EMMEANS`) and standard errors (adjusted for other factors and covariates) are available for all between-subjects and within-subjects factor combinations in the original variable metrics.
- Uncorrected pairwise comparisons among estimated marginal means for any main effect in the model, for both between- and within-subjects factors.
- Feature where post hoc or multiple comparison tests for unadjusted one-way factor means are available for between-subjects factors in ANOVA designs; twenty different types of comparisons are offered.
- Weighted least squares (WLS) estimation, including saving of weighted predicted values and residuals.
- Automatic handling of random effects in random-effects models and mixed models, including generation of expected mean squares and automatic assignment of proper error terms.
- Specification of several types of nested models via dialog boxes with proper use of the interaction operator (\*), due to the nonreparameterized approach.
- Univariate homogeneity-of-variance assumption, tested by using the Levene test.
- Between-subjects factors that do not require specification of levels.
- Profile (interaction) plots of estimated marginal means for visual exploration of interactions involving combinations of between-subjects and/or within-subjects factors.
- Saving of casewise temporary variables for model diagnosis:
  - Predicted values—unstandardized (raw), weighted unstandardized.
  - Residuals—unstandardized, weighted unstandardized, standardized, Studentized, deleted.
  - Standard error of prediction.
  - Cook's distance.
  - Leverage.
- Saving of a datafile in IBM® SPSS® Statistics format with parameter estimates and their degrees of freedom and significance level.

To simplify the presentation, GLM reference material is divided into three sections: *univariate* designs with one dependent variable, *multivariate* designs with several interrelated dependent variables, and *repeated measures* designs, in which the dependent variables represent the same types of measurements, taken at more than one time.

The full syntax diagram for GLM is presented here. The following GLM sections include partial syntax diagrams, showing the subcommands and specifications that are discussed in that section. Individually, those diagrams are incomplete. Subcommands that are listed for univariate designs

are available for any analysis, and subcommands that are listed for multivariate designs can be used in any multivariate analysis, including repeated measures.

## Models

The following examples are models that can be specified by using GLM:

### **Model 1: Univariate or Multivariate Simple and Multiple Regression**

```
GLM Y WITH X1 X2 .
```

```
GLM Y1 Y2 WITH X1 X2 X3 .
```

### **Model 2: Fixed-effects ANOVA and MANOVA**

```
GLM Y1 Y2 BY B .
```

### **Model 3: ANCOVA and Multivariate ANCOVA (MANCOVA)**

```
GLM Y1 Y2 BY B WITH X1 X2 X3 .
```

### **Model 4: Random-effects ANOVA and ANCOVA**

```
GLM Y1 BY C WITH X1 X2  
/RANDOM = C .
```

### **Model 5: Mixed-model ANOVA and ANCOVA**

```
GLM Y1 BY B, C WITH X1 X2  
/RANDOM = C .
```

### **Model 6: Repeated Measures Analysis Using a Split-plot Design**

(Univariate mixed models approach with subject as a random effect)

If *drug* is a between-subjects factor and *time* is a within-subjects factor,

```
GLM Y BY DRUG SUBJECT TIME  
/RANDOM = SUBJECT  
/DESIGN = DRUG SUBJECT*DRUG TIME DRUG*TIME .
```

### **Model 7: Repeated Measures Using the WSFACTOR Subcommand**

Use this model only when there is no random between-subjects effect in the model. For example, if *Y1*, *Y2*, *Y3*, and *Y4* are the dependent variables, measured at times 1 to 4,

```
GLM Y1 Y2 Y3 Y4 BY DRUG  
/WSFACTOR = TIME 4  
/DESIGN .
```

**Model 8: Repeated Measures Doubly Multivariate Model**

Repeated measures fixed-effects MANOVA is also called a doubly multivariate model. Varying or time-dependent covariates are not available. This model can be used only when there is no random between-subjects effect in the model.

```
GLM X11 X12 X13 X21 X22 X23
      Y11 Y12 Y13 Y21 Y22 Y23 BY C D
/MEASURE = X Y
/WSFACTOR = A 2 B 3
/WSDESIGN = A B A*B
/DESIGN = C D.
```

**Model 9: Means Model for ANOVA and MANOVA**

This model takes only fixed-effect factors (no random effects and covariates) and always assumes the highest order of the interactions among the factors. For example, *B*, *D*, and *E* are fixed factors, and *Y1* and *Y2* are two dependent variables. You can specify a means model by suppressing the intercept effect and specifying the highest order of interaction on the `DESIGN` subcommand.

```
GLM Y1 Y2 BY B, D, E
/INTERCEPT = EXCLUDE
/DESIGN = B*D*E.
```

**Custom Hypothesis Specifications**

GLM provides a flexible way to customize hypothesis testing based on the general linear hypothesis  $\mathbf{LBM} = \mathbf{K}$ , where  $\mathbf{B}$  is the parameter vector or matrix. You can specify a customized linear hypothesis by using one or more of the subcommands `LMATRIX`, `MMATRIX`, `KMATRIX`, and `CONTRAST`.

**LMATRIX, MMATRIX, and KMATRIX Subcommands**

- The **L** matrix is called the **contrast coefficients matrix**. This matrix specifies coefficients of contrasts, which can be used for studying the between-subjects effects in the model. One way to define the **L** matrix is by specifying the `CONTRAST` subcommand, on which you select a type of contrast. Another way is to specify your own **L** matrix directly by using the `LMATRIX` subcommand. [For more information, see the topic LMATRIX Subcommand on p. 900.](#)
- The **M** matrix is called the **transformation coefficients matrix**. This matrix provides a transformation for the dependent variables. This transformation can be used to construct contrasts among the dependent variables in the model. The **M** matrix can be specified on the `MMATRIX` subcommand. [For more information, see the topic MMATRIX Subcommand on p. 914.](#)
- The **K** matrix is called the **contrast results matrix**. This matrix specifies the results matrix in the general linear hypothesis. To define your own **K** matrix, use the `KMATRIX` subcommand. [For more information, see the topic KMATRIX Subcommand on p. 902.](#)

For univariate and multivariate models, you can specify one, two, or all three of the **L**, **M**, and **K** matrices. If only one or two types are specified, the unspecified matrices use the defaults that are shown in the following table (read across the rows).

Table 104-1

Default matrices for univariate and multivariate models if one matrix is specified

L matrix	M matrix	K matrix
If LMATRIX is used to specify the L matrix	Default = identity matrix*	Default = zero matrix
Default = intercept matrix†	If MMATRIX is used to specify the M matrix	Default = zero matrix
Default = intercept matrix†	Default = identity matrix*	If KMATRIX is used to specify the K matrix

\* The dimension of the identity matrix is the same as the number of dependent variables that are being studied.

† The intercept matrix is the matrix that corresponds to the estimable function for the intercept term in the model, provided that the intercept term is included in the model. If the intercept term is not included in the model, the L matrix is not defined, and this custom hypothesis test cannot be performed.

### Example

```
GLM Y1 Y2 BY A B
  /LMATRIX = A 1 -1
  /DESIGN A B.
```

Assume that factor *A* has two levels.

- Because there are two dependent variables, this model is a multivariate model with two main factor effects, *A* and *B*.
- A custom hypothesis test is requested by the LMATRIX subcommand.
- Because no MMATRIX or KMATRIX is specified, the M matrix is the default two-dimensional identity matrix, and the K matrix is a zero-row vector (0, 0).

For a repeated measures model, you can specify one, two, or all three of the L, M, and K matrices. If only one or two types are specified, the unspecified matrices use the defaults that are shown in the following table (read across the rows).

Table 104-2

Default matrices for repeated measures models if only one matrix is specified

L matrix	M matrix	K matrix
If LMATRIX is used to specify the L matrix	Default = average matrix*	Default = zero matrix
Default = intercept matrix†	If MMATRIX is used to specify the M matrix	Default = zero matrix
Default = intercept matrix†	Default = average matrix*	If KMATRIX is used to specify the K matrix

\* The average matrix is the transformation matrix that corresponds to the transformation for the between-subjects test. The dimension is the number of measures.

† The intercept matrix is the matrix that corresponds to the estimable function for the intercept term in the model, provided that the intercept term is included in the model. If the intercept term is not included in the model, the **L** matrix is not defined, and this custom hypothesis test cannot be performed.

### Example

```
GLM Y1 Y2 BY A B
  /WSFACTOR TIME (2)
  /MMATRIX Y1 1 Y2 1; Y1 1 Y2 -1
  /DESIGN A B.
```

- Because `WSFACTOR` is specified, this model is a repeated measures model with two between-subjects factors *A* and *B*, and a within-subjects factor, *TIME*.
- A custom hypothesis is requested by the `MMATRIX` subcommand. The **M** matrix is a  $2 \times 2$  matrix:

```
1    1
1   -1
```

- Because the **L** matrix and **K** matrix are not specified, their defaults are used. The default for the **L** matrix is the matrix that corresponds to the estimable function for the intercept term in the between-subjects model, and the default for the **K** matrix is a zero-row vector (0, 0).

### CONTRAST Subcommand

When the `CONTRAST` subcommand is used, an **L** matrix, which is used in custom hypothesis testing, is generated according to the chosen contrast. The **K** matrix is always taken to be the zero matrix. If the model is univariate or multivariate, the **M** matrix is always the identity matrix, and its dimension is equal to the number of dependent variables. For a repeated measures model, the **M** matrix is always the average matrix that corresponds to the average transformation for the dependent variable.



# GLM: Univariate

GLM is available in the Advanced Statistics option.

```
GLM dependent var [BY factor list [WITH covariate list]]
[/RANDOM=factor factor...]
[/REGWGT=varname]
[/METHOD=SSTYPE({1  })]
                {2  }
                {3**}
                {4  }
[/INTERCEPT=[INCLUDE**] [EXCLUDE]]
[/MISSING=[INCLUDE] [EXCLUDE**]]
[/CRITERIA=[EPS({1E-8**})] [ALPHA({0.05**})]
           {a      }      {a      }
[/PRINT = [DESCRIPTIVE] [HOMOGENEITY] [PARAMETER] [ETASQ]
          [GEF] [LOF] [OPOWER] [TEST(LMATRIX)]]
[/PLOT=[SPREADLEVEL] [RESIDUALS]
       [PROFILE (factor factor*factor factor*factor*factor ...)]
       [WITH(covariate={value} [...])]]]
       {MEAN }
[/TEST=effect VS {linear combination [DF(df)]}]
                {value DF (df)      }
[/LMATRIX={{["label"] effect list effect list ...;...}]
           {["label"] effect list effect list ...      }
           {["label"] ALL list; ALL...                  }
           {["label"] ALL list                          }
[/KMATRIX= {number      }
           {number;...}
[/CONTRAST (factor name)={DEVIATION[(refcat)]**      }
                    {SIMPLE [(refcat)]              }
                    {DIFFERENCE                      }
                    {HELMERT                         }
                    {REPEATED                        }
                    {POLYNOMIAL [({1,2,3...})]      }
                    {metric }
                    {SPECIAL (matrix)              }
[/POSTHOC =effect [effect...]
          ({[SNK] [TUKEY] [BTUKEY] [DUNCAN]
           [SCHEFFE] [DUNNETT(refcat)] [DUNNETTL(refcat)]
           [DUNNETTR(refcat)] [BONFERRONI] [LSD] [SIDAK]
           [GT2] [GABRIEL] [FREGW] [QREGW] [T2] [T3] [GH] [C]
           [WALLER ({100** })]])
          {kratio}
          [VS effect]
[/EMMEANS=TABLES({OVERALL      }) [WITH(covariate={value} [...])]]
              {factor      }      {MEAN }
              {factor*factor...}
          [COMPARE ADJ(LSD) (BONFERRONI) (SIDAK)]
[/SAVE=[tempvar [(name)]] [tempvar [(name)]]...]
[/OUTFILE={COVB('savfile'|'dataset')}
          {CORB('savfile'|'dataset')}
```

```

[EFFECT('savfile'|'dataset')] [DESIGN('savfile'|'dataset')]

[/DESIGN={{INTERCEPT...}}]
      {{effect effect...}}

```

\*\* Default if the subcommand or keyword is omitted.

*Temporary variables (tempvar) are:*

```
PRED, WPRED, RESID, WRESID, DRESID, ZRESID, SRESID, SEPREP, COOK, LEVER
```

### Example

```
GLM YIELD BY SEED FERT
  /DESIGN.
```

## Overview

This section describes the use of GLM for univariate analyses. However, most of the subcommands that are described here can be used in any type of analysis with GLM. For additional subcommands that are used in multivariate analysis, see *GLM: Multivariate*. For additional subcommands that are used in repeated measures analysis, see *GLM: Repeated Measures*. For basic specification, syntax rules, and limitations of the GLM procedures, see *GLM*.

### Options

**Design Specification.** You can use the DESIGN subcommand to specify which terms to include in the design. This allows you to estimate a model other than the default full factorial model, incorporate factor-by-covariate interactions or covariate-by-covariate interactions, and indicate nesting of effects.

**Contrast Types.** You can specify contrasts other than the default deviation contrasts on the CONTRAST subcommand.

**Optional Output.** You can choose from a variety of optional output on the PRINT subcommand. Output that is appropriate to univariate designs includes descriptive statistics for each cell, parameter estimates, Levene's test for equality of variance across cells, partial eta-squared for each effect and each parameter estimate, the general estimable function matrix, and a contrast coefficients table (**L'** matrix). The OUTFILE subcommand allows you to write out the covariance or correlation matrix, the design matrix, or the statistics from the between-subjects ANOVA table into a separate data file.

Using the EMMEANS subcommand, you can request tables of estimated marginal means of the dependent variable and their standard deviations. The SAVE subcommand allows you to save predicted values and residuals in weighted or unweighted and standardized or unstandardized forms. You can use the POSTHOC subcommand to specify different means comparison tests for comparing all possible pairs of cell means. In addition, you can specify your own hypothesis tests by specifying an **L** matrix and a **K** matrix to test the univariate hypothesis  $\mathbf{LB} = \mathbf{K}$ .

**Basic Specification**

- The basic specification is a variable list identifying the dependent variable, the factors (if any), and the covariates (if any).
- By default, GLM uses a model that includes the intercept term, the covariate (if any), and the full factorial model, which includes all main effects and all possible interactions among factors. The intercept term is excluded if it is excluded in the model by specifying the keyword `EXCLUDE` on the `INTERCEPT` subcommand. Sums of squares are calculated and hypothesis tests are performed by using type-specific estimable functions. Parameters are estimated by using the normal equation and a generalized inverse of the SSCP matrix.

**Subcommand Order**

- The variable list must be specified first.
- Subcommands can be used in any order.

**Syntax Rules**

- For many analyses, the GLM variable list and the `DESIGN` subcommand are the only specifications that are needed.
- If you do not enter a `DESIGN` subcommand, GLM uses a full factorial model, with main effects of covariates, if any.
- At least one dependent variable must be specified, and at least one of the following specifications must occur: `INTERCEPT`, a between-subjects factor, or a covariate. The design contains the intercept by default.
- If more than one `DESIGN` subcommand is specified, only the last subcommand is in effect.
- Dependent variables and covariates must be numeric, but factors can be numeric or string variables.
- If a string variable is specified as a factor, only the first eight bytes of each value are used in distinguishing among values.
- If more than one `MISSING` subcommand is specified, only the last subcommand is in effect.
- The following words are reserved as keywords or internal commands in the GLM procedure:  
`INTERCEPT`, `BY`, `WITH`, `ALL`, `OVERALL`, `WITHIN`  
Variable names that duplicate these words should be changed before you run GLM.

**Limitations**

- Any number of factors can be specified, but if the number of between-subjects factors plus the number of split variables exceeds 18, the Descriptive Statistics table is not printed even when you request it.
- Memory requirements depend primarily on the number of cells in the design. For the default full factorial model, this equals the product of the number of levels or categories in each factor.

**Example**

```
GLM YIELD BY SEED FERT WITH RAINFALL  
/PRINT=DESCRIPTIVE PARAMETER
```

```
/DESIGN.
```

- *YIELD* is the dependent variable; *SEED* and *FERT* are factors; *RAINFALL* is a covariate.
- The `PRINT` subcommand requests the descriptive statistics for the dependent variable for each cell and the parameter estimates, in addition to the default tables Between-Subjects Factors and Univariate Tests.
- The `DESIGN` subcommand requests the default design (a full factorial model with a covariate). This subcommand could have been omitted or could have been specified in full as

```
/DESIGN = INTERCEPT RAINFALL, SEED, FERT, SEED BY FERT.
```

## GLM Variable List

The variable list specifies the dependent variable, the factors, and the covariates in the model.

- The dependent variable must be the first specification on `GLM`.
- The names of the factors follow the dependent variable. Use the keyword `BY` to separate the factors from the dependent variable.
- Enter the covariates, if any, following the factors. Use the keyword `WITH` to separate covariates from factors (if any) and the dependent variable.

### Example

```
GLM DEPENDNT BY FACTOR1 FACTOR2, FACTOR3.
```

- In this example, three factors are specified.
- A default full factorial model is used for the analysis.

### Example

```
GLM Y BY A WITH X
/DESIGN.
```

- In this example, the `DESIGN` subcommand requests the default design, which includes the intercept term, the covariate *X*, and the factor *A*.

## RANDOM Subcommand

`RANDOM` allows you to specify which effects in your design are random. When the `RANDOM` subcommand is used, a table of expected mean squares for all effects in the design is displayed, and an appropriate error term for testing each effect is calculated and used automatically.

- Random always implies a univariate mixed-model analysis.
- If you specify an effect on `RANDOM`, higher-order effects containing the specified effect (excluding any effects containing covariates) are automatically treated as random effects.
- The keyword `INTERCEPT` and effects containing covariates are not allowed on this subcommand.

- The RANDOM subcommand cannot be used if there is any within-subjects factor in the model (that is, RANDOM cannot be specified if WSFACTOR is specified).
- When the RANDOM subcommand is used, the appropriate error terms for the hypothesis testing of all effects in the model are automatically computed and used.
- More than one RANDOM subcommand is allowed. The specifications are accumulated.

### Example

```
GLM DEP BY A B
  /RANDOM = B
  /DESIGN = A, B, A*B.
```

- In the example, effects  $B$  and  $A*B$  are considered as random effects. If only effect  $B$  is specified in the RANDOM subcommand,  $A*B$  is automatically considered as a random effect.
- The hypothesis testing for each effect in the design ( $A$ ,  $B$ , and  $A*B$ ) will be carried out by using the appropriate error term, which is calculated automatically.

## REGWGT Subcommand

The only specification on REGWGT is the name of the variable containing the weights to be used in estimating a weighted least-squares model.

- Specify a numeric weight variable name following the REGWGT subcommand. Only observations with positive values in the weight variable will be used in the analysis.
- If more than one REGWGT subcommand is specified, only the last subcommand is in effect.

### Example

```
GLM OUTCOME BY TREATMNT
  /REGWGT WT.
```

- The procedure performs a weighted least-squares analysis. The variable  $WT$  is used as the weight variable.

## METHOD Subcommand

METHOD controls the computational aspects of the GLM analysis. You can specify one of four different methods for partitioning the sums of squares. If more than one METHOD subcommand is specified, only the last subcommand is in effect.

- |                  |   |
|------------------|---|
| <b>SSTYPE(1)</b> | <i>Type I sum-of-squares method.</i> The Type I sum-of-squares method is also known as the hierarchical decomposition of the sum-of-squares method. Each term is adjusted only for the terms that precede it on the DESIGN subcommand. Under a balanced design, it is an orthogonal decomposition, and the sums of squares in the model add up to the total sum of squares. |
| <b>SSTYPE(2)</b> | <i>Type II sum-of-squares method.</i> This method calculates the sum of squares of an effect in the model, adjusted for all other “appropriate” effects. An appropriate effect is an effect that corresponds to all effects that do not contain the effect that is being examined.  |

For any two effects  $F1$  and  $F2$  in the model,  $F1$  is **contained** in  $F2$  under the following three conditions: Both effects  $F1$  and  $F2$  have the same covariate (if any),  $F2$  consists of more factors than  $F1$ , or all factors in  $F1$  also appear in  $F2$ . The intercept effect is treated as contained in all the pure factor effects. However, the intercept effect is not contained in any effect involving a covariate. No effect is contained in the intercept effect. Thus, for any one effect  $F$  of interest, all other effects in the model can be classified as being in one of the following two groups: the effects that do not contain  $F$  or the effects that contain  $F$ .

If the model is a main-effects design (that is, only main effects are in the model), the Type II sum-of-squares method is equivalent to the regression approach sums of squares, meaning that each main effect is adjusted for every other term in the model.

**SSTYPE(3)** *Type III sum-of-squares method.* This setting is the default. This method calculates the sum of squares of an effect  $F$  in the design as the sum of squares adjusted for any other effects that do not contain it, and *orthogonal* to any effects (if any) that contain it. The Type III sums of squares have one major advantage—they are invariant with respect to the cell frequencies as long as the general form of estimability remains constant. Hence, this type of sums of squares is often used for an unbalanced model with no missing cells. In a factorial design with no missing cells, this method is equivalent to the Yates' weighted squares of means technique, and it also coincides with the overparameterized  $\Sigma$ -restricted model.

**SSTYPE(4)** *Type IV sum-of-squares method.* This method is designed for a situation in which there are missing cells. For any effect  $F$  in the design, if  $F$  is not contained in any other effect, then Type IV = Type III = Type II. When  $F$  is contained in other effects, Type IV equitably distributes the contrasts being made among the parameters in  $F$  to all higher-level effects.

### Example

```
GLM DEP BY A B C
  /METHOD=SSTYPE(3)
  /DESIGN=A, B, C.
```

- The design is a main-effects model.
- The METHOD subcommand requests that the model be fitted with Type III sums of squares.

## INTERCEPT Subcommand

INTERCEPT controls whether an intercept term is included in the model. If more than one INTERCEPT subcommand is specified, only the last subcommand is in effect.

**INCLUDE** *Include the intercept term.* The intercept (constant) term is included in the model. This setting is the default.

**EXCLUDE** *Exclude the intercept term.* The intercept term is excluded from the model. Specification of the keyword INTERCEPT on the DESIGN subcommand overrides INTERCEPT = EXCLUDE.

## MISSING Subcommand

By default, cases with missing values for any of the variables on the GLM variable list are excluded from the analysis. The MISSING subcommand allows you to include cases with user-missing values.

- If MISSING is not specified, the default is EXCLUDE.

- Pairwise deletion of missing data is not available in GLM.
- Keywords INCLUDE and EXCLUDE are mutually exclusive.
- If more than one MISSING subcommand is specified, only the last subcommand is in effect.

**EXCLUDE**            *Exclude both user-missing and system-missing values.* This setting is the default when MISSING is not specified.

**INCLUDE**           *Treat user-missing values as valid.* System-missing values cannot be included in the analysis.

## **CRITERIA Subcommand**

CRITERIA controls the statistical criteria used to build the models.

- More than one CRITERIA subcommand is allowed. The specifications are accumulated. Conflicts across CRITERIA subcommands are resolved by using the conflicting specification that was given on the last CRITERIA subcommand.
- The keyword must be followed by a positive number in parentheses.

**EPS(n)**            *The tolerance level in redundancy detection.* This value is used for redundancy checking in the design matrix. The default value is 1E-8.

**ALPHA(n)**        *The alpha level.* This keyword has two functions. First, the keyword gives the alpha level at which the power is calculated for the  $F$  test. After the noncentrality parameter for the alternative hypothesis is estimated from the data, the power is the probability that the test statistic is greater than the critical value under the alternative hypothesis. (The observed power is displayed by default for GLM.) The second function of alpha is to specify the level of the confidence interval. If the specified alpha level is  $n$ , the value  $(1-n)\times 100$  indicates the level of confidence for all individual and simultaneous confidence intervals that are generated for the specified model. The value of  $n$  must be between 0 and 1, exclusive. The default value of alpha is 0.05, which means that the default power calculation is at the 0.05 level, and the default level of the confidence intervals is 95%, because  $(1-0.05)\times 100=95$ .

## **PRINT Subcommand**

PRINT controls the display of optional output.

- Some PRINT output applies to the entire GLM procedure and is displayed only once.
- Additional output can be obtained on the EMMEANS, PLOT, and SAVE subcommands.
- Some optional output may greatly increase the processing time. Request only the output that you want to see.
- If no PRINT command is specified, default output for a univariate analysis includes a factor information table and a Univariate Tests table (ANOVA) for all effects in the model.
- If more than one PRINT subcommand is specified, only the last subcommand is in effect.

The following keywords are available for GLM univariate analyses. For information about PRINT specifications that are appropriate for other GLM models, see *GLM: Multivariate* and *GLM: Repeated Measures*.

<b>DESCRIPTIVES</b>	<i>Basic information about each cell in the design.</i> This process determines observed means, standard deviations, and counts for the dependent variable in all cells. The cells are constructed from the highest-order crossing of the between-subjects factors. For a multivariate model, statistics are given for each dependent variable. If the number of between-subjects factors plus the number of split variables exceeds 18, the Descriptive Statistics table is not printed.
<b>HOMOGENEITY</b>	<i>Tests of homogeneity of variance.</i> Levene's test for equality of variances for the dependent variable across all level combinations of the between-subjects factors. If there are no between-subjects factors, this keyword is not valid. For a multivariate model, tests are displayed for each dependent variable.
<b>PARAMETER</b>	<i>Parameter estimates.</i> Parameter estimates, standard errors, <i>t</i> tests, and confidence intervals.
<b>ETASQ</b>	<i>Partial eta-squared (<math>\eta^2</math>).</i> This value is an overestimate of the actual effect size in an <i>F</i> test. It is defined as $\eta^2 = \frac{dfh \times F}{dfh \times F + dfe}$ where <i>F</i> is the test statistic and <i>dfh</i> and <i>dfe</i> are its degrees of freedom and degrees of freedom for error. The keyword EFSIZE can be used in place of ETASQ.
<b>GEF</b>	<i>General estimable function table.</i> This table shows the general form of the estimable functions.
<b>LOF</b>	<i>Instruction to perform a lack-of-fit test (which requires at least one cell to have multiple observations).</i> If the test is rejected, it implies that the current model cannot adequately account for the relationship between the response variable and the predictors. Either a variable is omitted or extra terms are needed in the model.
<b>OPOWER</b>	<i>Observed power for each test.</i> The observed power gives the probability that the <i>F</i> test would detect a population difference between groups that is equal to the difference that is implied by the sample difference.
<b>TEST(LMATRIX)</b>	<i>Set of contrast coefficients (L) matrices.</i> The transpose of the <b>L</b> matrix ( <b>L'</b> ) is displayed. This set always includes one matrix displaying the estimable function for each between-subjects effect that appears or is implied in the DESIGN subcommand. Also, any <b>L</b> matrices generated by the LMATRIX or CONTRAST subcommands are displayed. TEST(ESTIMABLE) can be used in place of TEST(LMATRIX).

### Example

```
GLM DEP BY A B WITH COV
  /PRINT=DESCRIPTIVE, TEST(LMATRIX), PARAMETER
  /DESIGN.
```

- Because the design in the DESIGN subcommand is not specified, the default design is used. In this case, the design includes the intercept term, the covariate *COV*, and the full factorial terms of *A* and *B*, which are *A*, *B*, and *A\*B*.
- For each combination of levels of *A* and *B*, the descriptive statistics of *DEP* are displayed.



- The set of L matrices that generates the sums of squares for testing each effect in the design is displayed.
- The parameter estimates, their standard errors,  $t$  tests, confidence intervals, and the observed power for each test are displayed.

## **PLOT Subcommand**

PLOT provides a variety of plots that are useful in checking the assumptions that are needed in the analysis. The PLOT subcommand can be specified more than once. All of the plots that are requested on each PLOT subcommand are produced.

Use the following keywords on the PLOT subcommand to request plots:

<b>SPREADLEVEL</b>	<i>Spread-versus-level plots.</i> Plots are produced that are plots of observed cell means versus standard deviations and versus variances.
<b>RESIDUALS</b>	<i>Observed by predicted by standardized residuals plot.</i> A plot is produced for each dependent variable. In a univariate analysis, a plot is produced for the single dependent variable.
<b>PROFILE</b>	<p><i>Line plots of dependent variable means for one-way, two-way, or three-way crossed factors.</i> The PROFILE keyword must be followed by parentheses containing a list of one or more factor combinations. All specified factors (either individual or crossed) must be composed of only valid factors on the factor list. Factor combinations on the PROFILE keyword may use an asterisk (*) or the keyword BY to specify crossed factors. A factor cannot occur in a single factor combination more than once.</p> <p>The order of factors in a factor combination is important, and there is no restriction on the order of factors. If a single factor is specified after the PROFILE keyword, a line plot of estimated means at each level of the factor is produced. If a two-way crossed factor combination is specified, the output includes a multiple-line plot of estimated means at each level of the first specified factor, with a separate line drawn for each level of the second specified factor. If a three-way crossed factor combination is specified, the output includes multiple-line plots of estimated means at each level of the first specified factor, with separate lines for each level of the second factor and separate plots for each level of the third factor.</p> <p>The PROFILE keyword may be followed by the WITH keyword and a parenthesized list of covariates and values. Plots are produced with the values of the covariates fixed at the specified values. The keyword MEAN causes the covariate to be fixed at its observed mean value.</p>

### **Example**

```
GLM DEP BY A B
  /PLOT = SPREADLEVEL PROFILE(A A*B A*B*C) WITH(X=1.0)
  /DESIGN.
```

Assume that each of the factors  $A$ ,  $B$ , and  $C$  has three levels.

- Spread-versus-level plots are produced, showing observed cell means versus standard deviations and observed cell means versus variances.
- Five profile plots are produced. For factor  $A$ , a line plot of estimated means at each level of  $A$  is produced (one plot). For the two-way crossed factor combination  $A*B$ , a multiple-line plot of estimated means at each level of  $A$  is produced (one plot), with a separate line for each

level of  $B$ . For the three-way crossed factor combination  $A*B*C$ , a multiple-line plot of estimated means at each level of  $A$  is produced for each of the three levels of  $C$  (three plots), with a separate line for each level of  $B$ . Plots are produced with the value of the covariate  $X$  fixed at 1.0.

## TEST Subcommand

The TEST subcommand allows you to test a hypothesis term against a specified error term.

- TEST is valid only for univariate analyses. Multiple TEST subcommands are allowed, with each subcommand being executed independently.
- You must specify both the hypothesis term and the error term. There is no default.
- The hypothesis term is specified before the keyword VS and must be a valid effect that is specified or implied on the DESIGN subcommand.
- The error term is specified after the keyword VS. You can specify either a linear combination or a value. The linear combination of effects takes the general form: coefficient\*effect +/- coefficient\*effect ...
- All effects in the linear combination must be specified or implied on the DESIGN subcommand. Effects that are specified or implied on DESIGN but not listed after VS are assumed to have a coefficient of 0.
- Duplicate effects are allowed. GLM adds coefficients associated with the same effect before performing the test. For example, the linear combination  $5*A-0.9*B-A$  is combined to  $4*A-0.9B$ .
- A coefficient can be specified as a fraction with a positive denominator (for example,  $1/3$  or  $-1/3$  are valid, but  $1/-3$  is invalid).
- If you specify a value for the error term, you must specify the degrees of freedom after the keyword DF. The degrees of freedom must be a positive real number. DF and the degrees of freedom are optional for a linear combination.

### Example

```
GLM DEP BY A B
  /TEST = A VS B + A*B
  /DESIGN = A, B, A*B.
```

- $A$  is tested against the pooled effect of  $B + A*B$ .

## LMATRIX Subcommand

The LMATRIX subcommand allows you to customize your hypotheses tests by specifying the  $\mathbf{L}$  matrix (contrast coefficients matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ , where  $\mathbf{K} = \mathbf{0}$  if it is not specified on the KMATRIX subcommand. The vector  $\mathbf{B}$  is the parameter vector in the linear model.

- The basic format for the LMATRIX subcommand is an optional label in quotation marks, one or more effect names or the keyword ALL, and one or more lists of real numbers.

- The optional label is a string with a maximum length of 255 bytes. Only one label can be specified.
- Only valid effects that appear or are implied on the `DESIGN` subcommand can be specified on the `LMATRIX` subcommand.
- The length of the list of real numbers must be equal to the number of parameters (including the redundant parameters) corresponding to that effect. For example, if the effect  $A*B$  uses six columns in the design matrix, the list after  $A*B$  must contain exactly six numbers.
- A number can be specified as a fraction with a positive denominator (for example,  $1/3$  or  $-1/3$  are valid, but  $1/-3$  is invalid).
- A semicolon (;) indicates the end of a row in the **L** matrix.
- When `ALL` is specified, the length of the list that follows `ALL` is equal to the total number of parameters (including the redundant parameters) in the model.
- Effects that appear or are implied on the `DESIGN` subcommand must be explicitly specified here.
- Multiple `LMATRIX` subcommands are allowed. Each subcommand is treated independently.

### Example

```
GLM DEP BY A B
/LMATRIX = "B1 vs B2 at A1"
           B 1 -1 0 A*B 1 -1 0 0 0 0 0 0 0
/LMATRIX = "Effect A"
           A 1 0 -1
           A*B 1/3 1/3 1/3
                0 0 0
                -1/3 -1/3 -1/3;
           A 0 1 -1
           A*B 0 0 0
                1/3 1/3 1/3
                -1/3 -1/3 -1/3
/LMATRIX = "B1 vs B2 at A2"
           ALL 0
                0 0 0
                1 -1 0
                0 0 0 1 -1 0 0 0 0
/DESIGN = A, B, A*B.
```

Assume that factors  $A$  and  $B$  each have three levels. There are three `LMATRIX` subcommands; each subcommand is treated independently.

- **B1 Versus B2 at A1.** In the first `LMATRIX` subcommand, the difference is tested between levels 1 and 2 of effect  $B$  when effect  $A$  is fixed at level 1. Because there are three levels each in effects  $A$  and  $B$ , the interaction effect  $A*B$  should use nine columns in the design matrix.
- **Effect A.** In the second `LMATRIX` subcommand, effect  $A$  is tested. Because there are three levels in effect  $A$ , no more than two independent contrasts can be formed; thus, there are two rows in the **L** matrix, which are separated by a semicolon (;). The first row tests the difference between levels 1 and 3 of effect  $A$ , while the second row tests the difference between levels 2 and 3 of effect  $A$ .
- **B1 Versus B2 at A2.** In the last `LMATRIX` subcommand, the keyword `ALL` is used. The first 0 corresponds to the intercept effect; the next three instances of 0 correspond to effect  $A$ .

## ***KMATRIX Subcommand***

The `KMATRIX` subcommand allows you to customize your hypothesis tests by specifying the **K** matrix (contrast results matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ . The vector **B** is the parameter vector in the linear model.

- The default **K** matrix is a zero matrix; that is,  $\mathbf{LB} = \mathbf{0}$  is assumed.
- For the `KMATRIX` subcommand to be valid, at least one of the following subcommands must be specified: the `LMATRIX` subcommand or the `INTERCEPT = INCLUDE` subcommand.
- If `KMATRIX` is specified but `LMATRIX` is not specified, the `LMATRIX` is assumed to take the row vector corresponding to the intercept in the estimable function, provided that the subcommand `INTERCEPT = INCLUDE` is specified. In this case, the **K** matrix can be only a scalar matrix.
- If `KMATRIX` and `LMATRIX` are specified, the number of rows in the requested **K** and **L** matrices must be equal. If there are multiple `LMATRIX` subcommands, all requested **L** matrices must have the same number of rows, and **K** must have the same number of rows as these **L** matrices.
- A semicolon (;) can be used to indicate the end of a row in the **K** matrix.
- If more than one `KMATRIX` subcommand is specified, only the last subcommand is in effect.

### ***Example***

```
GLM DEP BY A B
  /LMATRIX = "Effect A"
            A 1 0 -1; A 1 -1 0
  /LMATRIX = "Effect B"
            B 1 0 -1; B 1 -1 0
  /KMATRIX = 0; 0
  /DESIGN = A B.
```

In this example, assume that factors *A* and *B* each have three levels.

- There are two `LMATRIX` subcommands; both subcommands have two rows.
- The first `LMATRIX` subcommand tests whether the effect of *A* is 0, while the second `LMATRIX` subcommand tests whether the effect of *B* is 0.
- The `KMATRIX` subcommand specifies that the **K** matrix also has two rows, each row with value 0.

## ***CONTRAST Subcommand***

`CONTRAST` specifies the type of contrast that is desired among the levels of a factor. For a factor with *k* levels or values, the contrast type determines the meaning of its *k*−1 degrees of freedom.

- Specify the factor name in parentheses following the subcommand `CONTRAST`.
- You can specify only one factor per `CONTRAST` subcommand, but you can enter multiple `CONTRAST` subcommands.
- After closing the parentheses, enter an equals sign followed by one of the contrast keywords.
- This subcommand creates an **L** matrix where the columns corresponding to the factor match the contrast that is given. The other columns are adjusted so that the **L** matrix is estimable.

The following contrast types are available:

**DEVIATION** *Deviations from the grand mean.* This setting is the default for between-subjects factors. Each level of the factor except one is compared to the grand mean. One category (by default, the last category) must be omitted so that the effects will be independent of one another. To omit a category other than the last category, specify the number of the omitted category (which is not necessarily the same as its value) in parentheses after the keyword `DEVIATION`. An example is as follows:

```
GLM Y BY B /CONTRAST(B)=DEVIATION(1).
```

Suppose factor *B* has three levels, with values 2, 4, and 6. The specified contrast omits the first category, in which *B* has the value 2. Deviation contrasts are not orthogonal.

**POLYNOMIAL** *Polynomial contrasts.* This setting is the default for within-subjects factors. The first degree of freedom contains the linear effect across the levels of the factor, the second degree of freedom contains the quadratic effect, and so on. In a balanced design, polynomial contrasts are orthogonal. By default, the levels are assumed to be equally spaced; you can specify unequal spacing by entering a metric consisting of one integer for each level of the factor in parentheses after the keyword `POLYNOMIAL`. (All metrics that are specified cannot be equal; thus, (1, 1, . . . 1) is not valid.) An example is as follows:

```
GLM RESPONSE BY STIMULUS /CONTRAST(STIMULUS) =
POLYNOMIAL(1,2,4).
```

Suppose that factor *STIMULUS* has three levels. The specified contrast indicates that the three levels of *STIMULUS* are actually in the proportion 1:2:4. The default metric is always (1, 2, . . . *k*), where *k* levels are involved. Only the relative differences between the terms of the metric matter. (1, 2, 4) is the same metric as (2, 3, 5) or (20, 30, 50) because, in each instance, the difference between the second and third numbers is twice the difference between the first and second.

**DIFFERENCE** *Difference or reverse Helmert contrasts.* Each level of the factor (except the first level) is compared to the mean of the previous levels. In a balanced design, difference contrasts are orthogonal.

**HELMERT** *Helmert contrasts.* Each level of the factor (except the last level) is compared to the mean of subsequent levels. In a balanced design, Helmert contrasts are orthogonal.

**SIMPLE** *Contrast where each level of the factor (except the last level) is compared to the last level.* To use a category other than the last category as the omitted reference category, specify the category's number (which is not necessarily the same as its value) in parentheses following the keyword `SIMPLE`. An example is as follows:

```
GLM Y BY B /CONTRAST(B)=SIMPLE(1).
```

Suppose that factor *B* has three levels with values 2, 4, and 6. The specified contrast compares the other levels to the first level of *B*, in which *B* has the value 2. Simple contrasts are not orthogonal.

**REPEATED** *Comparison of adjacent levels.* Each level of the factor (except the last level) is compared to the next level. Repeated contrasts are not orthogonal.

**SPECIAL** *A user-defined contrast.* Values that are specified after this keyword are stored in a matrix in column major order. For example, if factor *A* has three levels, then `CONTRAST(A)=SPECIAL(1 1 1 1 -1 0 0 1 -1)` produces the following contrast matrix:

```
1  1  0
1 -1  1
1  0 -1
```

**Orthogonal contrasts** are particularly useful. In a balanced design, contrasts are orthogonal if the sum of the coefficients in each contrast row is 0 and if, for any pair of contrast rows, the products of corresponding coefficients sum to 0. DIFFERENCE, HELMERT, and POLYNOMIAL contrasts always meet these criteria in balanced designs.

### Example

```
GLM DEP BY FAC
  /CONTRAST (FAC) =DIFFERENCE
  /DESIGN.
```

- Suppose that the factor *FAC* has five categories and, therefore, has four degrees of freedom.
- CONTRAST requests DIFFERENCE contrasts, which compare each level (except the first level) with the mean of the previous levels.

## POSTHOC Subcommand

POSTHOC allows you to produce multiple comparisons between means of a factor. These comparisons are usually not planned at the beginning of the study but are suggested by the data during the course of study.

- Post hoc tests are computed for the dependent variable. The alpha value that is used in the tests can be specified by using the keyword ALPHA on the CRITERIA subcommand. The default alpha value is 0.05. The confidence level for any confidence interval that is constructed is  $(1-\alpha)\times 100$ . The default confidence level is 95. For a multivariate model, tests are computed for all specified dependent variables.
- Only between-subjects factors that appear in the factor list are valid in this subcommand. Individual factors can be specified.
- You can specify one or more effects to be tested. Only fixed main effects that appear or are implied on the DESIGN subcommand are valid test effects.
- Optionally, you can specify an effect defining the error term following the keyword VS after the test specification. The error effect can be any single effect in the design that is not the intercept or a main effect that is named on a POSTHOC subcommand.
- A variety of multiple comparison tests are available. Some tests are designed for detecting homogeneity subsets among the groups of means, some tests are designed for pairwise comparisons among all means, and some tests can be used for both purposes.
- For tests that are used for detecting homogeneity subsets of means, non-empty group means are sorted in ascending order. Means that are not significantly different are included together to form a homogeneity subset. The significance for each homogeneity subset of means is displayed. In a case where the numbers of valid cases are not equal in all groups, for most post hoc tests, the harmonic mean of the group sizes is used as the sample size in the calculation. For QREGW or FREGW, individual sample sizes are used.

- For tests that are used for pairwise comparisons, the display includes the difference between each pair of compared means, the confidence interval for the difference, and the significance. The sample sizes of the two groups that are being compared are used in the calculation.
- Output for tests that are specified on the `POSTHOC` subcommand is available according to their statistical purposes. The following table illustrates the statistical purpose of the post hoc tests:

Post Hoc Tests	Statistical Purpose	
	Homogeneity Subsets Detection	Pairwise Comparison and Confidence Interval
LSD		Yes
SIDAK		Yes
BONFERRONI		Yes
GH		Yes
T2		Yes
T3		Yes
C		Yes
DUNNETT		Yes*
DUNNETTL		Yes*
DUNNETTR		Yes*
SNK	Yes	
BTUKEY	Yes	
DUNCAN	Yes	
QREGW	Yes	
FREGW	Yes	
WALLER	Yes†	
TUKEY	Yes	Yes
SCHEFFE	Yes	Yes
GT2	Yes	Yes
GABRIEL	Yes	Yes

\* Only CIs for differences between test group means and control group means are given.

† No significance for Waller test is given.

- Tests that are designed for homogeneity subset detection display the detected homogeneity subsets and their corresponding significances.
- Tests that are designed for both homogeneity subset detection and pairwise comparisons display both kinds of output.
- For the `DUNNETT`, `DUNNETTL`, and `DUNNETTR` keywords, only individual factors can be specified.
- The default reference category for `DUNNETT`, `DUNNETTL`, and `DUNNETTR` is the last category. An integer that is greater than 0, specified within parentheses, can be used to specify a different reference category. For example, `POSTHOC = A (DUNNETT(2))` requests a `DUNNETT` test for factor *A*, using the second level of *A* as the reference category.
- The keywords `DUNCAN`, `DUNNETT`, `DUNNETTL`, and `DUNNETTR` must be spelled out in full; using the first three characters alone is not sufficient.

- If the `REGWGT` subcommand is specified, weighted means are used in performing post hoc tests.
- Multiple `POSTHOC` subcommands are allowed. Each specification is executed independently so that you can test different effects against different error terms.

<b>SNK</b>	<i>Student-Newman-Keuls procedure based on the Studentized range test.</i>
<b>TUKEY</b>	<i>Tukey's honestly significant difference.</i> This test uses the Studentized range statistic to make all pairwise comparisons between groups.
<b>BTUKEY</b>	<i>Tukey's b.</i> This procedure is a multiple comparison procedure based on the average of Studentized range tests.
<b>DUNCAN</b>	<i>Duncan's multiple comparison procedure based on the Studentized range test.</i>
<b>SCHEFFE</b>	<i>Scheffé's multiple comparison t test.</i>
<b>DUNNETT(refcat)</b>	<i>Dunnett's two-tailed t test.</i> Each level of the factor is compared to a reference category. A reference category can be specified in parentheses. The default reference category is the last category. This keyword must be spelled out in full.
<b>DUNNETTL(refcat)</b>	<i>Dunnett's one-tailed t test.</i> This test indicates whether the mean at any level (except the reference category) of the factor is <i>smaller</i> than the mean of the reference category. A reference category can be specified in parentheses. The default reference category is the last category. This keyword must be spelled out in full.
<b>DUNNETTR(refcat)</b>	<i>Dunnett's one-tailed t test.</i> This test indicates whether the mean at any level (except the reference category) of the factor is <i>larger</i> than the mean of the reference category. A reference category can be specified in parentheses. The default reference category is the last category. This keyword must be spelled out in full.
<b>BONFERRONI</b>	<i>Bonferroni t test.</i> This test is based on Student's <i>t</i> statistic and adjusts the observed significance level based on the fact that multiple comparisons are made.
<b>LSD</b>	<i>Least significant difference t test.</i> This test is equivalent to multiple <i>t</i> tests between all pairs of groups. This test does not control the overall probability of rejecting the hypotheses that some pairs of means are different, while in fact they are equal.
<b>SIDAK</b>	<i>Sidak t test.</i> This test provides tighter bounds than the Bonferroni test.
<b>GT2</b>	<i>Hochberg's GT2.</i> This test is a pairwise comparisons test based on the Studentized maximum modulus test. Unless the cell sizes are extremely unbalanced, this test is fairly robust even for unequal variances.
<b>GABRIEL</b>	<i>Gabriel's pairwise comparisons test based on the Studentized maximum modulus test.</i>
<b>FREGW</b>	<i>Ryan-Einot-Gabriel-Welsch's multiple stepdown procedure based on an F test.</i>
<b>QREGW</b>	<i>Ryan-Einot-Gabriel-Welsch's multiple stepdown procedure based on the Studentized range test.</i>
<b>T2</b>	<i>Tamhane's T2.</i> Tamhane's pairwise comparisons test based on a <i>t</i> test. This test can be applied in situations where the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.
<b>T3</b>	<i>Dunnett's T3.</i> Pairwise comparisons test based on the Studentized maximum modulus. This test is appropriate when the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.



<b>GH</b>	<i>Games and Howell's pairwise comparisons test based on the Studentized range test.</i> This test can be applied in situations where the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.
<b>C</b>	<i>Dunnett's C.</i> Pairwise comparisons based on the weighted average of Studentized ranges. This test can be applied in situations where the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.
<b>WALLER(kratio)</b>	<i>Waller-Duncan t test.</i> This test uses a Bayesian approach. The test is restricted to cases with equal sample sizes. For cases with unequal sample sizes, the harmonic mean of the sample size is used. The kratio is the Type 1/Type 2 error seriousness ratio. The default value is 100. You can specify an integer that is greater than 1, enclosed within parentheses.

## **EMMEANS Subcommand**

EMMEANS displays estimated marginal means of the dependent variable in the cells (with covariates held at their overall mean value) and their standard errors of the means for the specified factors. These means are predicted, not observed, means. The estimated marginal means are calculated by using a modified definition by Searle, Speed, and Milliken (1980).

- TABLES, followed by an option in parentheses, is required. COMPARE is optional; if specified, COMPARE must follow TABLES.
- Multiple EMMEANS subcommands are allowed. Each subcommand is treated independently.
- If identical EMMEANS subcommands are specified, only the last identical subcommand is in effect. EMMEANS subcommands that are redundant but not identical (for example, crossed factor combinations such as  $A*B$  and  $B*A$ ) are all processed.

<b>TABLES(option)</b>	<p><i>Table specification.</i> Valid options are the keyword OVERALL, factors appearing on the factor list, and crossed factors that are constructed of factors on the factor list. Crossed factors can be specified by using an asterisk (*) or the keyword BY. All factors in a crossed factor specification must be unique.</p> <p>If OVERALL is specified, the estimated marginal means of the dependent variable are displayed, collapsing over between-subjects factors.</p> <p>If a between-subjects factor, or a crossing of between-subjects factors, is specified on the TABLES keyword, GLM collapses over any other between-subjects factors before computing the estimated marginal means for the dependent variable. For a multivariate model, GLM collapses over any other between-subjects or within-subjects factors.</p> <p>The TABLES keyword may be followed by the WITH keyword and a parenthesized list of covariates and values. Tables are produced with the values of the covariates fixed at the specified values. The keyword MEAN causes the covariate to be fixed at its observed mean value.</p>
<b>COMPARE(factor) ADJ(method)</b>	<p><i>Main-effects or simple-main-effects omnibus tests and pairwise comparisons of the dependent variable.</i> This option gives the mean difference, standard error, significance, and confidence interval for each pair of levels for the effect that is specified in the TABLES command, as well as an omnibus test for that effect. If only one factor is specified on TABLES, COMPARE can be specified by itself; otherwise, the factor specification is required. In this case, levels of the specified factor are compared with each other for each level of the other factors in the interaction.</p>

The optional `ADJ` keyword allows you to apply an adjustment to the confidence intervals and significance values to account for multiple comparisons. Available methods are `LSD` (no adjustment), `BONFERRONI`, or `SIDAK`.

If `OVERALL` is specified on `TABLES`, `COMPARE` is invalid.

### Example

```
GLM DEP BY A B
  /EMMEANS = TABLES(A*B) WITH(X=1.0) COMPARE(A)
  /DESIGN.
```

- The output of this analysis includes a pairwise comparisons table for the dependent variable *DEP*.
- Assume that *A* has three levels and *B* has two levels. The first level of *A* is compared with the second and third levels, the second level is compared with the first and third levels, and the third level is compared with the first and second levels, with the value of the covariate *X* fixed at 1.0. The pairwise comparison is repeated for the two levels of *B*.

## SAVE Subcommand

Use `SAVE` to add one or more residual or fit values to the active dataset.

- Specify one or more temporary variables, each variable followed by an optional new name in parentheses. For a multivariate model, you can optionally specify a new name for the temporary variable related to each dependent variable.
- `WPRED` and `WRESID` can be saved only if `REGWGT` has been specified.
- Specifying a temporary variable on this subcommand results in a variable being added to the active data file for each dependent variable.
- You can specify variable names for the temporary variables. These names must be unique, valid variable names. For a multivariate model, there should be as many variable names specified as there are dependent variables, and names should be listed in the order of the dependent variables as specified on the `GLM` command. If you do not specify enough variable names, default variable names are used for any remaining variables.
- If new names are not specified, `GLM` generates a rootname by using a shortened form of the temporary variable name with a suffix. For a multivariate model, the suffix `_n` is added to the temporary variable name, where *n* is the ordinal number of the dependent variable as specified on the `GLM` command.
- If more than one `SAVE` subcommand is specified, only the last subcommand is in effect.

<b>PRED</b>	<i>Unstandardized predicted values.</i>
<b>WPRED</b>	<i>Weighted unstandardized predicted values.</i> This setting is available only if <code>REGWGT</code> has been specified.
<b>RESID</b>	<i>Unstandardized residuals.</i>
<b>WRESID</b>	<i>Weighted unstandardized residuals.</i> This setting is available only if <code>REGWGT</code> has been specified.
<b>DRESID</b>	<i>Deleted residuals.</i>
<b>ZRESID</b>	<i>Standardized residuals.</i>

<b>SRESID</b>	<i>Studentized residuals.</i>
<b>SEPRE</b>	<i>Standard errors of predicted value.</i>
<b>COOK</b>	<i>Cook's distances.</i>
<b>LEVER</b>	<i>Uncentered leverage values.</i>

## **OUTFILE Subcommand**

The **OUTFILE** subcommand writes data to a data file or previously declared dataset that can be used in other procedures.

- You must specify a keyword on **OUTFILE**. There is no default.
- You must specify a quoted file specification or previously declared dataset name (**DATASET DECLARE** command) in parentheses after a keyword. The asterisk (\*) is not allowed.
- If you specify more than one keyword, a different filename is required for each keyword.
- If more than one **OUTFILE** subcommand is specified, only the last subcommand is in effect.
- For **COVB** or **CORB**, the output will contain, in addition to the covariance or correlation matrix, three rows for each dependent variable: a row of parameter estimates, a row of residual degrees of freedom, and a row of significance values for the *t* statistics corresponding to the parameter estimates. All statistics are displayed separately by split.

<b>COVB</b> ('savfile' 'dataset')	<i>Writes the parameter covariance matrix.</i>
<b>CORB</b> ('savfile' 'dataset')	<i>Writes the parameter correlation matrix.</i>
<b>EFFECT</b> ('savfile' 'dataset')	<i>Writes the statistics from the between-subjects ANOVA table. This specification is invalid for repeated measures analyses.</i>
<b>DESIGN</b> ('savfile' 'dataset')	<i>Writes the design matrix. The number of rows equals the number of cases, and the number of columns equals the number of parameters. The variable names are <i>DES_1</i>, <i>DES_2</i>, ..., <i>DES_p</i>, where <i>p</i> is the number of the parameters.</i>

## **DESIGN Subcommand**

**DESIGN** specifies the effects included in a specific model. The cells in a design are defined by all of the possible combinations of levels of the factors in that design. The number of cells equals the product of the number of levels of all the factors. A design is *balanced* if each cell contains the same number of cases. **GLM** can analyze both balanced and unbalanced designs.

- Specify a list of terms to be included in the model, and separate the terms by spaces or commas.
- The default design, if the **DESIGN** subcommand is omitted or is specified by itself, is a design consisting of the following terms in order: the intercept term (if **INTERCEPT=INCLUDE** is specified), the covariates that are given in the covariate list, and the full factorial model defined by all factors on the factor list and excluding the intercept.
- To include a term for the main effect of a factor, enter the name of the factor on the **DESIGN** subcommand.
- To include the intercept term in the design, use the keyword **INTERCEPT** on the **DESIGN** subcommand. If **INTERCEPT** is specified on the **DESIGN** subcommand, the subcommand **INTERCEPT=EXCLUDE** is overridden.

- To include a term for an interaction between factors, use the keyword `BY` or the asterisk (`*`) to join the factors that are involved in the interaction. For example,  $A*B$  means a two-way interaction effect of  $A$  and  $B$ , where  $A$  and  $B$  are factors.  $A*A$  is not allowed because factors inside an interaction effect must be distinct.
- To include a term for nesting one effect within another effect, use the keyword `WITHIN` or use a pair of parentheses on the `DESIGN` subcommand. For example,  $A(B)$  means that  $A$  is nested within  $B$ . The expression  $A(B)$  is equivalent to the expression  $A$  *WITHIN*  $B$ . When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus,  $A(B)(C)$  is not valid.
- Multiple nesting is allowed. For example,  $A(B(C))$  means that  $B$  is nested within  $C$  and  $A$  is nested within  $B(C)$ .
- Interactions between nested effects are not valid. For example, neither  $A(C)*B(C)$  nor  $A(C)*B(D)$  is valid.
- To include a covariate term in the design, enter the name of the covariate on the `DESIGN` subcommand.
- Covariates can be connected—but not nested—through the `*` operator to form another covariate effect. Therefore, interactions among covariates such as  $X1*X1$  and  $X1*X2$  are valid but not  $X1(X2)$ . Using covariate effects such as  $X1*X1$ ,  $X1*X1*X1$ ,  $X1*X2$ , and  $X1*X1*X2*X2$  makes fitting a polynomial regression model easy in GLM.
- Factor and covariate effects can be connected only by the `*` operator. Suppose  $A$  and  $B$  are factors and  $X1$  and  $X2$  are covariates. Examples of valid factor-by-covariate interaction effects are  $A*X1$ ,  $A*B*X1$ ,  $X1*A(B)$ ,  $A*X1*X1$ , and  $B*X1*X2$ .
- If more than one `DESIGN` subcommand is specified, only the last subcommand is in effect.

### **Example**

```
GLM Y BY A B C WITH X
  /DESIGN A B(A) X*A.
```

- In this example, the design consists of a main effect  $A$ , a nested effect  $B$  within  $A$ , and an interaction effect of a covariate  $X$  with a factor  $A$ .

# GLM: Multivariate

GLM is available in the Advanced Statistics option.

```
GLM dependent varlist [BY factor list [WITH covariate list]]
[/REGWGT=varname]
[/METHOD=SSTYPE({1  })]
                {2  }
                {3**}
                {4  }
[/INTERCEPT=[INCLUDE**] [EXCLUDE]]
[/MISSING=[INCLUDE] [EXCLUDE**]]
[/CRITERIA=[EPS({1E-8**})] [ALPHA({0.05**})]
           {a      }      {a      }
[/PRINT = [DESCRIPTIVE] [HOMOGENEITY] [PARAMETER][ETASQ] [RSSCP]
          [GEF] [LOF] [OPOWER] [TEST [(SSCP] [LMATRIX] [MMATRIX])]]
[/PLOT=[SPREADLEVEL] [RESIDUALS]
       [PROFILE (factor factor*factor factor*factor*factor ...) [WITH(covariate={value} [...])]]
       {MEAN }
[/LMATRIX={{"label"} effect list effect list ...;...}}
           {{"label"} effect list effect list ...      }
           {{"label"} ALL list; ALL...                  }
           {{"label"} ALL list                          }
[/MMATRIX= {{"label"} depvar value depvar value ...;["label"]...}}
           {{"label"} depvar value depvar value ...      }
           {{"label"} ALL list; ["label"] ...              }
           {{"label"} ALL list                            }
[/KMATRIX= {list of numbers      }
           {list of numbers;...}
[/SAVE=[tempvar [(list of names)]] [tempvar [(list of names)]]...
       [DESIGN]
[/OUTFILE={{COVB('savfile' | 'dataset')}}
          {CORB('savfile' | 'dataset')}}
          [EFFECT('savfile' | 'dataset')] [DESIGN('savfile' | 'dataset')]
[/DESIGN={{INTERCEPT...      }
          {effect effect...}}]
```

\*\* Default if the subcommand or keyword is omitted.

*Temporary variables (tempvar) are:*

PRED, WPRED, RESID, WRESID, DRESID, ZRESID, SRESID, SEPREP, COOK, LEVER

## Example

```
GLM SCORE1 TO SCORE4 BY METHOD(1,3).
```

## Overview

This section discusses the subcommands that are used in multivariate general linear models and covariance designs with several interrelated dependent variables. The discussion focuses on subcommands and keywords that do not apply—or apply in different manners—to univariate analyses. The discussion does not contain information about all subcommands that you will need to specify the design. For subcommands that are not covered here, see *GLM: Univariate*.

### Options

**Optional Output.** In addition to the output that is described in *GLM: Univariate*, you can have both multivariate and univariate  $F$  tests. Using the `PRINT` subcommand, you can request the hypothesis and error sums-of-squares and cross-product matrices for each effect in the design, the transformation coefficient table (M matrix), Box's  $M$  test for equality of covariance matrices, and Bartlett's test of sphericity.

### Basic Specification

- The basic specification is a variable list identifying the dependent variables, with the factors (if any) named after `BY` and the covariates (if any) named after `WITH`.
- By default, `GLM` uses a model that includes the intercept term, the covariates (if any), and the full factorial model, which includes all main effects and all possible interactions among factors. The intercept term is excluded if it is excluded in the model by specifying `EXCLUDE` on the `INTERCEPT` subcommand. `GLM` produces multivariate and univariate  $F$  tests for each effect in the model. `GLM` also calculates the power for each test, based on the default alpha value.

### Subcommand Order

- The variable list must be specified first.
- Subcommands can be used in any order.

### Syntax Rules

- The syntax rules that apply to univariate analysis also apply to multivariate analysis.
- If you enter one of the multivariate specifications in a univariate analysis, `GLM` ignores it.

### Limitations

- Any number of factors can be specified, but if the number of between-subjects factors plus the number of split variables exceeds 18, the Descriptive Statistics table is not printed even when you request it.
- Memory requirements depend primarily on the number of cells in the design. For the default full factorial model, this equals the product of the number of levels or categories in each factor.

## Example

### Multivariate Analysis of Variance (MANOVA)

```
GLM
  los cost  BY clotsolv proc
  /CONTRAST (clotsolv)=Simple(1)
  /METHOD = SSTYPE(3)
  /INTERCEPT = INCLUDE
  /PRINT = ETASQ TEST(SSCP) HOMOGENEITY
  /PLOT = SPREADLEVEL
  /CRITERIA = ALPHA(.05)
  /DESIGN = clotsolv proc clotsolv*proc .
```

- The procedure fits a model for the dependent variables *los* and *cost* using *clotsolv* and *proc* as factors.
- The `CONTRAST` subcommand specifies simple contrasts for *clotsolv*, using the first category as the reference category, to test differences between the categories. No contrasts are specified for *proc*.
- The `PRINT` subcommand requests tabular output for estimates of effect size, SSCP matrices, and homogeneity tests.
- The `PLOT` subcommand requests spread vs. level plots.
- All other options are set to their default values.

## GLM Variable List

- Multivariate `GLM` calculates statistical tests that are valid for analyses of dependent variables that are correlated with one another. The dependent variables must be specified first.
- The factor and covariate lists follow the same rules as in univariate analyses.
- If the dependent variables are uncorrelated, the univariate significance tests have greater statistical power.

## PRINT Subcommand

By default, if no `PRINT` subcommand is specified, multivariate `GLM` produces multivariate tests (MANOVA) and univariate tests (ANOVA) for all effects in the model. All `PRINT` specifications that are described in *GLM: Univariate* are available in multivariate analyses. The following additional output can be requested:

<b>TEST(SSCP)</b>	<i>Sums-of-squares and cross-product matrices.</i> Hypothesis (HSSCP) and error (ESSCP) sums-of-squares and cross-product matrices for each effect in the design are displayed. Each between-subjects effect has a different HSSCP matrix, but there is a single ESSCP matrix for all between-subjects effects. For a repeated measures design, each within-subjects effect has an HSSCP matrix and an ESSCP matrix. If there are no within-subjects effects, the ESSCP matrix for the between-subjects effects is the same as the RSSCP matrix.
<b>TEST(MMATRIX)</b>	<i>Set of transformation coefficients (M) matrices.</i> Any M matrices that are generated by the <code>MMATRIX</code> subcommand are displayed. If no M matrix is specified on the <code>MMATRIX</code> subcommand, this specification is skipped, unless you are using a repeated measures design. In a repeated

measures design, this set always includes the M matrix that is determined by the `WSFACTOR` subcommand. The specification `TEST (TRANSFORM)` is equivalent to `TEST (MMATRIX)`.

### HOMOGENEITY

*Tests of homogeneity of variance.* In addition to Levene's test for equality of variances for each dependent variable, the display includes Box's *M* test of homogeneity of the covariance matrices of the dependent variables across all level combinations of the between-subjects factors.

### RSSCP

*Sums-of-squares and cross-products of residuals.* Three matrices are displayed:

*Residual SSCP matrix.* This matrix is a square matrix of sums of squares and cross-products of residuals. The dimension of this matrix is the same as the number of dependent variables in the model.

*Residual covariance matrix.* This matrix is the residual SSCP matrix divided by the degrees of freedom of the residual.

*Residual correlation matrix.* This matrix is the standardized form of the residual covariance matrix.

### Example

```
GLM Y1 Y2 Y3 BY A B
  /PRINT = HOMOGENEITY RSSCP
  /DESIGN.
```

- Since there are three dependent variables, this model is a multivariate model.
- The keyword `RSSCP` produces three matrices of sums of squares and cross-products of residuals. The output also contains the result of Bartlett's test of the sphericity of the residual covariance matrix.
- In addition to the Levene test for each dependent variable, the keyword `HOMOGENEITY` produces the result of Box's *M* test of homogeneity in the multivariate model.

## MMATRIX Subcommand

The `MMATRIX` subcommand allows you to customize your hypothesis tests by specifying the M matrix (transformation coefficients matrix) in the general form of the linear hypothesis  $LBM = K$ , where  $K = 0$  if it is not specified on the `KMATRIX` subcommand. The vector *B* is the parameter vector in the linear model.

- Specify an optional label in quotation marks. Then either list dependent variable names, each name followed by a real number, or specify the keyword `ALL` followed by a list of real numbers. Only variable names that appear on the dependent variable list can be specified on the `MMATRIX` subcommand.
- You can specify one label for each column in the M matrix.
- If you specify `ALL`, the length of the list that follows `ALL` should be equal to the number of dependent variables.
- There is no limit on the length of the label.
- For the `MMATRIX` subcommand to be valid, at least one of the following specifications must be made: the `LMATRIX` subcommand or `INTERCEPT=INCLUDE`. (Either of these specifications defines an L matrix.)



- If both `LMATRIX` and `MMATRIX` are specified, the **L** matrix is defined by the `LMATRIX` subcommand.
- If `MMATRIX` or `KMATRIX` is specified but `LMATRIX` is not specified, the **L** matrix is defined by the estimable function for the intercept effect, provided that the intercept effect is included in the model.
- If `LMATRIX` is specified but `MMATRIX` is not specified, the **M** matrix is assumed to be an  $r \times r$  identity matrix, where  $r$  is the number of dependent variables.
- A semicolon (;) indicates the end of a column in the **M** matrix.
- Dependent variables that do not appear on a list of dependent variable names and real numbers are assigned a value of 0.
- Dependent variables that do not appear in the `MMATRIX` subcommand will have a row of zeros in the **M** matrix.
- A number can be specified as a fraction with a positive denominator (for example,  $1/3$  or  $-1/3$  is valid, but  $1/-3$  is invalid).
- The number of columns must be greater than 0. You can specify as many columns as you need.
- If more than one `MMATRIX` subcommand is specified, only the last subcommand is in effect.

### **Example**

```
GLM Y1 Y2 Y3 BY A B
  /MMATRIX = "Y1-Y2" Y1 1 Y2 -1; "Y1-Y3" Y1 1 Y3 -1
            "Y2-Y3" Y2 1 Y3 -1
  /DESIGN.
```

- In the above example,  $Y1$ ,  $Y2$ , and  $Y3$  are the dependent variables.
- The `MMATRIX` subcommand requests all pairwise comparisons among the dependent variables.
- Because `LMATRIX` was not specified, the **L** matrix is defined by the estimable function for the intercept effect.

# GLM: Repeated Measures

GLM is available in the Advanced Statistics option.

```
GLM dependent varlist [BY factor list [WITH covariate list]]

/WSFACTOR=name levels [{DEVIATION [(refcat)]          }] name...
                      {SIMPLE [(refcat)]             }
                      {DIFFERENCE                    }
                      {HELMERT                        }
                      {REPEATED                       }
                      {POLYNOMIAL [({1,2,3...})]**}
                      {          {metric }           }
                      {SPECIAL (matrix)              }

[/MEASURE=newname newname...]

[/WSDESIGN=effect effect...]

[/REGWGT=varname]

[/METHOD=SSTYPE({1  })]
                  {2  }
                  {3**}
                  {4  }

[/INTERCEPT=[INCLUDE]** [EXCLUDE]]

[/MISSING=[INCLUDE] [EXCLUDE]**]

[/PRINT = [DESCRIPTIVE] [HOMOGENEITY] [PARAMETER][ETASQ] [RSSCP]
          [GEF] [LOF] [OPOWER] [TEST [({SSCP [LMATRIX] [MMATRIX])]}]]

[/SAVE=[tempvar [(list of names)]] [tempvar [(list of names)]]...]
      [DESIGN]

[/EMMEANS=TABLES({OVERALL          })] [COMPARE ADJ(LSD) (BONFERRONI) (SIDAK)]
      {factor                       }
      {factor*factor...             }
      {wsfactor                     }
      {wsfactor*wsfactor...        }
      {factor*...wsfactor*...      }
      {factor*factor...            }

[/DESIGN={ [INTERCEPT... ]}*
          {[effect effect...]}]
```

\* The DESIGN subcommand has the same syntax as is described in *GLM: Univariate*.

\*\* Default if the subcommand or keyword is omitted.

## Example

```
GLM Y1 TO Y4 BY GROUP
  /WSFACTOR=YEAR 4.
```

## Overview

This section discusses the subcommands that are used in repeated measures designs, in which the dependent variables represent measurements of the same variable (or variables) taken repeatedly. This section does not contain information on all of the subcommands that you will need to specify

the design. For some subcommands or keywords not covered here, such as `DESIGN`, see *GLM: Univariate*. For information on optional output and the multivariate significance tests available, see *GLM: Multivariate*.

- In a simple repeated measures analysis, all dependent variables represent different measurements of the same variable for different values (or levels) of a within-subjects factor. Between-subjects factors and covariates can also be included in the model, just as in analyses not involving repeated measures.
- A **within-subjects factor** is simply a factor that distinguishes measurements made on the same subject or case, rather than distinguishing different subjects or cases.
- GLM permits more complex analyses, in which the dependent variables represent levels of two or more within-subjects factors.
- GLM also permits analyses in which the dependent variables represent measurements of several variables for the different levels of the within-subjects factors. These are known as **doubly multivariate designs**.
- A repeated measures analysis includes a within-subjects design describing the model to be tested with the within-subjects factors, as well as the usual between-subjects design describing the effects to be tested with between-subjects factors. The default for the within-subjects factors design is a full factorial model which includes the main within-subjects factor effects and all their interaction effects.
- If a custom hypothesis test is required (defined by the `CONTRAST`, `LMATRIX`, or `KMATRIX` subcommands), the default transformation matrix (**M** matrix) is taken to be the average transformation matrix, which can be displayed by using the keyword `TEST(MMATRIX)` on the `PRINT` subcommand. The default contrast result matrix (**K** matrix) is the zero matrix.
- If the contrast coefficient matrix (**L** matrix) is not specified, but a custom hypothesis test is required by the `MMATRIX` or the `KMATRIX` subcommand, the contrast coefficient matrix (**L** matrix) is taken to be the **L** matrix which corresponds to the estimable function for the intercept in the between-subjects model. This matrix can be displayed by using the keyword `TEST(LMATRIX)` on the `PRINT` subcommand.

### **Basic Specification**

- The basic specification is a variable list followed by the `WSFACTOR` subcommand.
- Whenever `WSFACTOR` is specified, GLM performs special repeated measures processing. The multivariate and univariate tests are provided. In addition, for any within-subjects effect involving more than one transformed variable, the Mauchly test of sphericity is displayed to test the assumption that the covariance matrix of the transformed variables is constant on the diagonal and zero off the diagonal. The Greenhouse-Geisser epsilon and the Huynh-Feldt epsilon are also displayed for use in correcting the significance tests in the event that the assumption of sphericity is violated.

### **Subcommand Order**

- The list of dependent variables, factors, and covariates must be first.

**Syntax Rules**

- The `WSFACTOR` (within-subjects factors), `WSDESIGN` (within-subjects design), and `MEASURE` subcommands are used only in repeated measures analysis.
- `WSFACTOR` is required for any repeated measures analysis.
- If `WSDESIGN` is not specified, a full factorial within-subjects design consisting of all main effects and all interactions among within-subjects factors is used by default.
- The `MEASURE` subcommand is used for doubly multivariate designs, in which the dependent variables represent repeated measurements of more than one variable.

**Limitations**

- Any number of factors can be specified, but if the number of between-subjects factors plus the number of split variables exceeds 18, the Descriptive Statistics table is not printed even when you request it.
- Maximum of 18 within-subjects factors.
- Memory requirements depend primarily on the number of cells in the design. For the default full factorial model, this equals the product of the number of levels or categories in each factor.

**Examples****Repeated Measures ANOVA**

```
GLM Y1 TO Y4 BY GROUP
  /WSFACTOR=YEAR 4 POLYNOMIAL
  /WSDESIGN=YEAR
  /PRINT=PARAMETER
  /DESIGN=GROUP.
```

- `WSFACTOR` specifies a repeated measures analysis in which the four dependent variables represent a single variable measured at four levels of the within-subjects factor. The within-subjects factor is called *YEAR* for the duration of the GLM procedure.
- `POLYNOMIAL` requests polynomial contrasts for the levels of *YEAR*. Because the four variables, *Y1*, *Y2*, *Y3*, and *Y4*, in the active dataset represent the four levels of *YEAR*, the effect is to perform an orthonormal polynomial transformation of these variables.
- `PRINT` requests that the parameter estimates be displayed.
- `WSDESIGN` specifies a within-subjects design that includes only the effect of the *YEAR* within-subjects factor. Because *YEAR* is the only within-subjects factor specified, this is the default design, and `WSDESIGN` could have been omitted.
- `DESIGN` specifies a between-subjects design that includes only the effect of the *GROUP* between-subjects factor. This subcommand could have been omitted.

**Repeated Measures ANOVA, Unbalanced Design with Missing Cells**

```
GLM
  sales.1 sales.2 sales.3 sales.4 BY promo marketid
  /WSFACTOR = week 4
  /MEASURE = sales
  /METHOD = SSTYPE(4)
```

```

/PRINT = ETASQ TEST(SSCP) PARAMETER RSSCP HOMOGENEITY
/PLOT = PROFILE ( week*promo )
/PLOT = PROFILE ( week*marketid )
/CRITERIA = ALPHA(.05)
/WSDESIGN = week
/DESIGN = promo marketid promo*marketid.

```

- The procedure fits a model to the dependent variables *sales.1* through *sales.4* using *promo* and *marketid* as factors.
- *WSFACTOR* specifies a repeated measures analysis in which the four dependent variables represent a single variable measured at four levels of the within-subjects factor. The within-subjects factor is called *week* for the duration of the GLM procedure.
- *METHOD* specifies the Type IV method for partitioning sums of squares. This is recommended because the cell frequencies for the layout of *Market ID* and *Promotion \* Market ID* are unbalanced and *Promotion \* Market ID* has empty cells.
- *PRINT* requests that the estimates of effect size, SSCP matrices, and homogeneity tests be displayed.
- The first *PLOT* subcommand requests profile plots for *week\*promo*, which will result in a single plot for each dependent variable. Each plot will feature the levels of *week* on the *x*-axis and the estimated marginal means for the levels of *week* on the *y*-axis. A separate line is produced for each level of *promo*.
- The first *PLOT* subcommand requests profile plots for *week\*marketid*, which will result in a single plot for each dependent variable. Each plot will feature the levels of *week* on the *x*-axis and the estimated marginal means for the levels of *week* on the *y*-axis. A separate line is produced for each level of *marketid*.
- All other options are set to their default values.

### **Doubly Multivariate ANOVA**

```

GLM
  tg0 tg1 tg2 tg3 tg4 wgt0 wgt1 wgt2 wgt3 wgt4 BY gender
/WSFACTOR = time 5 Repeated
/MEASURE = tg wgt
/METHOD = SSTYPE(3)
/PLOT = PROFILE( time*gender )
/EMMEANS = TABLES( gender*time )
/PRINT = ETASQ TEST(SSCP)
/CRITERIA = ALPHA(.05)
/WSDESIGN = time
/DESIGN = gender .

```

- The procedure fits a model to the dependent variables *tg0* through *wgt4* using *gender* as a factor.
- *WSFACTOR* and *MEASURE* specify a repeated measures analysis in which the ten dependent variables represent two variables measured at five levels of the within-subjects factor. The within-subjects factor is called *time* and the measures are called *tg* and *wgt* for the duration of the GLM procedure. Repeated contrasts are specified for the within-subjects factor.
- *PLOT* requests profile plots for *time\*gender*, which will result in a single plot for each measure. Each plot will feature the levels of *week* on the *x*-axis and the estimated marginal means for the measure at the levels of *week* on the *y*-axis. A separate line is produced for each level of *gender*.

- EMMEANS requests estimated marginal means for *gender\*time*, which will result in a tabular representation of the profile plots.
- PRINT requests that the estimates of effect size and SSCP matrices be displayed.
- All other options are set to their default values.

## GLM Variable List

The list of dependent variables, factors, and covariates must be specified first.

- WSFACTOR determines how the dependent variables on the GLM variable list will be interpreted.
- The number of dependent variables on the GLM variable list must be a multiple of the number of cells in the within-subjects design. If there are six cells in the within-subjects design, each group of six dependent variables represents a single within-subjects variable that has been measured in each of the six cells.
- Normally, the number of dependent variables should equal the number of cells in the within-subjects design multiplied by the number of variables named on the MEASURE subcommand (if one is used). If you have more groups of dependent variables than are accounted for by the MEASURE subcommand, GLM will choose variable names to label the output, which may be difficult to interpret.
- Covariates are specified after keyword WITH. You can specify constant covariates. Constant covariates represent variables whose values remain the same at each within-subjects level.

### Example

```
GLM MATH1 TO MATH4 BY METHOD WITH SES
  /WSFACTOR=SEMESTER 4.
```

- The four dependent variables represent a score measured four times (corresponding to the four levels of *SEMESTER*).
- *SES* is a constant covariate. Its value does not change over the time covered by the four levels of *SEMESTER*.
- Default contrast (*POLYNOMIAL*) is used.

## WSFACTOR Subcommand

WSFACTOR names the within-subjects factors, specifies the number of levels for each, and specifies the contrast for each.

- Presence of the WSFACTOR subcommand implies that the repeated measures model is being used.
- Mauchly's test of sphericity is automatically performed when WSFACTOR is specified.
- Names and number levels for the within-subjects factors are specified on the WSFACTOR subcommand. Factor names must not duplicate any of the dependent variables, factors, or covariates named on the GLM variable list. A type of contrast can also be specified for each within-subjects factor in order to perform comparisons among its levels. This contrast amounts to a transformation on the dependent variables.

- If there are more than one within-subjects factors, they must be named in the order corresponding to the order of the dependent variables on the GLM variable list. GLM varies the levels of the last-named within-subjects factor most rapidly when assigning dependent variables to within-subjects cells (see the example below).
- The number of cells in the within-subjects design is the product of the number of levels for all within-subjects factors.
- Levels of the factors must be represented in the data by the dependent variables named on the GLM variable list.
- The number of levels of each factor must be at least two. Enter an integer equal to or greater than 2 after each factor to indicate how many levels the factor has. Optionally, you can enclose the number of levels in parentheses.
- Enter only the *number of levels* for within-subjects factors, not a range of values.
- If more than one WSFACOR subcommand is specified, only the last one is in effect.

### **Contrasts for WSFACOR**

The levels of a within-subjects factor are represented by different dependent variables. Therefore, contrasts between levels of such a factor compare these dependent variables. Specifying the type of contrast amounts to specifying a transformation to be performed on the dependent variables.

- In testing the within-subjects effects, an orthonormal transformation is automatically performed on the dependent variables in a repeated measures analysis.
- The contrast for each within-subjects factor is entered after the number of levels. If no contrast keyword is specified, POLYNOMIAL ( 1 , 2 , 3 . . . ) is the default. This contrast is used in comparing the levels of the within-subjects factors. Intrinsically orthogonal contrast types are recommended for within-subjects factors if you wish to examine each degree-of-freedom test, provided compound symmetry is assumed within each within-subjects factor. Other orthogonal contrast types are DIFFERENCE and HELMERT.
- If there are more than one within-subjects factors, the transformation matrix (**M** matrix) is computed as the Kronecker product of the matrices generated by the contrasts specified.
- The transformation matrix (**M** matrix) generated by the specified contrasts can be displayed by using the keyword TEST (MMATRIX) on the subcommand PRINT.
- The contrast types available for within-subjects factors are the same as those on the CONTRAST subcommand for between-subjects factors, described in CONTRAST Subcommand on p. 902 in *GLM: Univariate*.

The following contrast types are available:

- DEVIATION**      *Deviations from the grand mean.* This is the default for between-subjects factors. Each level of the factor except one is compared to the grand mean. One category (by default the last) must be omitted so that the effects will be independent of one another. To omit a category other than the last, specify the number of the omitted category in parentheses after the keyword DEVIATION. For example, GLM Y1 Y2 Y3 BY GROUP /WSFACTOR = Y 3 DEVIATION (1)  
Deviation contrasts are not orthogonal.

<b>POLYNOMIAL</b>	<p><i>Polynomial contrasts.</i> This is the default for within-subjects factors. The first degree of freedom contains the linear effect across the levels of the factor, the second contains the quadratic effect, and so on. In a balanced design, polynomial contrasts are orthogonal. By default, the levels are assumed to be equally spaced; you can specify unequal spacing by entering a metric consisting of one integer for each level of the factor in parentheses after the keyword <code>POLYNOMIAL</code>. (All metrics specified cannot be equal; thus (1,1,...,1) is not valid.) For example, <code>/WSFACTOR=D 3 POLYNOMIAL(1,2,4)</code>.</p> <p>Suppose that factor D has three levels. The specified contrast indicates that the three levels of D are actually in the proportion 1:2:4. The default metric is always (1,2,...,k), where k levels are involved. Only the relative differences between the terms of the metric matter (1,2,4) is the same metric as (2,3,5) or (20,30,50) because, in each instance, the difference between the second and third numbers is twice the difference between the first and second.</p>									
<b>DIFFERENCE</b>	<p><i>Difference or reverse Helmert contrasts.</i> Each level of the factor except the first is compared to the mean of the previous levels. In a balanced design, difference contrasts are orthogonal.</p>									
<b>HELMERT</b>	<p><i>Helmert contrasts.</i> Each level of the factor except the last is compared to the mean of subsequent levels. In a balanced design, Helmert contrasts are orthogonal.</p>									
<b>SIMPLE</b>	<p><i>Each level of the factor except the last is compared to the last level.</i> To use a category other than the last as the omitted reference category, specify its number in parentheses following keyword <code>SIMPLE</code>. For example, <code>/WSFACTOR=B 3 SIMPLE (1)</code>.</p> <p>Simple contrasts are not orthogonal.</p>									
<b>REPEATED</b>	<p><i>Comparison of adjacent levels.</i> Each level of the factor except the last is compared to the next level. Repeated contrasts are not orthogonal.</p>									
<b>SPECIAL</b>	<p><i>A user-defined contrast.</i> Values specified after this keyword are stored in a matrix in column major order. For example, if factor A has three levels, then <code>WSFACTOR(A)=SPECIAL(1 1 1 1 -1 0 0 1 -1)</code> produces the following contrast matrix:</p> <table border="0" style="margin-left: 2em;"> <tr><td>1</td><td>1</td><td>0</td></tr> <tr><td>1</td><td>-1</td><td>1</td></tr> <tr><td>1</td><td>0</td><td>-1</td></tr> </table>	1	1	0	1	-1	1	1	0	-1
1	1	0								
1	-1	1								
1	0	-1								

**Example**

```
GLM X1Y1 X1Y2 X2Y1 X2Y2 X3Y1 X3Y2 BY TREATMNT GROUP
  /WSFACTOR=X 3 Y 2
  /DESIGN.
```

- The GLM variable list names six dependent variables and two between-subjects factors, *TREATMNT* and *GROUP*.
- `WSFACTOR` identifies two within-subjects factors whose levels distinguish the six dependent variables. *X* has three levels, and *Y* has two. Thus, there are  $3 \times 2 = 6$  cells in the within-subjects design, corresponding to the six dependent variables.
- Variable *X1Y1* corresponds to levels 1,1 of the two within-subjects factors; variable *X1Y2* corresponds to levels 1,2; *X2Y1* to levels 2,1; and so on up to *X3Y2*, which corresponds to levels 3,2. The first within-subjects factor named, *X*, varies most slowly, and the last within-subjects factor named, *Y*, varies most rapidly on the list of dependent variables.
- Because there is no `WSDSIGN` subcommand, the within-subjects design will include all main effects and interactions: *X*, *Y*, and *X* by *Y*.



- Likewise, the between-subjects design includes all main effects and interactions (*TREATMNT*, *GROUP*, and *TREATMNT* by *GROUP*) plus the intercept.
- In addition, a repeated measures analysis always includes interactions between the within-subjects factors and the between-subjects factors. There are three such interactions for each of the three within-subjects effects.

### Example

```
GLM SCORE1 SCORE2 SCORE3 BY GROUP
  /WSFACTOR=ROUND 3 DIFFERENCE
  /CONTRAST (GROUP)=DEVIATION
  /PRINT=PARAMETER TEST (LMATRIX) .
```

- This analysis has one between-subjects factor, *GROUP*, and one within-subjects factor, *ROUND*, with three levels that are represented by the three dependent variables.
- The *WSFACTOR* subcommand also specifies difference contrasts for *ROUND*, the within-subjects factor.
- There is no *WSDSIGN* subcommand, so a default full factorial within-subjects design is assumed. This could also have been specified as *WSDSIGN=ROUND*, or simply *WSDSIGN*.
- The *CONTRAST* subcommand specifies deviation contrasts for *GROUP*, the between-subjects factor. This subcommand could have been omitted because deviation contrasts are the default.
- *PRINT* requests the display of the parameter estimates for the model and the **L** matrix.
- There is no *DESIGN* subcommand, so a default full factorial between-subjects design is assumed. This could also have been specified as *DESIGN=GROUP*, or simply *DESIGN*.

## WSDSIGN Subcommand

*WSDSIGN* specifies the design for within-subjects factors. Its specifications are like those of the *DESIGN* subcommand, but it uses the within-subjects factors rather than the between-subjects factors.

- The default *WSDSIGN* is a full factorial design, which includes all main effects and all interactions for within-subjects factors. The default is in effect whenever a design is processed without a preceding *WSDSIGN* or when the preceding *WSDSIGN* subcommand has no specifications.
- A *WSDSIGN* specification cannot include between-subjects factors or terms based on them, nor does it accept interval-level variables.
- The keyword *INTERCEPT* is not allowed on *WSDSIGN*.
- Nested effects are not allowed. Therefore, the symbols ( ) are not allowed here.
- If more than one *WSDSIGN* subcommand is specified, only the last one is in effect.

### Example

```
GLM JANLO, JANHI, FEBLO, FEBHI, MARLO, MARHI BY SEX
  /WSFACTOR MONTH 3 STIMULUS 2
  /WSDSIGN MONTH, STIMULUS
  /DESIGN SEX.
```

- There are six dependent variables, corresponding to three months and two different levels of stimulus.
- The dependent variables are named on the GLM variable list in an order such that the level of stimulus varies more rapidly than the month. Thus, *STIMULUS* is named last on the *WSFACTOR* subcommand.
- The *WSDESIGN* subcommand specifies only the main effects for within-subjects factors. There is no *MONTH-by-STIMULUS* interaction term.

## MEASURE Subcommand

In a doubly multivariate analysis, the dependent variables represent multiple variables measured under the different levels of the within-subjects factors. Use *MEASURE* to assign names to the variables that you have measured for the different levels of within-subjects factors.

- Specify a list of one or more variable names to be used in labeling the averaged results. If no within-subjects factor has more than two levels, *MEASURE* has no effect. You can use up to 255 bytes for each name.
- The number of dependent variables in the dependent variables list should equal the product of the number of cells in the within-subjects design and the number of names on *MEASURE*.
- If you do not enter a *MEASURE* subcommand and there are more dependent variables than cells in the within-subjects design, GLM assigns names (normally *MEASURE\_1*, *MEASURE\_2*, and so on) to the different measures.
- All of the dependent variables corresponding to each measure should be listed together and ordered so that the within-subjects factor named last on the *WSFACTORS* subcommand varies most rapidly.

### Example

```
GLM TEMP11 TEMP12 TEMP21 TEMP22 TEMP31 TEMP32,
      WEIGHT11 WEIGHT12 WEIGHT21 WEIGHT22 WEIGHT31 WEIGHT32 BY GROUP
  /WSFACTOR=DAY 3 AMPM 2
  /MEASURE=TEMP WEIGHT
  /WSDESIGN=DAY, AMPM, DAY BY AMPM
  /DESIGN.
```

- There are 12 dependent variables: six temperatures and six weights, corresponding to morning and afternoon measurements on three days.
- *WSFACTOR* identifies the two factors (*DAY* and *AMPM*) that distinguish the temperature and weight measurements for each subject. These factors define six within-subjects cells.
- *MEASURE* indicates that the first group of six dependent variables correspond to *TEMP* and the second group of six dependent variables correspond to *WEIGHT*.
- These labels, *TEMP* and *WEIGHT*, are used on the output as the measure labels.
- *WSDESIGN* requests a full factorial within-subjects model. Because this is the default, *WSDESIGN* could have been omitted.

## ***EMMEANS Subcommand***

EMMEANS displays estimated marginal means of the dependent variables in the cells, adjusted for the effects of covariates at their overall means, for the specified factors. Note that these are predicted, not observed, means. The standard errors are also displayed. [For more information, see the topic EMMEANS Subcommand on p. 907.](#)

- For the TABLES and COMPARE keywords, valid options include the within-subjects factors specified in the WSFACOR subcommand, crossings among them, and crossings among factors specified in the factor list and factors specified on the WSFACOR subcommand.
- All factors in a crossed-factors specification must be unique.
- If a between- or within-subjects factor, or a crossing of between- or within-subjects factors, is specified on the TABLES keyword, then GLM will collapse over any other between- or within-subjects factors before computing the estimated marginal means for the dependent variables.

# GRAPH

GRAPH is available in the Statistics Base option.

GRAPH

```
[/TITLE='line 1' ['line 2']]
[/SUBTITLE='line 1']
[/FOOTNOTE='line 1' ['line 2']]

{/BAR [{(SIMPLE)}]=function/variable specificationt }
      {(GROUPED)}
      {(STACKED)}
      {(RANGE)}

{/LINE [{(SIMPLE)}]=function/variable specificationt }
      {(MULTIPLE)}
      {(DROP)}
      {(AREA)}
      {(DIFFERENCE)}

{/PIE }

{/PARETO[ {(CUM)} ] [ {(SIMPLE)} ] =function/variable specificationt}
      {(NOCUM)} {(STACKED)}

{/HILO[ {(SIMPLE)} ] =function/variable specificationt}
      {(GROUPED)}

{/HISTOGRAM[ (NORMAL) ] =var }

{/SCATTERPLOT[ {(BIVARIATE)} ] =variable specificationt}
      {(OVERLAY)}
      {(MATRIX)}
      {(XYZ)}

{/ERRORBAR[ {(CI[ {95} ] ) } ] = {var [var var ...] [BY var]} }
      {n} {var BY var BY var }
      {(STERRIR[ {12} ] ) }
      {n}
      {(STDDEV[ {2} ] ) }
      {n}

[/PANEL COLVAR=varlist COLOP={CROSS**} ROWVAR=varlist ROWOP={CROSS**}]
      {NEST } {NEST }

[/INTERVAL {CI { (95) }}]
      { (n) }
      {STDDEV { (2) }}
      { (n) }
      {SE { (2) }}
      { (n) }

[/TEMPLATE=file]

[/MISSING={ {LISTWISE**} } [ {NOREPORT**} } [ {EXCLUDE**} ] ]
      {VARIABLE } ] {REPORT } {INCLUDE }
```

\*\* Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

**Release History**

Release 13.0

- PANEL subcommand introduced.
- INTERVAL subcommand introduced.

The following table shows all possible function/variable specifications for BAR, LINE, PIE, BLOCK, and PARETO subcommands. For special restrictions, see the individual subcommands. In the table, *valuef* refers to the value function, *countf* refers to the count functions, and *sumf* refers to the summary functions.

	Simple bar, simple or area line, pie, simple high-low, and simple Pareto charts	Grouped or stacked bar, multiple, drop or difference area, and stacked Pareto charts
<b>Categorical charts</b>	[countf BY] var	[countf BY] var BY var
	sumf(var) BY var	sumf(var) BY var BY var
	sumf(varlist)	sumf(varlist) BY var
	sumf(var) sumf(var)...	sumf(var) sumf(var)... BY var
<b>Noncategorical charts</b>	valuef(var) [BY var]	valuef(varlist) [BY var]

The following table shows all possible function/variable specifications for the HILO subcommand. Categorical variables for simple high-low-close charts must be dichotomous or trichotomous.

Simple range bar and simple high-low-close charts	Clustered range bar and clustered high-low-close charts
[countf BY] var	(sumf(var) sumf(var) [sumf(var)]) (. . .) . . . BY var
sumf(var) sumf(var) sumf(var) BY var	sumf(var) sumf(var) [sumf(var)] BY var BY var
sumf(var) BY var BY var	
valuef(varlist) [BY var]	valuef(varlist) (...) ... [BY var]

Variable specification is required on all types of scatterplots. The following table shows all possible specifications:

BIVARIATE	var WITH var [BY var] [BY var ( {NAME } )] { IDENTIFY }
OVERLAY	varlist WITH varlist [ ( PAIR ) ] [BY var ( {NAME } )] { IDENTIFY }
MATRIX	varlist [BY var] [BY var ( {NAME } )] { IDENTIFY }
XYZ	var WITH var WITH var [BY var] [BY var ( {NAME } )] { IDENTIFY }

*Value function:*

The VALUE function yields the value of the specified variable for each case. It always produces one bar, point, or slice for each case. The VALUE (X) specification implies the value of X by n, where n is the number of each case. You can specify multiple variables, as in:

```
GRAPH /BAR = VALUE(SALARY BONUS BENEFIT) .
```

This command draws a bar chart with the values of *SALARY*, *BONUS*, and *BENEFIT* for each employee (case). A *BY* variable can be used to supply case labels, but it does not affect the layout of the chart, even if values of the *BY* variable are the same for multiple cases.

*Aggregation functions:*

Two groups of aggregation functions are available: count functions and summary functions.

*Count functions:*

<b>COUNT</b>	<i>Frequency of cases in each category.</i>
<b>PCT</b>	<i>Frequency of cases in each category expressed as a percentage of the whole.</i>
<b>CUPCT</b>	<i>Cumulative percentage sorted by category value.</i>
<b>CUFREQ</b>	<i>Cumulative frequency sorted by category value.</i>

- Count functions yield the count or percentage of valid cases within categories determined by one or more *BY* variables, as in:

```
GRAPH /BAR (SIMPLE) = PCT BY REGION.
```

- Count functions do not have any arguments.
- You can omit the keyword **COUNT** and the subsequent keyword **BY** and specify just a variable, as in

```
GRAPH /BAR = DEPT.
```

This command is interpreted as

```
GRAPH /BAR = COUNT BY DEPT.
```

*Summary functions:*

<b>MINIMUM</b>	<i>Minimum value of the variable.</i>
<b>MAXIMUM</b>	<i>Maximum value of the variable.</i>
<b>N</b>	<i>Number of cases for which the variable has a nonmissing value.</i>
<b>SUM</b>	<i>Sum of the values of the variable.</i>
<b>CUSUM</b>	<i>Sum of the summary variable accumulated across values of the category variable.</i>
<b>MEAN</b>	<i>Mean.</i>
<b>STDDEV</b>	<i>Standard deviation.</i>
<b>VARIANCE</b>	<i>Variance.</i>
<b>MEDIAN</b>	<i>Median.</i>
<b>GMEDIAN</b>	<i>Group median.</i>
<b>MODE</b>	<i>Mode.</i>
<b>PTILE(x)</b>	<i>Xth percentile value of the variable. X must be greater than 0 and less than 100.</i>
<b>PLT(x)</b>	<i>Percentage of cases for which the value of the variable is less than x.</i>
<b>PGT(x)</b>	<i>Percentage of cases for which the value of the variable is greater than x.</i>
<b>NLT(x)</b>	<i>Number of cases for which the value of the variable is less than x.</i>
<b>NGT(x)</b>	<i>Number of cases for which the value of the variable is greater than x.</i>

**PIN(x1,x2)** *Percentage of cases for which the value of the variable is greater than or equal to x1 and less than or equal to x2. x1 cannot exceed x2.*

**NIN(x1,x2)** *Number of cases for which the value of the variable is greater than or equal to x1 and less than or equal to x2. x1 cannot exceed x2.*

- Summary functions are usually used with summary variables (variables that record continuous values, such as age or expenses). To use a summary function, specify the name of one or more variables in parentheses after the name of the function, as in:

```
GRAPH /BAR = SUM(SALARY) BY DEPT.
```

- You can specify multiple summary functions for more chart types. For example, the same function can be applied to a list of variables, as in:

```
GRAPH /BAR = SUM(SALARY BONUS BENEFIT) BY DEPT.
```

This syntax is equivalent to:

```
GRAPH /BAR = SUM(SALARY) SUM(BONUS) SUM(BENEFIT) BY DEPT.
```

Different functions can be applied to the same variable, as in:

```
GRAPH /BAR = MEAN(SALARY) MEDIAN(SALARY) BY DEPT.
```

Different functions and variables can be combined, as in:

```
GRAPH /BAR = MIN(SALARY81) MAX(SALARY81)
              MIN(SALARY82) MAX(SALARY82) BY JOBCAT.
```

The effect of multiple summary functions on the structure of the charts is illustrated under the discussion of specific chart types.

## Overview

GRAPH generates a high-resolution chart by computing statistics from variables in the active dataset and constructing the chart according to your specification. The chart can be a bar chart, pie chart, line chart, error bar chart, high-low-close histogram, scatterplot, or Pareto chart. The chart is displayed where high-resolution display is available and can be edited with a chart editor and saved as a chart file.

### Options

**Titles and Footnotes.** You can specify a title, subtitle, and footnote for the chart using the `TITLE`, `SUBTITLE`, and `FOOTNOTE` subcommands.

**Chart Type.** You can request a specific type of chart using the `BAR`, `LINE`, `PIE`, `ERRORBAR`, `HILO`, `HISTOGRAM`, `SCATTERPLOT`, or `PARETO` subcommand.

**Chart Content.** You can specify an aggregated categorical chart using various aggregation functions or a nonaggregated categorical chart using the `VALUE` function.

**Templates.** You can specify a template, using the `TEMPLATE` subcommand, to override the default chart attribute settings on your system.

### **Basic Specification**

The basic specification is a chart type subcommand. By default, the generated chart will have no title, subtitle, or footnote.

### **Subcommand Order**

Subcommands can be specified in any order.

### **Syntax Rules**

- Only one chart type subcommand can be specified.
- The function/variable specification is required for all subtypes of bar, line, error bar, hilo, and Pareto charts; the variable specification is required for histograms and all subtypes of scatterplots.
- The function/variable or variable specifications should match the subtype keywords. If there is a discrepancy, GRAPH produces the default chart for the function/variable or variable specification regardless of the specified keyword.

### **Operations**

- GRAPH computes aggregated functions to obtain the values needed for the requested chart and calculates an optimal scale for charting.
- The chart title, subtitle, and footnote are assigned as they are specified on the TITLE, SUBTITLE, and FOOTNOTE subcommands. If you do not use these subcommands, the chart title, subtitle, and footnote are null. The split-file information is displayed as a subtitle if split-file is in effect.
- GRAPH creates labels that provide information about the source of the values being plotted. Labeling conventions vary for different subtypes. Where variable or value labels are defined in the active dataset, GRAPH uses the labels; otherwise, variable names or values are used.

### **Limitations**

Categorical charts cannot display fewer than 2 or more than 3,000 categories.

## **Examples**

```
GRAPH /BAR=SUM (MURDER) BY CITY.
```

- This command generates a simple (default) bar chart showing the number of murders in each city.
- The category axis (*x* axis) labels are defined by the value labels (or values if no value labels exist) of the variable *CITY*.
- The default span (2) and sigma value (3) are used.
- Since no BY variable is specified, the *x* axis is labeled by sequence numbers.



## ***TITLE, SUBTITLE, and FOOTNOTE Subcommands***

TITLE, SUBTITLE, and FOOTNOTE specify lines of text placed at the top or bottom of the chart.

- One or two lines of text can be specified for TITLE or FOOTNOTE, and one line of text can be specified for SUBTITLE.
- Each line of text must be enclosed in quotes. The maximum length of any line is 72 characters.
- The default font sizes and types are used for the title, subtitle, and footnote.
- By default, the title, subtitle, and footnote are left-aligned with the y axis.
- If you do not specify TITLE, the default title, subtitle, and footnote are null, which leaves more space for the chart. If split-file processing is in effect, the split-file information is provided as a default subtitle.

### ***Example***

```
GRAPH TITLE = 'Murder in Major U.S. Cities'
/SUBTITLE='per 100,000 people'
/FOOTNOTE='The above data was reported on August 26, 1987'
/BAR=SUM(MURDER) BY CITY.
```

## ***BAR Subcommand***

BAR creates one of five types of bar charts using the keywords SIMPLE, COMPOSITIONAL, GROUPED, STACKED, or RANGE.

- Only one keyword can be specified, and it must be specified in parentheses.
- When no keyword is specified, the default is either SIMPLE or GROUPED, depending on the type of function/variable specification.

<b>SIMPLE</b>	<i>Simple bar chart.</i> This is the default if no keyword is specified on the BAR subcommand and the variables define a simple bar chart. A simple bar chart can be defined by a single summary or count function and a single BY variable or by multiple summary functions and no BY variable.
<b>GROUPED</b>	<i>Clustered bar chart.</i> A clustered bar chart is defined by a single function and two BY variables or by multiple functions and a single BY variable. This is the default if no keyword is specified on the BAR subcommand and the variables define a clustered bar chart.
<b>STACKED</b>	<i>Stacked bar chart.</i> A stacked bar chart displays a series of bars, each divided into segments stacked one on top of the other. The height of each segment represents the value of the category. Like a clustered bar chart, it is defined by a single function and two BY variables or by multiple functions and a single BY variable.
<b>RANGE</b>	<i>Range bar chart.</i> A range bar chart displays a series of floating bars. The height of each bar represents the range of the category and its position in the chart indicates the minimum and maximum values. A range bar chart can be defined by a single function and two BY variables or by multiple functions and a single BY variable. If a variable list is used as the argument for a function, the list must be of an even number. If a second BY variable is used to define the range, the variable must be dichotomous.

## **LINE Subcommand**

LINE creates one of five types of line charts using the keywords SIMPLE, MULTIPLE, DROP, AREA, or DIFFERENCE.

- Only one keyword can be specified, and it must be specified in parentheses.
- When no keyword is specified, the default is either SIMPLE or MULTIPLE, depending on the type of function/variable specification.

<b>SIMPLE</b>	<i>Simple line chart.</i> A simple line chart is defined by a single function and a single BY variable or by multiple functions and no BY keyword. This is the default if no keyword is specified on LINE and the data define a simple line.
<b>MULTIPLE</b>	<i>Multiple line chart.</i> A multiple line chart is defined by a single function and two BY variables or by multiple functions and a single BY variable. This is the default if no keyword is specified on LINE and the data define a multiple line.
<b>DROP</b>	<i>Drop-line chart.</i> A drop-line chart shows the difference between two or more fluctuating variables. It is defined by a single function and two BY variables or by multiple functions and a single BY variable.
<b>AREA</b>	<i>Area line chart.</i> An area line chart fills the area beneath each line with a color or pattern. When multiple lines are specified, the second line is the sum of the first and second variables, the third line is the sum of the first, second, and third variables, and so on. The specification is the same as that for a simple or multiple line chart.
<b>DIFFERENCE</b>	<i>Difference area chart.</i> A difference area chart fills the area between a pair of lines. It highlights the difference between two variables or two groups. A difference area chart is defined by a single function and two BY variables or by two summary functions and a single BY variable. If a second BY variable is used to define the two groups, the variable must be dichotomous.

## **PIE Subcommand**

PIE creates pie charts. A pie chart can be defined by a single function and a single BY variable or by multiple summary functions and no BY variable. A pie chart divides a circle into slices. The size of each slice indicates the value of the category relative to the whole. Cumulative functions (CUPCT, CUFREQ, and CUSUM) are inappropriate for pie charts but are not prohibited. When specified, all cases except those in the last category are counted more than once in the resulting pie.

## **HILO Subcommand**

HILO creates one of two types of high-low-close charts using the keywords SIMPLE or GROUPED. High-low-close charts show the range and the closing (or average) value of a series.

- Only one keyword can be specified.

- When a keyword is specified, it must be specified in parentheses.
- When no keyword is specified, the default is either `SIMPLE` or `GROUPED`, depending on the type of function/variable specification.

**SIMPLE** *Simple high-low-close chart.* A simple high-low-close chart can be defined by a single summary or count function and two `BY` variables, by three summary functions and one `BY` variable, or by three values with one or no `BY` variable. When a second `BY` variable is used to define a high-low-close chart, the variable must be dichotomous or trichotomous. If dichotomous, the first value defines low and the second value defines high; if trichotomous, the first value defines high, the second defines low, and the third defines close.

**GROUPED** *Grouped high-low-close chart.* A grouped high-low-close chart is defined by a single function and two `BY` variables or by multiple functions and a single `BY` variable. When a variable list is used for a single function, the list must contain two or three variables. If it contains two variables, the first defines the high value and the second defines the low value. If it contains three variables, the first defines the high value, the second defines the low value, and the third defines the close value. Likewise, if multiple functions are specified, they must be in groups of either two or three. The first function defines the high value, the second defines the low value, and the third, if specified, defines the close value.

## ***ERRORBAR Subcommand***

`ERRORBAR` creates either a simple or a clustered error bar chart, depending on the variable specification on the subcommand. A simple error bar chart is defined by one numeric variable with or without a `BY` variable or a variable list. A clustered error bar chart is defined by one numeric variable with two `BY` variables or a variable list with a `BY` variable.

Error bar charts can display confidence intervals, standard deviations, or standard errors of the mean. To specify the statistics to be displayed, one of the following keywords is required:

- CI value** *Display confidence intervals for mean.* You can specify a confidence level between 50 and 99.9. The default is 95.
- STERROR n** *Display standard errors of mean.* You can specify any positive number for *n*. The default is 2.
- STDDEV n** *Display standard deviations.* You can specify any positive number for *n*. The default is 2.

## ***SCATTERPLOT Subcommand***

`SCATTERPLOT` produces two- or three-dimensional scatterplots. Multiple two-dimensional plots can be plotted within the same frame or as a scatterplot matrix. Only variables can be specified; aggregated functions cannot be plotted. When `SCATTERPLOT` is specified without keywords, the default is `BIVARIATE`.

- BIVARIATE** *One two-dimensional scatterplot.* A basic scatterplot is defined by two variables separated by the keyword `WITH`. This is the default when `SCATTERPLOT` is specified without keywords.
- OVERLAY** *Multiple plots drawn within the same frame.* Specify a variable list on both sides of `WITH`. By default, one scatterplot is drawn for each combination of variables on the left of `WITH` with variables on the right. You can specify `PAIR` in parentheses to indicate that the first variable on the left is paired with the first variable on the right, the second variable on the left with the second variable on the right, and so on. All plots are drawn

within the same frame and are differentiated by color or pattern. The axes are scaled to accommodate the minimum and maximum values across all variables.

- MATRIX** *Scatterplot matrix.* Specify at least two variables. One scatterplot is drawn for each combination of the specified variables above the diagonal and a second below the diagonal in a square matrix.
- XYZ** *One three-dimensional plot.* Specify three variables, each separated from the next with the keyword `WITH`.

- If you specify a control variable using `BY`, `GRAPH` produces a control scatterplot where values of the `BY` variable are indicated by different colors or patterns. A control variable cannot be specified for overlay plots.
- You can display the value label of an identification variable at the plotting position for each case by adding `BY var (NAME)` or `BY var (IDENTIFY)` to the end of any valid scatterplot specification. When the chart is created, `NAME` turns the labels on, while `IDENTIFY` turns the labels off. You can use the Point Selection tool to turn individual labels off or on in the scatterplot.

## ***HISTOGRAM Subcommand***

`HISTOGRAM` creates a histogram.

- Only one variable can be specified on this subcommand.
- `GRAPH` divides the values of the variable into several evenly spaced intervals and produces a bar chart showing the number of times the values for the variable fall within each interval.
- You can request a normal distribution line by specifying the keyword `NORMAL` in parentheses.

## ***PARETO Subcommand***

`PARETO` creates one of two types of Pareto charts. A Pareto chart is used in quality control to identify the few problems that create the majority of nonconformities. Only `SUM`, `VALUE`, and `COUNT` can be used with the `PARETO` subcommand.

Before plotting, `PARETO` sorts the plotted values in descending order by category. The right axis is always labeled by the cumulative percentage from 0 to 100. By default, a cumulative line is displayed. You can eliminate the cumulative line or explicitly request it by specifying one of the following keywords:

- CUM** *Display the cumulative line.* This is the default.
- NOCUM** *Do not display the cumulative line.*

You can request a simple or a stacked Pareto chart by specifying one of the following keywords and define it with appropriate function/variable specifications:

- SIMPLE**     *Simple Pareto chart.* Each bar represents one type of nonconformity. A simple Pareto chart can be defined by a single variable, a single `VALUE` function, a single `SUM` function with a `BY` variable, or a `SUM` function with a variable list as an argument with no `BY` variable.
- STACKED**     *Stacked Pareto chart.* Each bar represents one or more types of nonconformity within the category. A stacked Pareto chart can be defined by a single `SUM` function with two `BY` variables, a single variable with a `BY` variable, a `VALUE` function with a variable list as an argument, or a `SUM` function with a variable list as an argument and a `BY` variable.

## **PANEL Subcommand**

The `PANEL` subcommand specifies the variables and method used for paneling. Each keyword in the subcommand is followed by an equals sign (=) and the value for that keyword.

## **COLVAR and ROWVAR Keywords**

The `COLVAR` and `ROWVAR` keywords identify the column and row variables, respectively. Each category in a column variable appears as a vertical column in the resulting chart. Each category in a row variable appears as a horizontal row in the resulting chart.

- If multiple variables are specified for a keyword, the `COLOP` and `ROWOP` keywords can be used to change the way in which variable categories are rendered in the chart.
- The `ROWVAR` keyword is not available for population pyramids.

**varlist**             *The list of variables used for paneling.*

### **Examples**

```
GRAPH
  /BAR(SIMPLE)=COUNT BY educ
  /PANEL COLVAR=gender COLOP=CROSS
```

- There are two columns in the resulting paneled chart, one for each gender.
- Because there is only one paneling variable, there are only as many panels as there are variable values. Therefore, there are two panels.

```
GRAPH
  /BAR(SIMPLE)=COUNT BY educ
  /PANEL COLVAR=minority ROWVAR=jobcat.
```

- There are two columns in the resulting paneled chart (for the *gender* variable values) and three rows (for the *jobcat* variable values).

## **COLOP and ROWOP Keywords**

The COLOP and ROWOP keywords specify the paneling method for the column and row variables, respectively. These keywords have no effect on the chart if there is only one variable in the rows and/or columns. They also have no effect if the data are not nested.

- CROSS** *Cross variables in the rows or columns.* When the variables are crossed, a panel is created for every combination of categories in the variables. For example, if the categories in one variable are A and B and the categories in another variable are 1 and 2, the resulting chart will display a panel for the combinations of A and 1, A and 2, B and 1, and B and 2. A panel can be empty if the categories in that panel do not cross (for example, if there are no cases in the B category *and* the 1 category). This is the default.
- NEST** *Nest variables in the rows or columns.* When the variables are nested, a panel is created for each category that is nested in the parent category. For example, if the data contain variables for states and cities, a panel is created for each city and the relevant state. However, panels are not created for cities that are not in certain states, as would happen with CROSS. When nesting, make sure that the variables specified for ROWVAR or COLVAR are in the correct order. Parent variables precede child variables.

### **Example**

Assume you have the following data:

Table 108-1  
*Nested data*

State	City	Temperature
NJ	Springfield	70
MA	Springfield	60
IL	Springfield	50
NJ	Trenton	70
MA	Boston	60

You can create a paneled chart from these data with the following syntax:

```
GRAPH
  /HISTOGRAM=temperature
  /PANEL COLVAR=state city COLOP=CROSS.
```

The command crosses every variable value to create the panels. Because not every state contains every city, the resulting paneled chart will contain blank panels. For example, there will be a blank panel for Springfield and New Jersey. In this dataset, the *city* variable is really nested in the *state* variable. To nest the variables in the panels and eliminate any blank panels, use the following syntax:

```
GRAPH
  /HISTOGRAM=temperature
  /PANEL COLVAR=state city COLOP=NEST.
```

## **INTERVAL Subcommand**

The `INTERVAL` subcommand adds error bars to the chart. This is different from the `ERRORBAR` subcommand. The `ERRORBAR` subcommand adds error bar data elements. `INTERVAL` adds error bars to other data elements (for example, areas, bars, and lines).

Error bars indicate the variability of the summary statistic being displayed. The length of the error bar on either side of the summary statistic represents a confidence interval or a specified number of standard errors or standard deviations. `GRAPH` supports error bars for simple or clustered categorical charts displaying means, medians, counts, and percentages.

The keywords are not followed by an equals sign (=). They are followed by a value in parentheses.

### **Example**

```
GRAPH
  /BAR(SIMPLE)=COUNT BY jobcat
  /INTERVAL CI(95).
```

### **CI Keyword**

**(value)**            *The percentage of the confidence interval to use as the length of the error bars.*

### **STDDEV Keyword**

**(value)**            *A multiplier indicating the number of standard deviations to use as the length of the error bars.*

### **SE Keyword**

**(value)**            *A multiplier indicating the number of standard errors to use as the length of the error bars.*

## **TEMPLATE Subcommand**

`TEMPLATE` uses an existing chart as a template and applies it to the chart requested by the current `GRAPH` command.

- The specification on `TEMPLATE` is a chart file saved during a previous session.
- The general rule of application is that the template overrides the default setting, but the specifications on the current `GRAPH` command override the template. Nonapplicable elements and attributes are ignored.
- Three types of elements and attributes can be applied from a chart template: those dependent on data, those dependent on the chart type, and those dependent on neither.

### ***Elements and Attributes Independent of Chart Types or Data***

Elements and attributes common to all chart types are always applied unless overridden by the specifications on the current GRAPH command.

- The title, subtitle, and footnote, including text, color, font type and size, and line alignment are always applied. To give your chart a new title, subtitle, or footnote, specify the text on the TITLE, SUBTITLE, or FOOTNOTE subcommand. You cannot change other attributes.
- The outer frame of the chart, including line style, color, and fill pattern, is always applied. The inner frame is applied except for those charts that do not have an inner frame. The template overrides the system default.
- Label formats are applied wherever applicable. The template overrides the system default. Label text, however, is not applied. GRAPH automatically provides axis labels according to the function/variable specification.
- Legends and the legend title attributes, including color, font type and size, and alignment, are applied provided the current chart requires legends. The legend title text, however, is not applied. GRAPH provides the legend title according to the function/variable specification.

### ***Elements and Attributes Dependent on Chart Type***

Elements and attributes dependent on the chart type are those that exist only in a specific chart type. They include bars (in bar charts), lines and areas (in line charts), markers (in scatterplots), boxes (in boxplots), and pie sectors (in pie charts). These elements and their attributes are usually applied only when the template chart and the requested chart are of the same type. Some elements or their attributes may override the default settings across chart type.

- Color and pattern are always applied except for pie charts. The template overrides the system default.
- Scale axis lines are applied across chart types.
- Interval axis lines are applied from interval axis to interval axis. Interval axis bins are never applied.
- If the template is a 3-D bar chart and you request a chart with one category axis, attributes of the first axis are applied from the template. If you request a 3-D bar chart and the template is not a 3-D chart, no category axis attributes are applied.

### ***Elements and Attributes Dependent on Data***

Data-dependent elements and attributes are applied only when the template and the requested chart are of the same type and the template has at least as many series assigned to the same types of chart elements as the requested chart.

- Category attributes and elements, including fill, border, color, pattern, line style, weight of pie sectors, pie sector explosion, reference lines, projection lines, and annotations, are applied only when category values in the requested chart match those in the template.
- The attributes of data-related elements with on/off states are always applied. For example, the line style, weight, and color of a quadratic fit in a simple bivariate scatterplot are applied if the requested chart is also a simple bivariate scatterplot. The specification on the GRAPH



command, for example, `HISTOGRAM(NORMAL)`, overrides the applied on/off status; in this case, a normal curve is displayed regardless of whether the template displays a normal curve.

- In bar, line, and area charts, the assignment of series to bars, lines, and areas is not applied.

## ***MISSING Subcommand***

`MISSING` controls the treatment of missing values in the chart drawn by `GRAPH`.

- The default is `LISTWISE`.
- The `MISSING` subcommand has no effect on variables used with the `VALUE` function to create nonaggregated charts. User-missing and system-missing values create empty cells.
- `LISTWISE` and `VARIABLE` are alternatives and apply to variables used in summary functions for a chart or to variables being plotted in a scatterplot.
- `REPORT` and `NOREPORT` are alternatives and apply only to category variables. They control whether categories and series with missing values are created. `NOREPORT` is the default.
- `INCLUDE` and `EXCLUDE` are alternatives and apply to both summary and category variables. `EXCLUDE` is the default.
- When a case has a missing value for the name variable but contains valid values for the dependent variable in a scatterplot, the case is always included. User-missing values are displayed as point labels; system-missing values are not displayed.
- For an aggregated categorical chart, if every aggregated series is empty in a category, the empty category is excluded.
- A nonaggregated categorical chart created with the `VALUE` function can contain completely empty categories. There are always as many categories as rows of data. However, at least one nonempty cell must be present; otherwise the chart is not created.

<b>LISTWISE</b>	<i>Listwise deletion of cases with missing values. A case with a missing value for any dependent variable is excluded from computations and graphs.</i>
<b>VARIABLE</b>	<i>Variable-wise deletion. A case is deleted from the analysis only if it has a missing value for the dependent variable being analyzed.</i>
<b>NOREPORT</b>	<i>Suppress missing-value categories. This is the default.</i>
<b>REPORT</b>	<i>Report and graph missing-value categories.</i>
<b>EXCLUDE</b>	<i>Exclude user-missing values. Both user- and system-missing values for dependent variables are excluded from computations and graphs. This is the default.</i>
<b>INCLUDE</b>	<i>Include user-missing values. Only system-missing values for dependent variables are excluded from computations and graphs.</i>

# HILOGLINEAR

HILOGLINEAR is available in the Advanced Statistics option.

```
HILOGLINEAR {varlist} (min,max) [varlist ...]
             {ALL      }

[/METHOD [= BACKWARD]]

[/MAXORDER = k]

[/CRITERIA = [CONVERGE({0.25**})] [ITERATE({20**})] [P({0.05**})]
             {n          } {n          } {prob  }
             [DELTA({0.5**})] [MAXSTEPS({10**})]
             {d          } {n          }
             [DEFAULT] ]

[/CWEIGHT = {varname }
            {(matrix)}]

[/PRINT = {[FREQ**] [RESID**] [ESTIM**] [ASSOCIATION**]}]
          {DEFAULT**      }
          {ALL             }
          {NONE            }

[/PLOT = [{NONE**          } ]
         {DEFAULT         }
         {[RESID] [NORMPROB]}
         {ALL             } ]

[/MISSING = [{EXCLUDE**}]
            {INCLUDE  } ]

[/DESIGN = effectname effectname*effectname ...]
```

\*\* Default if subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
HILOGLINEAR V1(1,2) V2(1,2)
            /DESIGN=V1*V2.
```

## Overview

HILOGLINEAR fits hierarchical loglinear models to multidimensional contingency tables using an iterative proportional-fitting algorithm. HILOGLINEAR also estimates parameters for saturated models. These techniques are described elsewhere in (Everitt, 1977), (Bishop, Feinberg, and Holland, 1975), and (Goodman, 1978). HILOGLINEAR is much more efficient for these models than the LOGLINEAR procedure because HILOGLINEAR uses an iterative proportional-fitting algorithm rather than the Newton-Raphson method used in LOGLINEAR.

### **Options**

**Design Specification.** You can request automatic model selection using backward elimination with the `METHOD` subcommand. You can also specify any hierarchical design and request multiple designs using the `DESIGN` subcommand.

**Design Control.** You can control the criteria used in the iterative proportional-fitting and model-selection routines with the `CRITERIA` subcommand. You can also limit the order of effects in the model with the `MAXORDER` subcommand and specify structural zeros for cells in the tables you analyze with the `CWEIGHT` subcommand.

**Display and Plots.** You can select the display for each design with the `PRINT` subcommand. For saturated models, you can request tests for different orders of effects as well. With the `PLOT` subcommand, you can request residuals plots or normal probability plots of residuals.

### **Basic Specification**

- The basic specification is a variable list with at least two variables followed by their minimum and maximum values.
- HILOGLINEAR estimates a saturated model for all variables in the analysis.
- By default, HILOGLINEAR displays parameter estimates, measures of partial association, goodness of fit, and frequencies for the saturated model.

### **Subcommand Order**

- The variable list must be specified first.
- Subcommands affecting a given `DESIGN` must appear before the `DESIGN` subcommand. Otherwise, subcommands can appear in any order.
- `MISSING` can be placed anywhere after the variable list.

### **Syntax Rules**

- `DESIGN` is optional. If `DESIGN` is omitted or the last specification is not a `DESIGN` subcommand, a default saturated model is estimated.
- You can specify multiple `PRINT`, `PLOT`, `CRITERIA`, `MAXORDER`, and `CWEIGHT` subcommands. The last of each type specified is in effect for subsequent designs.
- `PRINT`, `PLOT`, `CRITERIA`, `MAXORDER`, and `CWEIGHT` specifications remain in effect until they are overridden by new specifications on these subcommands.
- You can specify multiple `METHOD` subcommands, but each one affects only the next design.
- `MISSING` can be specified only once.

### **Operations**

- HILOGLINEAR builds a contingency table using all variables on the variable list. The table contains a cell for each possible combination of values within the range specified for each variable.

- HILOGLINEAR assumes that there is a category for every integer value in the range of each variable. Empty categories waste space and can cause computational problems. If there are empty categories, use the RECODE command to create consecutive integer values for categories.
- Cases with values outside the range specified for a variable are excluded.
- If the last subcommand is not a DESIGN subcommand, HILOGLINEAR displays a warning and generates the default model. This is the saturated model unless MAXORDER is specified. This model is in addition to any that are explicitly requested.
- If the model is not saturated (for example, when MAXORDER is less than the number of factors), only the goodness of fit and the observed and expected frequencies are given.
- The display uses the WIDTH subcommand defined on the SET command. If the defined width is less than 132, some portions of the display may be deleted.

### Limitations

The HILOGLINEAR procedure cannot estimate all possible frequency models, and it produces limited output for unsaturated models.

- It can estimate only hierarchical loglinear models.
- It treats all table variables as nominal. (You can use LOGLINEAR to fit nonhierarchical models to tables involving variables that are ordinal.)
- It can produce parameter estimates for saturated models only (those with all possible main-effect and interaction terms).
- It can estimate partial associations for saturated models only.
- It can handle tables with no more than 10 factors.

### Example

```
HILOGLINEAR V1(1,2) V2(1,2) V3(1,3) V4(1,3)
  /DESIGN=V1*V2*V3, V4.
```

- HILOGLINEAR builds a  $2 \times 2 \times 3 \times 3$  contingency table for analysis.
- DESIGN specifies the generating class for a hierarchical model. This model consists of main effects for all four variables, two-way interactions among  $V1$ ,  $V2$ , and  $V3$ , and the three-way interaction term  $V1$  by  $V2$  by  $V3$ .

### Backward Elimination

```
HILOGLINEAR
  inccat(1 4) news(0 1) response(0 1) /METHOD=BACKWARD
  /CRITERIA MAXSTEPS(10) P(.05) ITERATION(20) DELTA(.5)
  /PRINT=FREQ RESID
  /DESIGN .
```

- HILOGLINEAR builds a  $4 \times 2 \times 2$  contingency table for analysis.

- `METHOD` specifies that backward elimination should be used to choose a final model.
- All other options are set to their default values. The empty `DESIGN` subcommand indicates that the procedure will start with a saturated model.

## Variable List

The required variable list specifies the variables in the analysis. The variable list must precede all other subcommands.

- Variables must be numeric and have integer values. If a variable has a fractional value, the fractional portion is truncated.
- Keyword `ALL` can be used to refer to all user-defined variables in the active dataset.
- A range must be specified for each variable, with the minimum and maximum values separated by a comma and enclosed in parentheses.
- If the same range applies to several variables, the range can be specified once after the last variable to which it applies.
- If `ALL` is specified, all variables must have the same range.

## METHOD Subcommand

By default, HILOGLINEAR tests the model specified on the `DESIGN` subcommand (or the default model) and does not perform any model selection. All variables are entered and none are removed. Use `METHOD` to specify automatic model selection using backward elimination for the next design specified.

- You can specify `METHOD` alone or with the keyword `BACKWARD` for an explicit specification.
- When the backward-elimination method is requested, a step-by-step output is displayed regardless of the specification on the `PRINT` subcommand.
- `METHOD` affects only the next design.

**BACKWARD**                      *Backward elimination.* Perform backward elimination of terms in the model. All terms are entered. Those that do not meet the `P` criterion specified on the `CRITERIA` subcommand (or the default `P`) are removed one at a time.

## MAXORDER Subcommand

`MAXORDER` controls the maximum order of terms in the model estimated for subsequent designs. If `MAXORDER` is specified, HILOGLINEAR tests a model only with terms of that order or less.

- `MAXORDER` specifies the highest-order term that will be considered for the next design. `MAXORDER` can thus be used to abbreviate computations for the `BACKWARD` method.
- If the integer on `MAXORDER` is less than the number of factors, parameter estimates and measures of partial association are not available. Only the goodness of fit and the observed and expected frequencies are displayed.

- You can use `MAXORDER` with backward elimination to find the best model with terms of a certain order or less. This is computationally much more efficient than eliminating terms from the saturated model.

### Example

```
HILOGLINEAR V1 V2 V3 (1,2)
  /MAXORDER=2
  /DESIGN=V1 V2 V3
  /DESIGN=V1*V2*V3.
```

- `HILOGLINEAR` builds a  $2 \times 2 \times 2$  contingency table for  $V1$ ,  $V2$ , and  $V3$ .
- `MAXORDER` has no effect on the first `DESIGN` subcommand because the design requested considers only main effects.
- `MAXORDER` restricts the terms in the model specified on the second `DESIGN` subcommand to two-way interactions and main effects.

## CRITERIA Subcommand

Use the `CRITERIA` subcommand to change the values of constants in the iterative proportional-fitting and model-selection routines for subsequent designs.

- The default criteria are in effect if the `CRITERIA` subcommand is omitted (see below).
- You cannot specify the `CRITERIA` subcommand without any keywords.
- Specify each `CRITERIA` keyword followed by a criterion value in parentheses. Only those criteria specifically altered are changed.
- You can specify more than one keyword on `CRITERIA`, and they can be in any order.

<b>DEFAULT</b>	<i>Reset parameters to their default values.</i> If you have specified criteria other than the defaults for a design, use this keyword to restore the defaults for subsequent designs.
<b>CONVERGE(n)</b>	<i>Convergence criterion.</i> The default is $10^{-3}$ times the largest cell size, or 0.25, whichever is larger.
<b>ITERATE(n)</b>	<i>Maximum number of iterations.</i> The default is 20.
<b>P(n)</b>	<i>Probability for change in chi-square if term is removed.</i> Specify a value between (but not including) 0 and 1 for the significance level. The default is 0.05. <code>P</code> is in effect only when you request <code>BACKWARD</code> on the <code>METHOD</code> subcommand.
<b>MAXSTEPS(n)</b>	<i>Maximum number of steps for model selection.</i> Specify an integer between 1 and 99, inclusive. The default is 10.
<b>DELTA(d)</b>	<i>Cell delta value.</i> The value of delta is added to each cell frequency for the first iteration when estimating saturated models; it is ignored for unsaturated models. The default value is 0.5. You can specify any decimal value between 0 and 1 for $d$ . <code>HILOGLINEAR</code> does not display parameter estimates or the covariance matrix of parameter estimates if any zero cells (either structural or sampling) exist in the expected table after delta is added.

## ***CWEIGHT Subcommand***

**CWEIGHT** specifies cell weights for a model. **CWEIGHT** is typically used to specify structural zeros in the table. You can also use **CWEIGHT** to adjust tables to fit new margins.

- You can specify the name of a variable whose values are cell weights, or provide a matrix of cell weights enclosed in parentheses.
- If you use a variable to specify cell weights, you are allowed only one **CWEIGHT** subcommand.
- If you specify a matrix, you must provide a weight for every cell in the contingency table, where the number of cells equals the product of the number of values of all variables.
- Cell weights are indexed by the values of the variables in the order in which they are specified on the variable list. The index values of the rightmost variable change the most quickly.
- You can use the notation  $n*cw$  to indicate that cell weight  $cw$  is repeated  $n$  times in the matrix.

### ***Example***

```
HILOGLINEAR V1(1,2) V2(1,2) V3(1,3)
/CWEIGHT=CELLWGT
/DESIGN=V1*V2, V2*V3, V1*V3.
```

- This example uses the variable *CELLWGT* to assign cell weights for the table. Only one **CWEIGHT** subcommand is allowed.

### ***Example***

```
HILOGLINEAR V4(1,3) V5(1,3)
/CWEIGHT=(0 1 1 1 0 1 1 1 0)
/DESIGN=V4, V5.
```

- The **HILOGLINEAR** command sets the diagonal cells in the model to structural zeros. This type of model is known as a **quasi-independence model**.
- Because both *V4* and *V5* have three values, weights must be specified for nine cells.
- The first cell weight is applied to the cell in which *V4* is 1 and *V5* is 1; the second weight is applied to the cell in which *V4* is 1 and *V5* is 2; and so on.

### ***Example***

```
HILOGLINEAR V4(1,3) V5(1,3)
/CWEIGHT=(0 3*1 0 3*1 0)
/DESIGN=V4,V5.
```

- This example is the same as the previous example except that the  $n*cw$  notation is used.

### ***Example***

\* An Incomplete Rectangular Table

```
DATA LIST FREE / LOCULAR RADIAL FREQ.
WEIGHT BY FREQ.
BEGIN DATA
1 1 462
1 2 130
1 3 2
```

*HILOGLINEAR*

```

1 4 1
2 1 103
2 2 35
2 3 1
2 4 0
3 5 614
3 6 138
3 7 21
3 8 14
3 9 1
4 5 443
4 6 95
4 7 22
4 8 8
4 9 5
END DATA.
HILOGLINEAR LOCULAR (1,4) RADIAL (1,9)
/CWEIGHT=(4*1 5*0 4*1 5*0 4*0 5*1 4*0 5*1)
/DESIGN LOCULAR RADIAL.

```

- This example uses aggregated table data as input.
- The `DATA LIST` command defines three variables. The values of *LOCULAR* and *RADIAL* index the levels of those variables, so that each case defines a cell in the table. The values of *FREQ* are the cell frequencies.
- The `WEIGHT` command weights each case by the value of the variable *FREQ*. Because each case represents a cell in this example, the `WEIGHT` command assigns the frequencies for each cell.
- The `BEGIN DATA` and `END DATA` commands enclose the inline data.
- The `HILOGLINEAR` variable list specifies two variables. *LOCULAR* has values 1, 2, 3, and 4. *RADIAL* has integer values 1 through 9.
- The `CWEIGHT` subcommand identifies a block rectangular pattern of cells that are logically empty. There is one weight specified for each cell of the 36-cell table.
- In this example, the matrix form needs to be used in `CWEIGHT` because the structural zeros do not appear in the actual data. (For example, there is no case corresponding to *LOCULAR* = 1, *RADIAL* = 5.)
- The `DESIGN` subcommand specifies main effects only for *LOCULAR* and *RADIAL*. Lack of fit for this model indicates an interaction of the two variables.
- Because there is no `PRINT` or `PLOT` subcommand, `HILOGLINEAR` produces the default output for an unsaturated model.

***PRINT Subcommand***

`PRINT` controls the display produced for the subsequent designs.

- If `PRINT` is omitted or included with no specifications, the default display is produced.
- If any keywords are specified on `PRINT`, only output specifically requested is displayed.
- `HILOGLINEAR` displays Pearson and likelihood-ratio chi-square goodness-of-fit tests for models. For saturated models, it also provides tests that the *k*-way effects and the *k*-way and higher-order effects are 0.



- Both adjusted and unadjusted degrees of freedom are displayed for tables with sampling or structural zeros. *K*-way and higher-order tests use the unadjusted degrees of freedom.
- The unadjusted degrees of freedom are not adjusted for zero cells, and they estimate the upper bound of the true degrees of freedom. These are the same degrees of freedom you would get if all cells were filled.
- The adjusted degrees of freedom are calculated from the number of non-zero-fitted cells minus the number of parameters that would be estimated if all cells were filled (that is, unadjusted degrees of freedom minus the number of zero-fitted cells). This estimate of degrees of freedom may be too low if some parameters do not exist because of zeros.

<b>DEFAULT</b>	<i>Default displays.</i> This option includes <code>FREQ</code> and <code>RESID</code> output for nonsaturated models, and <code>FREQ</code> , <code>RESID</code> , <code>ESTIM</code> , and <code>ASSOCIATION</code> output for saturated models. For saturated models, the observed and expected frequencies are equal, and the residuals are zeros.
<b>FREQ</b>	<i>Observed and expected cell frequencies.</i>
<b>RESID</b>	<i>Raw and standardized residuals.</i>
<b>ESTIM</b>	<i>Parameter estimates for a saturated model.</i>
<b>ASSOCIATION</b>	<i>Partial associations.</i> You can request partial associations of effects only when you specify a saturated model. This option is computationally expensive for tables with many factors.
<b>ALL</b>	<i>All available output.</i>
<b>NONE</b>	<i>Design information and goodness-of-fit statistics only.</i> Use of this option overrides all other specifications on <code>PRINT</code> .

## ***PLOT Subcommand***

Use `PLOT` to request residuals plots.

- If `PLOT` is included without specifications, standardized residuals and normal probability plots are produced.
- No plots are displayed for saturated models.
- If `PLOT` is omitted, no plots are produced.

<b>RESID</b>	<i>Standardized residuals by observed and expected counts.</i>
<b>NORMPLOT</b>	<i>Normal probability plots of adjusted residuals.</i>
<b>NONE</b>	<i>No plots.</i> Specify <code>NONE</code> to suppress plots requested on a previous <code>PLOT</code> subcommand. This is the default if <code>PLOT</code> is omitted.
<b>DEFAULT</b>	<i>Default plots.</i> Includes <code>RESID</code> and <code>NORMPLOT</code> . This is the default when <code>PLOT</code> is specified without keywords.
<b>ALL</b>	<i>All available plots.</i>

## ***MISSING Subcommand***

By default, a case with either system-missing or user-missing values for any variable named on the `HILOGLINEAR` variable list is omitted from the analysis. Use `MISSING` to change the treatment of cases with user-missing values.

- `MISSING` can be named only once and can be placed anywhere following the variable list.

- MISSING cannot be used without specifications.
  - A case with a system-missing value for any variable named on the variable list is always excluded from the analysis.
- EXCLUDE**                      *Delete cases with missing values.* This is the default if the subcommand is omitted. You can also specify keyword `DEFAULT`.
- INCLUDE**                      *Include user-missing values as valid.* Only cases with system-missing values are deleted.

## ***DESIGN Subcommand***

By default, HILOGLINEAR uses a saturated model that includes all variables on the variable list. The model contains all main effects and interactions for those variables. Use `DESIGN` to specify a different generating class for the model.

- If `DESIGN` is omitted or included without specifications, the default model is estimated. When `DESIGN` is omitted, a warning message is issued.
- To specify a design, list the highest-order terms, using variable names and asterisks (\*) to indicate interaction effects.
- In a hierarchical model, higher-order interaction effects imply lower-order interaction and main effects.  $V1*V2*V3$  implies the three-way interaction  $V1$  by  $V2$  by  $V3$ , two-way interactions  $V1$  by  $V2$ ,  $V1$  by  $V3$ , and  $V2$  by  $V3$ , and main effects for  $V1$ ,  $V2$ , and  $V3$ . The highest-order effects to be estimated are the generating class.
- Any `PRINT`, `PLOT`, `CRITERIA`, `METHOD`, and `MAXORDER` subcommands that apply to a `DESIGN` subcommand must appear before it.
- All variables named on `DESIGN` must be named or implied on the variable list.
- You can specify more than one `DESIGN` subcommand. One model is estimated for each `DESIGN` subcommand.
- If the last subcommand on HILOGLINEAR is not `DESIGN`, the default model will be estimated in addition to models explicitly requested. A warning message is issued for a missing `DESIGN` subcommand.

## ***References***

- Bishop, Y. M., S. E. Feinberg, and P. W. Holland. 1975. *Discrete multivariate analysis: Theory and practice*. Cambridge, Mass.: MIT Press.
- Everitt, B. S. 1977. *The Analysis of Contingency Tables*. London: Chapman & Hall.
- Goodman, L. A. 1978. *Analyzing qualitative/categorical data*. New York: University Press of America.

# HOMALS

HOMALS is available in the Categories option.

```
HOMALS VARIABLES=varlist(max)

[/ANALYSIS=varlist]

[/NOBSERVATIONS=value]

[/DIMENSION={2** }]
           {value}

[/MAXITER={100**}]
           {value}

[/CONVERGENCE={.00001**}]
           {value }

[/PRINT={DEFAULT**} [FREQ**] [EIGEN**] [DISCRIM**]
        [QUANT**] [OBJECT] [HISTORY] [ALL] [NONE]]

[/PLOT=[NDIM=({1, 2 }**)]
       {value, value}
       {ALL, MAX }
       [QUANT**[(varlist)][(n)]] [OBJECT**[(varlist)][(n)]]
       [DEFAULT**[(n)]] [DISCRIM[(n)]] [ALL[(n)]] [NONE]]

[/SAVE=[rootname] [(value)]]

[/MATRIX=OUT({* })]
           {'savfile'|'dataset'}
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
HOMALS VARIABLES=ACOLA(2) BCOLA(2) CCOLA(2) DCOLA(2).
```

## Overview

HOMALS (*homogeneity analysis by means of alternating least squares*) estimates category quantifications, object scores, and other associated statistics that separate categories (levels) of nominal variables as much as possible and divide cases into homogeneous subgroups.

### Options

**Data and variable selection.** You can use a subset of the variables in the analysis and restrict the analysis to the first  $n$  observations.

**Number of dimensions.** You can specify the number of dimensions HOMALS should compute.

**Iterations and convergence.** You can specify the maximum number of iterations and the value of a convergence criterion.

**Display output.** The output can include all available statistics; just the default frequencies, eigenvalues, discrimination measures and category quantifications; or just the specific statistics you request. You can also control which statistics are plotted and specify the number of characters used in plot labels.

**Saving scores.** You can save object scores in the working data file.

**Writing matrices.** You can write a matrix data file containing category quantifications for use in further analyses.

### ***Basic Specification***

- The basic specification is `HOMALS` and the `VARIABLES` subcommand. By default, `HOMALS` analyzes all of the variables listed for all cases and computes two solutions. Frequencies, eigenvalues, discrimination measures, and category quantifications are displayed, and category quantifications and object scores are plotted.

### ***Subcommand Order***

- Subcommands can appear in any order.

### ***Syntax Rules***

- If `ANALYSIS` is specified more than once, `HOMALS` is not executed. For all other subcommands, if a subcommand is specified more than once, only the last occurrence is executed.

### ***Operations***

- `HOMALS` treats every value in the range of 1 to the maximum value specified on `VARIABLES` as a valid category. If the data are not sequential, the empty categories (categories with no valid data) are assigned zeros for all statistics. You may want to use `RECODE` or `AUTORECODE` before `HOMALS` to get rid of these empty categories and avoid the unnecessary output (see `RECODE` and `AUTORECODE` for more information).

### ***Limitations***

- String variables are not allowed; use `AUTORECODE` to recode string variables into numeric variables.
- The data (category values) must be positive integers. Zeros and negative values are treated as system-missing, which means that they are excluded from the analysis. Fractional values are truncated after the decimal and are included in the analysis. If one of the levels of a variable has been coded 0 or a negative value and you want to treat it as a valid category, use the `AUTORECODE` or `RECODE` command to recode the values of that variable.
- `HOMALS` ignores user-missing value specifications. Positive user-missing values less than the maximum value specified on the `VARIABLES` subcommand are treated as valid category values and are included in the analysis. If you do not want the category included, use `COMPUTE` or `RECODE` to change the value to something outside of the valid range. Values outside of the range (less than 1 or greater than the maximum value) are treated as system-missing and are excluded from the analysis.

## Example

```
HOMALS VARIABLES=ACOLA(2) BCOLA(2) CCOLA(2) DCOLA(2)
/PRINT=FREQ EIGEN QUANT OBJECT.
```

- The four variables are analyzed using all available observations. Each variable has two categories, 1 and 2.
- The PRINT subcommand lists the frequencies, eigenvalues, category quantifications, and object scores.
- By default, plots of the category quantifications and the object scores are produced.

## VARIABLES Subcommand

VARIABLES specifies the variables that will be used in the analysis.

- The VARIABLES subcommand is required. The actual word VARIABLES can be omitted.
- After each variable or variable list, specify in parentheses the maximum number of categories (levels) of the variables.
- The number specified in parentheses indicates the number of categories *and* the maximum category value. For example, *VARI(3)* indicates that *VARI* has three categories coded 1, 2, and 3. However, if a variable is not coded with consecutive integers, the number of categories used in the analysis will differ from the number of observed categories. For example, if a three-category variable is coded {2, 4, 6}, the maximum category value is 6. The analysis treats the variable as having six categories, three of which (categories 1, 3, and 5) are not observed and receive quantifications of 0.
- To avoid unnecessary output, use the AUTORECODE or RECODE command before HOMALS to recode a variable that does not have sequential values (see AUTORECODE and RECODE for more information).

### Example

```
DATA LIST FREE/V1 V2 V3.
BEGIN DATA
3 1 1
6 1 1
3 1 3
3 2 2
3 2 2
6 2 2
6 1 3
6 2 2
3 2 2
6 2 1
END DATA.
AUTORECODE V1 /INTO NEWVAR1.
HOMALS VARIABLES=NEWVAR1 V2(2) V3(3).
```

- DATA LIST defines three variables, *V1*, *V2*, and *V3*.
- *V1* has two levels, coded 3 and 6, *V2* has two levels, coded 1 and 2, and *V3* has three levels, coded 1, 2, and 3.

- The `AUTORECODE` command creates *NEWVARI* containing recoded values of *V1*. Values of 3 are recoded to 1; values of 6 are recoded to 2.
- The maximum category value for both *NEWVARI* and *V2* is 2. A maximum value of 3 is specified for *V3*.

## ***ANALYSIS Subcommand***

`ANALYSIS` limits the analysis to a specific subset of the variables named on the `VARIABLES` subcommand.

- If `ANALYSIS` is not specified, all variables listed on the `VARIABLES` subcommand are used.
- `ANALYSIS` is followed by a variable list. The variables on the list must be specified on the `VARIABLES` subcommand.
- Variables listed on the `VARIABLES` subcommand but not on the `ANALYSIS` subcommand can still be used to label object scores on the `PLOT` subcommand.

### ***Example***

```
HOMALS VARIABLES=ACOLA(2) BCOLA(2) CCOLA(2) DCOLA(2)
/ANALYSIS=ACOLA BCOLA
/PRINT=OBJECT QUANT
/PLOT=OBJECT(CCOLA) .
```

- The `VARIABLES` subcommand specifies four variables.
- The `ANALYSIS` subcommand limits analysis to the first two variables. The `PRINT` subcommand lists the object scores and category quantifications from this analysis.
- The plot of the object scores is labeled with variable *CCOLA*, even though this variable is not included in the computations.

## ***NOBSERVATIONS Subcommand***

`NOBSERVATIONS` specifies how many cases are used in the analysis.

- If `NOBSERVATIONS` is not specified, all available observations in the working data file are used.
- `NOBSERVATIONS` is followed by an integer indicating that the first *n* cases are to be used.

## ***DIMENSION Subcommand***

`DIMENSION` specifies the number of dimensions you want `HOMALS` to compute.

- If you do not specify the `DIMENSION` subcommand, `HOMALS` computes two dimensions.
- The specification on `DIMENSION` is a positive integer indicating the number of dimensions.

- The minimum number of dimensions is 1.
- The maximum number of dimensions is equal to the smaller of the two values below:

### **MAXITER Subcommand**

MAXITER specifies the maximum number of iterations HOMALS can go through in its computations.

- If MAXITER is not specified, HOMALS will iterate up to 100 times.
- The specification on MAXITER is a positive integer indicating the maximum number of iterations.

### **CONVERGENCE Subcommand**

CONVERGENCE specifies a convergence criterion value. HOMALS stops iterating if the difference in total fit between the last two iterations is less than the CONVERGENCE value.

- If CONVERGENCE is not specified, the default value is 0.00001.
- The specification on CONVERGENCE is a positive value.

### **PRINT Subcommand**

PRINT controls which statistics are included in your display output. The default display includes the frequencies, eigenvalues, discrimination measures, and category quantifications.

The following keywords are available:

<b>FREQ</b>	<i>Marginal frequencies for the variables in the analysis.</i>
<b>HISTORY</b>	<i>History of the iterations.</i>
<b>EIGEN</b>	<i>Eigenvalues.</i>
<b>DISCRIM</b>	<i>Discrimination measures for the variables in the analysis.</i>
<b>OBJECT</b>	<i>Object scores.</i>
<b>QUANT</b>	<i>Category quantifications for the variables in the analysis.</i>
<b>DEFAULT</b>	<i>FREQ, EIGEN, DISCRIM, and QUANT. These statistics are also displayed when you omit the PRINT subcommand.</i>
<b>ALL</b>	<i>All available statistics.</i>
<b>NONE</b>	<i>No statistics.</i>

### **PLOT Subcommand**

PLOT can be used to produce plots of category quantifications, object scores, and discrimination measures.

- If PLOT is not specified, plots of the object scores and of the quantifications are produced.
- No plots are produced for a one-dimensional solution.

The following keywords can be specified on PLOT:

<b>DISCRIM</b>	<i>Plots of the discrimination measures.</i>
<b>OBJECT</b>	<i>Plots of the object scores.</i>
<b>QUANT</b>	<i>Plots of the category quantifications.</i>
<b>DEFAULT</b>	<i>QUANT and OBJECT.</i>
<b>ALL</b>	<i>All available plots.</i>
<b>NONE</b>	<i>No plots.</i>

- Keywords OBJECT and QUANT can each be followed by a variable list in parentheses to indicate that plots should be labeled with those variables. For QUANT, the labeling variables must be specified on both the VARIABLES and ANALYSIS subcommands. For OBJECT, the variables must be specified on the VARIABLES subcommand but need not appear on the ANALYSIS subcommand. This means that variables not used in the computations can be used to label OBJECT plots. If the variable list is omitted, the default object and quantification plots are produced.
- Object score plots labeled with variables that appear on the ANALYSIS subcommand use category labels corresponding to all categories within the defined range. Objects in a category that is outside the defined range are labeled with the label corresponding to the category immediately following the defined maximum category value.
- Object score plots labeled with variables not included on the ANALYSIS subcommand use all category labels, regardless of whether or not the category value is inside the defined range.
- All keywords except NONE can be followed by an integer value in parentheses to indicate how many characters of the variable or value label are to be used on the plot. (If you specify a variable list after OBJECT or QUANT, specify the value in parentheses after the list.) The value can range from 1 to 20; the default is to use 12 characters. Spaces between words count as characters.
- DISCRIM plots use variable labels; all other plots use value labels.
- If a variable label is not supplied, the variable name is used for that variable. If a value label is not supplied, the actual value is used.
- Variable and value labels should be unique.
- When points overlap, the points involved are described in a summary following the plot.

### **Example**

```
HOMALS VARIABLES COLA1 (4) COLA2 (4) COLA3 (4) COLA4 (2)
/ANALYSIS COLA1 COLA2 COLA3 COLA4
/PLOT OBJECT(COLA4) .
```

- Four variables are included in the analysis.
- OBJECT requests a plot of the object scores labeled with the values of COLA4. Any object whose COLA4 value is not 1 or 2, is labeled 3 (or the value label for category 3, if supplied).

### **Example**

```
HOMALS VARIABLES COLA1 (4) COLA2 (4) COLA3 (4) COLA4 (2)
/ANALYSIS COLA1 COLA2 COLA3
/PLOT OBJECT(COLA4) .
```



- Three variables are included in the analysis.
- OBJECT requests a plot of the object scores labeled with the values of COLA4, a variable not included in the analysis. Objects are labeled using all values of COLA4.

In addition to the plot keywords, the following can be specified:

**NDIM**      *Dimension pairs to be plotted.* NDIM is followed by a pair of values in parentheses. If NDIM is not specified, plots are produced for dimension 1 versus dimension 2.

- The first value indicates the dimension that is plotted against all higher dimensions. This value can be any integer from 1 to the number of dimensions minus 1.
- The second value indicates the highest dimension to be used in plotting the dimension pairs. This value can be any integer from 2 to the number of dimensions.
- Keyword ALL can be used instead of the first value to indicate that all dimensions are paired with higher dimensions.
- Keyword MAX can be used instead of the second value to indicate that plots should be produced up to and including the highest dimension fit by the procedure.

### **Example**

```
HOMALS COLA1 COLA2 COLA3 COLA4 (4)
/PLOT NDIM(1,3) QUANT(5) .
```

- The NDIM(1,3) specification indicates that plots should be produced for two dimension pairs—dimension 1 versus dimension 2 and dimension 1 versus dimension 3.
- QUANT requests plots of the category quantifications. The (5) specification indicates that the first five characters of the value labels are to be used on the plots.

### **Example**

```
HOMALS COLA1 COLA2 COLA3 COLA4 (4)
/PLOT NDIM(ALL,3) QUANT(5) .
```

- This plot is the same as above except for the ALL specification following NDIM. This indicates that all possible pairs up to the second value should be plotted, so QUANT plots will be produced for dimension 1 versus dimension 2, dimension 2 versus dimension 3, and dimension 1 versus dimension 3.

## **SAVE Subcommand**

SAVE lets you add variables containing the object scores computed by HOMALS to the working data file.

- If SAVE is not specified, object scores are not added to the working data file.
- A variable rootname can be specified on the SAVE subcommand to which HOMALS adds the number of the dimension. Only one rootname can be specified and it can contain up to six characters.

- If a rootname is not specified, unique variable names are automatically generated. The variable names are *HOM<sub>n</sub>\_m*, where *n* is a dimension number and *m* is a set number. If three dimensions are saved, the first set of names is *HOM1\_1*, *HOM2\_1*, and *HOM3\_1*. If another HOMALS is then run, the variable names for the second set are *HOM1\_2*, *HOM2\_2*, *HOM3\_2*, and so on.
- Following the rootname, the number of dimensions for which you want to save object scores can be specified in parentheses. The number cannot exceed the value on the DIMENSION subcommand.
- If the number of dimensions is not specified, the SAVE subcommand saves object scores for all dimensions.
- If you replace the working data file by specifying an asterisk (\*) on a MATRIX subcommand, the SAVE subcommand is not executed.

### Example

```
HOMALS CAR1 CAR2 CAR3 CAR4 (5)
  /DIMENSION=3
  /SAVE=DIM(2) .
```

- Four variables, each with five categories, are analyzed.
- The DIMENSION subcommand specifies that results for three dimensions will be computed.
- SAVE adds the object scores from the first two dimensions to the working data file. The names of these new variables will be *DIM00001* and *DIM00002*, respectively.

## MATRIX Subcommand

The MATRIX subcommand is used to write category quantifications to a matrix data file or a previously declared dataset name (DATASET DECLARE command).

- The specification on MATRIX is keyword OUT and a quoted file specification of dataset name, enclosed in parentheses.
- You can specify an asterisk (\*) replace the active dataset.
- The matrix data file has one case for each value of each original variable.

The variables of the matrix data file and their values are:

<b>ROWTYPE_</b>	<i>String variable containing value QUANT for all cases.</i>
<b>LEVEL</b>	<i>String variable LEVEL containing the values (or value labels if present) of each original variable.</i>
<b>VARNAME_</b>	<i>String variable containing the original variable names.</i>
<b>DIM1...DIMn</b>	<i>Numeric variable containing the category quantifications for each dimension. Each variable is labeled DIMn, where n represents the dimension number.</i>

# HOST

*Note:* Square brackets used in the `HOST` syntax chart are required parts of the syntax and are not used to indicate optional elements. Equals signs (=) used in the syntax chart are required elements.

```
HOST COMMAND=['command' 'command'...'command']  
TIMELIMIT=n.
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 13.0

- Command introduced.

## **Example**

```
HOST COMMAND=['dir c:\myfiles\*.sav'].
```

## **Overview**

The `HOST` command executes external commands at the operating system level. For a Windows operating system, for example, this is equivalent to running commands from a command prompt in a command window.

- No output is displayed in a command window. Output is either displayed in the Viewer or redirected as specified in the operating system command.
- Standard output is either displayed in a text object in the Viewer window or redirected as specified in the operating system command.
- Standard errors are displayed as text objects in the Viewer.
- Commands that return a prompt for user input result in an EOF condition without waiting for any user input (unless input has been redirected to read from a file).
- A command that generates an error condition terminates the `HOST` command, and no subsequent commands specified on the `HOST` command are executed.
- The `HOST` command runs synchronously. Commands that launch applications result in the suspension of further IBM® SPSS® Statistics processing until the application finishes execution, unless you also specify a time limit (see keyword `TIMELIMIT` on p. 958). For example, in Windows operating systems, if a file extension is associated with an application, simply specifying a file name and extension on the command line will launch the associated application, and no further commands will be executed until the application is closed.

- The `HOST` command starts in the current working directory. By default, the initial working directory is the installation directory.
- In distributed analysis mode (available with SPSS Statistics Server), file paths in command specifications are relative to the remote server.

## Syntax

The minimum specification is the command name `HOST`, followed by the keyword `COMMAND`, an equals sign (=), and one or more operating system level commands, each enclosed in quotes, with the entire set of commands enclosed in square brackets.

### Example

```
HOST COMMAND=['dir c:\myfiles\*.sav'
             'dir c:\myfiles\*.sps > c:\myfiles\command_files.txt'
             'copy c:\myfiles\file1.txt > c:\myfiles\file2.txt'
             'dur c:\myfiles\*.xml > c:\myfiles\xmlfiles.txt'
             'c:\myfiles\myjobs\report.bat'].
```

- The directory listing for all `.sav` files is displayed in a text output object in the Viewer window.
- The directory listing for `.sps` files is redirected to a text file; so no output is displayed in the Viewer window.
- If `file2.txt` does not already exist, the `copy` command will copy the contents of `file1.txt` to a new file called `file2.txt`. If `file2.txt` exists, the `copy` command will not be executed since this would result in a user prompt asking for the user to confirm overwriting the file.
- The invalid `dur` command generates an error, which is displayed in the Viewer, and no output for that command is redirected to specified text file.
- The error condition caused by the invalid `dur` command terminates the `HOST` command, and `report.bat` is not run.

## Quoted Strings

If the command at the operating system level uses quoted strings, the standard rules for quoted strings within quoted strings apply. In general, use double-quotes to enclose a string that includes a string enclosed in single quotes, and vice-versa. [For more information, see the topic String Values in Command Specifications on p. 40.](#)

## TIMELIMIT Keyword

The optional `TIMELIMIT` keyword sets a time limit in seconds for execution of the bracketed list of commands. Fractional time values are rounded to the nearest integer.

### Example

```
HOST COMMAND=['c:\myfiles\report.bat'] TIMELIMIT=10.
```

## Using *TIMELIMIT* to Return Control

Since the `HOST` command runs synchronously, commands that launch applications result in the suspension of further IBM® SPSS® Statistics processing until the application finishes execution. That means that any commands that follow the `HOST` command will not be executed until any applications launched by the command are closed.

### Example

```
OMS /DESTINATION FORMAT=HTML
  OUTFILE='c:\temp\temp.htm'.
FREQUENCIES VARIABLES=ALL.
OMSEND.
HOST COMMAND=['c:\temp\temp.htm'].
DESCRIPTIVES VARIABLES=ALL.
```

- On Windows operating systems, if the *.htm* extension is associated with an application (typically Internet Explorer), the `HOST` command in this example will launch the associated application.
- In the absence of a `TIMELIMIT` specification, the subsequent `DESCRIPTIVES` command will not be executed until the application launched by the `HOST` command is closed.

To make sure control is automatically returned to SPSS Statistics and subsequent commands are executed, include a `TIMELIMIT` value, as in:

```
OMS /DESTINATION FORMAT=HTML
  OUTFILE='c:\temp\temp.htm'.
FREQUENCIES VARIABLES=ALL.
OMSEND.
HOST COMMAND=['c:\temp\temp.htm'] TIMELIMIT=5.
DESCRIPTIVES VARIABLES=ALL.
```

## Working Directory

The `HOST` command starts in the current working directory. By default, the initial working directory is the installation directory. So, for example, `HOST COMMAND=['dir']` executed at the start of a session would typically return a directory listing of the installation directory. The working directory can be changed, however, by the `CD` command and the `CD` keyword of the `INSERT` command.

### Example

```
*start of session.
HOST COMMAND=['dir']. /*lists contents of install directory.
CD 'c:\temp'.
HOST COMMAND=['dir']. /*lists contents of c:\temp directory.
```

### ***UNC Paths on Windows Operating Systems***

To start in the IBM® SPSS® Statistics working directory, the `HOST` command actually issues an OS-level `CD` command that specifies the SPSS Statistics working directory. On Windows operating systems, if you use UNC path specifications of the general form:

```
\\servername\sharename\path
```

on SPSS Statistics commands such as `CD` or `INSERT` to set the working directory location, the `HOST` command will fail because UNC paths are not valid on the Windows `CD` command.

#### ***Example***

```
INSERT FILE='\\hqserver\public\report.sps' CD=YES.  
HOST ['dir'].
```

- The `INSERT` command uses a UNC path specification, and `CD=YES` makes that directory the working directory.
- The subsequent `HOST` command will generate an OS-level error message that says the current directory path is invalid because UNC paths are not supported.

# IF

```
IF [(logical expression)] target variable=expression
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

*The following relational operators can be used in logical expressions:*

Symbol	Definition
EQ or =	Equal to
NE or $\neq$ or $\neq$ or $\langle \rangle$	Not equal to
LT or <	Less than
LE or $\leq$	Less than or equal to
GT or >	Greater than
GE or $\geq$	Greater than or equal to

*The following logical operators can be used in logical expressions:*

Symbol	Definition
AND or &	Both relations must be true
OR or	Either relation can be true
NOT	Reverses the outcome of an expression

### Example

```
IF (AGE > 20 AND SEX = 1) GROUP=2.
```

## Overview

IF conditionally executes a single transformation command based upon logical conditions found in the data. The transformation can create a new variable or modify the values of an existing variable for each case in the active dataset. You can create or modify the values of both numeric and string variables. If you create a new string variable, you must first declare it on the `STRING` command.

IF has three components: a *logical expression* that sets up the logical criteria, a *target variable* (the one to be modified or created), and an *assignment expression*. The target variable's values are modified according to the assignment expression.

IF is most efficient when used to execute a single, conditional, `COMPUTE`-like transformation. If you need multiple IF statements to define the condition, it is usually more efficient to use the `RECODE` command or a `DO IF-END IF` structure.

### Basic Specification

The basic specification is a logical expression followed by a target variable, a required equals sign, and the assignment expression. The assignment is executed only if the logical expression is true.

**Syntax Rules**

- Logical expressions can be simple logical variables or relations, or complex logical tests involving variables, constants, functions, relational operators, and logical operators. Both the logical expression and the assignment expression can use any of the numeric or string functions allowed in `COMPUTE` transformations.
- Parentheses can be used to enclose the logical expression. Parentheses can also be used within the logical expression to specify the order of operations. Extra blanks or parentheses can be used to make the expression easier to read.
- A relation can compare variables, constants, or more complicated arithmetic expressions. Relations cannot be abbreviated. For example, `(A EQ 2 OR A EQ 5)` is valid, while `(A EQ 2 OR 5)` is not. Blanks (not commas) must be used to separate relational operators from the expressions being compared.
- A relation cannot compare a string variable to a numeric value or variable, or vice versa. A relation cannot compare the result of the logical functions `SYSMIS`, `MISSING`, `ANY`, or `RANGE` to a number.
- String values used in expressions must be specified in quotes and must include any leading or trailing blanks. Lowercase letters are considered distinct from uppercase letters.
- String variables that are used as target variables must already exist. To declare a new string variable, first create the variable with the `STRING` command and then specify the new variable as the target variable on `IF`.

**Examples****IF with Numeric Values**

```
IF (AGE > 20 AND SEX = 1) GROUP=2.
```

- The numeric variable `GROUP` is set to 2 for cases where `AGE` is greater than 20 and `SEX` is equal to 1.
- When the expression is false or missing, the value of `GROUP` remains unchanged. If `GROUP` has not been previously defined, it contains the system-missing value.

**IF with String Values**

```
IF (SEX EQ 'F') EEO=QUOTA+GAIN.
```

- The logical expression tests the string variable `SEX` for the value `F`.
- When the expression is true (when `SEX` equals `F`), the value of the numeric variable `EEO` is assigned the value of `QUOTA` plus `GAIN`. Both `QUOTA` and `GAIN` must be previously defined numeric variables.
- When the expression is false or missing (for example, if `SEX` equals `F`), the value of `EEO` remains unchanged. If `EEO` has not been previously defined, it contains the system-missing value.



### Conditional Expressions with Arithmetic Operations

```
COMPUTE V3=0.
IF ((V1-V2) LE 7) V3=V1**2.
```

- COMPUTE assigns *V3* the value 0.
- The logical expression tests whether *V1* minus *V2* is less than or equal to 7. If it is, the value of *V3* is assigned the value of *V1* squared. Otherwise, the value of *V3* remains at 0.

### Conditional Expressions with Arithmetic Operations and Functions

```
IF (ABS(A-C) LT 100) INT=100.
```

- IF tests whether the absolute value of the variable *A* minus the variable *C* is less than 100. If it is, *INT* is assigned the value 100. Otherwise, the value is unchanged. If *INT* has not been previously defined, it is system-missing.

### Testing for Missing Values

\* Test for listwise deletion of missing values.

```
DATA LIST /V1 TO V6 1-6.
STRING SELECT(A1).
COMPUTE SELECT='V'.
VECTOR V=V1 TO V6.

LOOP #I=1 TO 6.
IF MISSING(V(#I)) SELECT='M'.
END LOOP.

BEGIN DATA
123456
   56
1 3456
123456
123456
END DATA.

FREQUENCIES VAR=SELECT.
```

- STRING creates the string variable *SELECT* with an A1 format and COMPUTE sets the value of *SELECT* to *V*.
- VECTOR defines the vector *V* as the original variables *V1* to *V6*. Variables on a single vector must be all numeric or all string variables. In this example, because the vector *V* is used as an argument on the MISSING function of IF, the variables must be numeric (MISSING is not available for string variables).
- The loop structure executes six times: once for each VECTOR element. If a value is missing for any element, SELECT is set equal to *M*. In effect, if any case has a missing value for any of the variables *V1* to *V6*, SELECT is set to *M*.
- FREQUENCIES generates a frequency table for SELECT. The table gives a count of how many cases have missing values for at least one variable and how many cases have valid values for all variables. This table can be used to determine how many cases would be dropped from an analysis that uses listwise deletion of missing values.

**Example**

```
IF YRHIRED LT 1980 RATE=0.02.
IF DEPT='SALES' DIVISION='TRANSFERRED'.
```

- The logical expression on the first IF command tests whether *YRHIRED* is less than 1980 (hired before 1980). If so, the variable *RATE* is set to 0.02.
- The logical expression on the second IF command tests whether *DEPT* equals *SALES*. When the condition is true, the value for the string variable *DIVISION* is changed to *TRANSFERRED* but is truncated if the format for *DIVISION* is not at least 11 characters wide. For any other value of *DEPT*, the value of *DIVISION* remains unchanged.
- Although there are two IF statements, each defines a separate and independent condition. The IF command is used rather than the DO IF-END IF structure in order to test both conditions on every case. If DO IF-END IF is used, control passes out of the structure as soon as the first logical condition is met.

**Example**

```
IF (STATE EQ 'IL' AND CITY EQ 13) COST=1.07 * COST.
```

- The logical expression tests whether *STATE* equals *IL* and *CITY* equals 13.
- If the logical expression is true, the numeric variable *COST* is increased by 7%.
- For any other value of *STATE* or *CITY*, the value of *COST* remains unchanged.

**Example**

```
STRING GROUP (A18).
IF (HIRED GE 1988) GROUP='Hired after merger'.
```

- STRING declares the string variable *GROUP* and assigns it a width of 18 characters.
- When *HIRED* is greater than or equal to 1988, *GROUP* is assigned the value *Hired after merger*. When *HIRED* is less than 1988, *GROUP* remains blank.

**Example**

```
IF (RECV GT DUE OR (REVNUES GE EXPNS AND BALNCE GT 0)) STATUS='SOLVENT'.
```

- First, the program tests whether *REVNUES* is greater than or equal to *EXPNS* and whether *BALNCE* is greater than 0.
- Second, the program evaluates if *RECV* is greater than *DUE*.
- If either of these expressions is true, *STATUS* is assigned the value *SOLVENT*.
- If both expressions are false, *STATUS* remains unchanged.
- *STATUS* is an existing string variable in the active dataset. Otherwise, it would have to be declared on a preceding STRING command.

## Operations

- Each `IF` command evaluates every case in the data. Compare `IF` with `DO IF`, which passes control for a case out of the `DO IF-END IF` structure as soon as a logical condition is met.
- The logical expression is evaluated as true, false, or missing. The assignment is executed only if the logical expression is true. If the logical expression is false or missing, the assignment is not made. Existing target variables remain unchanged; new numeric variables retain their initial (system-missing) values.
- In general, a logical expression is evaluated as missing if any one of the variables used in the logical expression is system- or user-missing. However, when relations are joined by the logical operators `AND` or `OR`, the expression can sometimes be evaluated as true or false even when variables have missing values. [For more information, see the topic Missing Values and Logical Operators on p. 965.](#)

## Numeric Variables

- Numeric variables created with `IF` are initially set to the system-missing value. By default, they are assigned an `F8.2` format.
- Logical expressions are evaluated in the following order: functions, followed by exponentiation, arithmetic operations, relations, and logical operators. When more than one logical operator is used, `NOT` is evaluated first, followed by `AND`, and then `OR`. You can change the order of operations using parentheses.
- Assignment expressions are evaluated in the following order: functions, then exponentiation, and then arithmetic operators.

## String Variables

- New string variables declared on `IF` are initially set to a blank value and are assigned the format specified on the `STRING` command that creates them.
- Logical expressions are evaluated in the following order: string functions, then relations, and then logical operators. When more than one logical operator is used, `NOT` is evaluated first, followed by `AND`, and then `OR`. You can change the order of operations using parentheses.
- If the transformed value of a string variable exceeds the variable's defined width, the transformed value is truncated. If the transformed value is shorter than the defined width, the string is right-padded with blanks.

## Missing Values and Logical Operators

When two or more relations are joined by logical operators `AND` or `OR`, the program always returns a missing value if all of the relations in the expression are missing. However, if any one of the relations can be determined, the program interprets the expression as true or false according to the logical outcomes below. The asterisk flags expressions where the program can evaluate the outcome with incomplete information.

Table 112-1  
*Logical outcomes*

<b>Expression</b>	<b>Outcome</b>	<b>Expression</b>	<b>Outcome</b>
true AND true	= true	true OR true	= true
true AND false	= false	true OR false	= true
false AND false	= false	false OR false	= false
true AND missing	= missing	true OR missing	= true*
missing AND missing	= missing	missing OR missing	= missing
false AND missing	= false*	false OR missing	= missing

# ***IGRAPH***

IGRAPH is no longer supported. Use the syntax converter to convert existing IGRAPH syntax. For more information, see the topic [TABLES and IGRAPH Command Syntax Converter](#) on p. 2234.

If you are writing new graph syntax, use the GGRAPH command. For more information, see the topic [GGRAPH](#) on p. 864.

# IMPORT

```
IMPORT FILE='file'  
  
  [/TYPE={COMM}  
    {TAPE}]  
  
  [/KEEP={ALL** } ] [/DROP=varlist]  
    {varlist}]  
  
  [/RENAME=(old varnames=new varnames)...]  
  
  [/MAP]
```

**\*\***Default if the subcommand is omitted.

## **Example**

```
IMPORT FILE=' /data/newdata.por' .
```

## **Overview**

IMPORT reads portable data files created with the EXPORT command. A portable data file is a data file created by the program and used to transport data between different types of computers and operating systems (such as between IBM CMS and Digital VAX/VMS) or between IBM® SPSS® Statistics and other software using the same portable file format. A portable file contains all of the data and dictionary information stored in the active dataset from which it was created.

The program can also read data files created by other software programs. See GET DATA and GET TRANSLATE for information on reading files created by spreadsheet and database programs such as dBASE, Lotus, and Excel.

## **Options**

**Format.** You can specify the format of the portable file (magnetic tape or communications program) on the TYPE subcommand.

**Variables.** You can read a subset of variables from the active dataset with the DROP and KEEP subcommands. You can rename variables using RENAME. You can also produce a record of all variables and their names in the active dataset with the MAP subcommand.

## **Basic Specification**

The basic specification is the FILE subcommand with a file specification. All variables from the portable file are copied into the active dataset with their original names, variable and value labels, missing-value flags, and print and write formats.

## **Subcommand Order**

- FILE and TYPE must precede all other subcommands.
- No specific order is required between FILE and TYPE or among other subcommands.

**Operations**

- The portable data file and dictionary become the active dataset and dictionary.
- A file saved with weighting in effect (using the `WEIGHT` command) automatically uses the case weights when the file is read.

**Examples**

```
IMPORT FILE="/data/newdata.por"
  /RENAME=(V1 TO V3=ID,SEX,AGE) /MAP.
```

- The active dataset is generated from the portable file *newdata.por*.
- Variables *V1*, *V2*, and *V3* are renamed *ID*, *SEX*, and *AGE* in the active dataset. Their names remain *V1*, *V2*, and *V3* in the portable file. None of the other variables copied into the active dataset are renamed.
- `MAP` requests a display of the variables in the active dataset.

**FILE Subcommand**

`FILE` specifies the portable file. `FILE` is the only required subcommand on `IMPORT`.

**TYPE Subcommand**

`TYPE` indicates whether the portable file is formatted for magnetic tape or for a communications program. `TYPE` can specify either `COMM` or `TAPE`. For more information on magnetic tapes and communications programs, see `EXPORT`.

<b>COMM</b>	<i>Communications-formatted file.</i> This is the default.
<b>TAPE</b>	<i>Tape-formatted file.</i>

**Example**

```
IMPORT TYPE=TAPE /FILE='hubout.por'.
```

- The file *hubout.por* is read as a tape-formatted portable file.

**DROP and KEEP Subcommands**

`DROP` and `KEEP` are used to read a subset of variables from the portable file.

- `DROP` excludes a variable or list of variables from the active dataset. All variables not named are included in the file.
- `KEEP` includes a variable or list of variables in the active dataset. All variables not specified on `KEEP` are excluded.
- `DROP` and `KEEP` cannot precede the `FILE` or `TYPE` subcommands.

- Variables can be specified in any order. The order of variables on `KEEP` determines the order of variables in the active dataset. The order on `DROP` does not affect the order of variables in the active dataset.
- If a variable is referred to twice on the same subcommand, only the first mention is recognized.
- Multiple `DROP` and `KEEP` subcommands are allowed; the effect is cumulative. Specifying a variable named on a previous `DROP` or not named on a previous `KEEP` results in an error and the command is not executed.
- The keyword `TO` can be used to specify a group of consecutive variables in the portable file.
- The portable file is not affected by `DROP` or `KEEP`.

### Example

```
IMPORT FILE='/data/newsum.por'
  /DROP=DEPT TO DIVISION.
```

- The active dataset is generated from the portable file *newsum.por*. Variables between and including *DEPT* and *DIVISION* in the portable file are excluded from the active dataset.
- All other variables are copied into the active dataset.

## RENAME Subcommand

`RENAME` renames variables being read from the portable file. The renamed variables retain the variable and value labels, missing-value flags, and print formats contained in the portable file.

- To rename a variable, specify the name of the variable in the portable file, a required equals sign, and the new name.
- A variable list can be specified on both sides of the equals sign. The number of variables on both sides must be the same, and the entire specification must be enclosed in parentheses.
- The keyword `TO` can be used for both variable lists.
- Any `DROP` or `KEEP` subcommand after `RENAME` must use the new variable names.

### Example

```
IMPORT FILE='/data/newsum.por'
  /DROP=DEPT TO DIVISION
  /RENAME=(NAME, WAGE=LNAME, SALARY) .
```

- `RENAME` renames *NAME* and *WAGE* to *LNAME* and *SALARY*.
- *LNAME* and *SALARY* retain the variable and value labels, missing-value flags, and print formats assigned to *NAME* and *WAGE*.

## MAP Subcommand

`MAP` displays a list of variables in the active dataset, showing all changes that have been specified on the `RENAME`, `DROP`, or `KEEP` subcommands.

- `MAP` can be specified as often as desired.



- `MAP` confirms only the changes specified on the subcommands that precede the `MAP` request.
- Results of subcommands that follow `MAP` are not mapped. When `MAP` is specified last, it also produces a description of the file.

**Example**

```
IMPORT FILE=' /data/newsun.por '  
  /DROP=DEPT TO DIVISION /MAP  
  /RENAME NAME=LNAME WAGE=SALARY /MAP.
```

- The first `MAP` subcommand produces a listing of the variables in the file after `DROP` has dropped the specified variables.
- `RENAME` renames *NAME* and *WAGE*.
- The second `MAP` subcommand shows the variables in the file after renaming.

# INCLUDE

```
INCLUDE FILE='file'  
      [ENCODING = 'encoding specification']
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- ENCODING keyword added for Unicode support.

## **Example**

```
INCLUDE FILE='/data/gsslabs.sps'.
```

## **Overview**

INCLUDE includes a file of commands in a session. INCLUDE is especially useful for including a long series of data definition statements or transformations. Another use for INCLUDE is to set up a library of commonly used commands and include them in the command sequence as they are needed.

*Note:* The newer INSERT provides equivalent functionality, plus additional features not available with INCLUDE. [For more information, see the topic INSERT on p. 979.](#)

INCLUDE allows you to run multiple commands together during a session and can save time. Complex or repetitive commands can be stored in a command file and included in the session, while simpler commands or commands unique to the current analysis can be entered during the session, before and after the included file.

## **Basic Specification**

The only specification is the FILE subcommand, which specifies the file to include. When INCLUDE is executed, the commands in the specified file are processed.

## **Syntax Rules**

- Commands in an included file must begin in column 1, and continuation lines for each command must be indented at least one column.
- The maximum line length for a command syntax file run via the INCLUDE command is 256 characters. Any characters beyond this limit are truncated.
- As many INCLUDE commands as needed can be used in a session.

- INCLUDE commands can be nested so that one set of included commands includes another set of commands. This nesting can go to five levels. However, a file cannot be included that is still open from a previous step.

### **Operations**

- If an included file contains a FINISH command, the session ends and no further commands are processed.
- If a journal file is created for the session, INCLUDE is copied to the journal file. Commands from the included file are also copied to the journal file but are treated like printed messages. Thus, INCLUDE can be executed from the journal file if the journal file is later used as a command file. Commands from the included file are executed only once.

## **ENCODING Keyword**

ENCODING specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is UTF8. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), UTF8, UTF16, UTF16BE (big endian), UTF16LE (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).

## **Examples**

```
INCLUDE FILE='/data/gsslabs.sps'.
```

- INCLUDE includes the file *gsslabs.sps* in the prompted session. When INCLUDE is executed, the commands in *gsslabs.sps* are processed.
- Assume that the include file *gsslabs.sps* contains the following:

```
DATA LIST FILE='/data/data52.txt'
  /RELIGION 5 OCCUPAT 7 SES 12 ETHNIC 15
  PARTY 19 VOTE48 33 VOTE52 41.
```

The active dataset will be defined and ready for analysis after INCLUDE is executed.

## **FILE Subcommand**

FILE identifies the file containing commands. FILE is the only specification on INCLUDE and is required.

# ***INFO***

This command is obsolete and no longer supported.

# INPUT PROGRAM-END INPUT PROGRAM

```
INPUT PROGRAM
commands to create or define cases
END INPUT PROGRAM
```

## Example

```
INPUT PROGRAM.
DATA LIST FILE=PRICES /YEAR 1-4 QUARTER 6 PRICE 8-12(2).

DO IF (YEAR GE 1881). /*Stop reading before 1881
END FILE.
END IF.
END INPUT PROGRAM.
```

## Overview

The INPUT PROGRAM and END INPUT PROGRAM commands enclose data definition and transformation commands that build cases from input records. The input program often encloses one or more DO IF-END IF or LOOP-END LOOP structures, and it must include at least one file definition command, such as DATA LIST. One of the following utility commands is also usually used:

<b>END CASE</b>	<i>Build cases from the commands within the input program and pass the cases to the commands immediately following the input program.</i>
<b>END FILE</b>	<i>Terminate processing of a data file before the actual end of the file or define the end of the file when the input program is used to read raw data.</i>
<b>REREAD</b>	<i>Reread the current record using a different DATA LIST.</i>
<b>REPEATING DATA</b>	<i>Read repeating groups of data from the same input record.</i>

For more information on the commands used in an input program, see the discussion of each command.

Input programs create a dictionary and data for an active dataset from raw data files; they cannot be used to read IBM® SPSS® Statistics data files. They can be used to process direct-access and keyed data files. For details, see KEYED DATA LIST.

## Basic Specification

The basic specification is INPUT PROGRAM, the commands used to create cases and define the active dataset, and END INPUT PROGRAM.

- INPUT PROGRAM and END INPUT PROGRAM each must be specified on a separate line and have no additional specifications.

---

*INPUT PROGRAM-END INPUT PROGRAM*

- To define an active dataset, the input program must include at least one `DATA LIST` or `END FILE` command.

**Operations**

- The `INPUT PROGRAM-END INPUT PROGRAM` structure defines an active dataset and is not executed until the program encounters a procedure or the `EXECUTE` command.
- `INPUT PROGRAM` clears the current active dataset.

**Examples****Select Cases with an Input Program**

```
INPUT PROGRAM.
DATA LIST FILE=PRICES /YEAR 1-4 QUARTER 6 PRICE 8-12(2).

DO IF (YEAR GE 1881). /*Stop reading when reaching 1881
END FILE.
END IF.
END INPUT PROGRAM.

LIST.
```

- The input program is defined between the `INPUT PROGRAM` and `END INPUT PROGRAM` commands.
- This example assumes that data records are entered chronologically by year. The `DO IF-END IF` structure specifies an end of file when the first case with a value of 1881 or later for *YEAR* is reached.
- `LIST` executes the input program and lists cases in the active dataset. The case that causes the end of the file is not included in the active dataset generated by the input program.
- As an alternative to this input program, you can use `N OF CASES` to select cases if you know the exact number of cases. Another alternative is to use `SELECT IF` to select cases before 1881, but then the program would unnecessarily read the entire input file.

**Skip the First n Records in a File**

```
INPUT PROGRAM.
NUMERIC          #INIT.
DO IF            NOT (#INIT).
+ LOOP          #I = 1 TO 5.
+ DATA LIST    NOTABLE/. /* No data - just skip record
+ END LOOP.
+ COMPUTE      #INIT = 1.
END IF.
DATA LIST        NOTABLE/ X 1.
END INPUT PROGRAM.

BEGIN DATA
A                /* The first 5 records are skipped
B
C
D
E
1
2
```

```

3
4
5
END DATA.
LIST.

```

- NUMERIC declares the scratch variable #INIT, which is initialized to system-missing.
- The DO IF structure is executed as long as #INIT does not equal 1.
- LOOP is executed five times. Within the loop, DATA LIST is specified without variable names, causing the program to read records in the data file without copying them into the active dataset. LOOP is executed five times, so the program reads five records in this manner. END LOOP terminates this loop.
- COMPUTE creates the scratch variable #INIT and sets it equal to 1. The DO IF structure is therefore not executed again.
- END IF terminates the DO IF structure.
- The second DATA LIST specifies numeric variable X, which is located in column 1 of each record. Because the program has already read five records, the first value for X that is copied into the active dataset is read from record 6.

## Input Programs

The program builds the active dataset dictionary when it encounters commands that create and define variables. At the same time, the program builds an *input program* that constructs cases and an optional *transformation program* that modifies cases prior to analysis or display. By the time the program encounters a procedure command that tells it to read the data, the active dataset dictionary is ready, and the programs that construct and modify the cases in the active dataset are built.

The internal input program is usually built from either a single DATA LIST command or from any of the commands that read or combine IBM® SPSS® Statistics data files (for example, GET, ADD FILES, MATCH FILES, UPDATE, and so on). The input program can also be built from the FILE TYPE-END FILE TYPE structure used to define nested, mixed, or grouped files. The third type of input program is specified with the INPUT PROGRAM-END INPUT PROGRAM commands.

With INPUT PROGRAM-END INPUT PROGRAM, you can create your own input program to perform many different operations on raw data. You can use transformation commands to build cases. You can read nonrectangular files, concatenate raw data files, and build cases selectively. You can also create an active dataset without reading any data at all.

## Input State

There are four program states in the program: the *initial state*, in which there is no active dataset dictionary; the *input state*, in which cases are created from the input file; the *transformation state*, in which cases are transformed; and the *procedure state*, in which procedures are executed. When you specify INPUT PROGRAM-END INPUT PROGRAM, you must pay attention to which commands are allowed within the input state, which commands can appear only within the input state, and which are not allowed within the input state.

## ***More Examples***

For additional examples of input programs, refer to DATA LIST, DO IF, DO REPEAT, END CASE, END FILE, LOOP, NUMERIC, POINT, REPEATING DATA, REREAD, and VECTOR.



# INSERT

*Note:* Equals signs (=) used in the syntax chart are required elements.

```
INSERT FILE='file specification'  
  [SYNTAX = {INTERACTIVE*}]  
    {BATCH      }  
  [ERROR = {CONTINUE*}]  
    {STOP      }  
  [CD = {NO*}]  
    {YES      }  
  [ENCODING = 'encoding specification']
```

\*Default if keyword omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 13.0

- Command introduced.

Release 16.0

- ENCODING keyword added for Unicode support.

## **Example**

```
INSERT FILE=' /examples/commands/file1.sps '  
  SYNTAX=BATCH ERROR=STOP CD=YES ENCODING='UTF8'.
```

## **OVERVIEW**

INSERT includes a file of commands in a session. INSERT is especially useful for including a long series of data definition statements or transformations. Another use for INSERT is to set up a library of commonly used commands and include them in the command sequence as they are needed.

INSERT allows you to run multiple commands together during a session and can save time. Complex or repetitive commands can be stored in a command file and included in the session, while simpler commands or commands unique to the current analysis can be entered during the session, before and after the included file.

INSERT provides the same basic functionality as INCLUDE, plus the ability to:

- Insert files that use either batch or interactive syntax rules.
- Control treatment of error conditions in inserted files.
- Change the working directory to the directory containing an inserted file.

### **Limitations**

The maximum line length for a command syntax file run via the `INSERT` command is 256 characters. Any characters beyond this limit are truncated.

## **FILE Keyword**

The minimum specification is the `FILE` keyword, followed by an equals sign and a quoted file specification (or quoted file handle) that specifies the file to insert. When the `INSERT` command is run, the commands in the specified file are processed.

### **Example**

```
INSERT FILE= '/examples/commands/file1.sps' .
```

## **SYNTAX Keyword**

The optional `SYNTAX` keyword specifies the syntax rules that apply to the inserted file. The keyword is followed by an equals sign (=) and one of the following alternatives:

**INTERACTIVE** *Each command must end with a period.* Periods can appear anywhere within the command, and commands can continue on multiple lines, but a period as the last non-blank character on a line is interpreted as the end of the command. Continuation lines and new commands can start anywhere on a new line. These are the “interactive” rules in effect when you select and run commands in a syntax window. This is the default if the `SYNTAX` keyword is omitted.

**BATCH** *Each command must start at the beginning of a new line (no blank spaces before the start of the command), and continuation lines must be indented at least one space.* If you want to indent new commands, you can use a plus sign, dash, or period as the first character at the start of the line and then indent the actual command. The period at the end of the command is optional. This setting is compatible with the syntax rules for command files included with the `INCLUDE` command.

Command syntax created with the Paste button in dialogs will work in either interactive or batch modes. For more information on interactive and batch syntax rules, see *Running Commands* on p. 38.

## **ERROR Keyword**

The optional `ERROR` keyword controls the handling of error conditions in inserted files. The keyword is followed by an equals sign (=) and one of the following alternatives:

**CONTINUE** *Errors in inserted files do not automatically stop command processing.* The inserted commands are treated as part of the normal command stream, and command processing continues in the normal fashion. This is the default if the `ERROR` keyword is omitted.

**STOP** *Command processing stops when the first error in an inserted file is encountered.* This is compatible with the behavior of command files included with the `INCLUDE` command.

## CD Keyword

The optional `CD` keyword can specify the directory containing the inserted file as the working directory, making it possible to use relative paths for file specifications within the inserted file. The keyword is followed by an equals sign (=) and one of the following alternatives:

- |            |   |
|------------|---|
| <b>NO</b>  | <i>The working directory is not changed.</i> This is the default if the <code>CD</code> keyword is omitted.   |
| <b>YES</b> | <i>The working directory is changed to the directory containing the inserted file.</i> Subsequent relative paths in command file specifications are interpreted as being relative to the location of the inserted file. |

The change in the working directory remains in effect until some other condition occurs that changes the working directory during the session, such as explicitly changing the working directory on another `INSERT` command with a `CD` keyword or a `CD` command that specifies a different directory (see `CD` on p. 300).

The `CD` keyword has no effect on the relative directory location for `SET` command file specifications, including `JOURNAL`, `CTEMPLATE`, and `TLOOK`. File specifications on the `SET` command should include complete path information.

The original working directory can be preserved with the `PRESERVE` command and later restored with the `RESTORE` command, as in:

```
PRESERVE.
INSERT FILE= '/commands/examples/file1.sps'
  CD=YES.
INSERT FILE='file2.sps'.
RESTORE.
```

- `PRESERVE` retains the original working directory location.
- The first `INSERT` command changes the working directory.
- The second `INSERT` command will look for `file2.sps` in `/commands/examples`.
- `RESTORE` resets the working directory to whatever it was prior to the first `INSERT` command.

For more information, see the `PRESERVE` and `RESTORE` commands.

## ENCODING Keyword

`ENCODING` specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is `UTF8`. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), `UTF8`, `UTF16`, `UTF16BE` (big endian), `UTF16LE` (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).

***INSERT vs. INCLUDE***

`INSERT` is a newer, more powerful and flexible alternative to `INCLUDE`. Files included with `INCLUDE` must always adhere to batch syntax rules, and command processing stops when the first error in an included file is encountered. You can effectively duplicate the `INCLUDE` behavior with `SYNTAX=BATCH` and `ERROR=STOP` on the `INSERT` command.

# KEYED DATA LIST

```
KEYED DATA LIST KEY=varname IN=varname  
  
FILE='file' [{TABLE } ] [ENCODING='encoding specification']  
          {NOTABLE}  
  
/varname {col location [(format)]} [varname ..]  
          {(FORTRAN-like format) }
```

## Release History

Release 16.0

- ENCODING subcommand added for Unicode support.

## Example

```
FILE HANDLE EMPL/ file specifications.  
KEYED DATA LIST FILE=EMPL KEY=#NXTCASE IN=#FOUND  
/YRHIRED 1-2 SEX 3 JOBCLASS 4.
```

## Overview

KEYED DATA LIST reads raw data from two types of nonsequential files: direct-access files, which provide direct access by a record number, and keyed files, which provide access by a record key. An example of a direct-access file is a file of 50 records, each corresponding to one of the United States. If you know the relationship between the states and the record numbers, you can retrieve the data for any specific state. An example of a keyed file is a file containing social security numbers and other information about a firm's employees. The social security number can be used to identify the records in the file.

### Direct-Access Files

There are various types of direct-access files. This program's concept of a direct-access file, however, is very specific. The file must be one from which individual records can be selected according to their number. The records in a 100-record direct-access file, for example, are numbered from 1 to 100.

Although the concept of record number applies to almost any file, not all files can be treated by this program as direct-access files. In fact, some operating systems provide no direct-access capabilities at all, and others permit only a narrowly defined subset of all files to be treated as direct access.

Very few files turn out to be good candidates for direct-access organization. In the case of an inventory file, for example, the usual large gaps in the part numbering sequence would result in large amounts of wasted file space. Gaps are not a problem, however, if they are predictable. For example, if you recognize that telephone area codes have first digits of 2 through 9, second digits

of 0 or 1, and third digits of 0 through 9, you can transform an area code into a record number by using the following COMPUTE statement:

```
COMPUTE RECNUM = 20*(DIGIT1-2) + 10*DIGIT2 + DIGIT3 + 1.
```

where *DIGIT1*, *DIGIT2*, and *DIGIT3* are variables corresponding to the respective digits in the area code, and *RECNUM* is the resulting record number. The record numbers would range from 1, for the nonexistent area code 200, through 160, for area code 919. The file would then have a manageable number of unused records.

### **Keyed Files**

Of the many kinds of keyed files, the ones to which the program can provide access are generally known as *indexed sequential files*. A file of this kind is basically a sequential file in which an index is maintained so that the file can be processed either sequentially or selectively. In effect, there is an underlying data file that is accessed through a file of index entries. The file of index entries may, for example, contain the fact that data record 797 is associated with social security number 476-77-1359. Depending on the implementation, the underlying data may or may not be maintained in sequential order.

The key for each record in the file generally comprises one or more pieces of information found within the record. An example of a complex key is a customer's last name and house number, plus the consonants in the street name, plus the zip code, plus a unique digit in case there are duplicates. Regardless of the information contained in the key, the program treats it as a character string.

On some systems, more than one key is associated with each record. That is, the records in a file can be identified according to different types of information. Although the primary key for a file normally must be unique, sometimes the secondary keys need not be. For example, the records in an employee file might be identified by social security number and job classification.

### **Options**

**Data Source.** You can specify the name of the keyed file on the FILE subcommand. By default, the last file that was specified on an input command, such as DATA LIST or REPEATING DATA, is read.

**Summary Table.** You can display a table that summarizes the variable definitions.

### **Basic Specification**

- The basic specification requires FILE, KEY, and IN, each of which specifies one variable, followed by a slash and variable definitions.
- FILE specifies the direct-access or keyed file. The file must have a file handle already defined.
- KEY specifies the variable whose value will be used to read a record. For direct-access files, the variable must be numeric; for keyed files, it must be string.
- IN creates a logical variable that flags whether a record was successfully read.
- Variable definitions follow all subcommands; the slash preceding them is required. Variable definitions are similar to those specified on DATA LIST.

**Subcommand Order**

- Subcommands can be named in any order.
- Variable definitions must follow all specified subcommands.

**Syntax Rules**

- Specifications for the variable definitions are the same as those described for DATA LIST. The only difference is that only one record can be defined per case.
- The FILE HANDLE command must be used if the FILE subcommand is specified on KEYED DATA LIST.
- KEYED DATA LIST can be specified in an input program, or it can be used as a transformation language to change an existing active dataset. This differs from all other input commands, such as GET and DATA LIST, which create new active datasets.

**Operations**

- Variable names are stored in the active dataset dictionary.
- Formats are stored in the active dataset dictionary and are used to display and write the values. To change output formats of numeric variables, use the FORMATS command.

**Examples****Specifying a Key Variable**

```
FILE HANDLE EMPL/ file specifications.
KEYED DATA LIST FILE=EMPL KEY=#NXTCASE IN=#FOUND
  /YRHIRED 1-2 SEX 3 JOBCLASS 4.
```

- FILE HANDLE defines the handle for the data file to be read by KEYED DATA LIST. The handle is specified on the FILE subcommand of KEYED DATA LIST.
- KEY on KEYED DATA LIST specifies the variable to be used as the access key. For a direct-access file, the value of the variable must be between 1 and the number of records in the file. For a keyed file, the value must be a string.
- IN creates the logical scratch variable *#FOUND*, whose value will be 1 if the record is successfully read, or 0 if the record is not found.
- The variable definitions are the same as those used for DATA LIST.

**Reading a Direct-Access File**

\* Reading a direct-access file: sampling 1 out of every 25 records.

```
FILE HANDLE      EMPL/ file specifications.
INPUT PROGRAM.
COMPUTE #INTRVL = TRUNC(UNIF(48))+1. /* Mean interval = 25
COMPUTE #NXTCASE = #NXTCASE+#INTRVL. /* Next record number
COMPUTE #EOF = #NXTCASE > 1000.     /* End of file check
DO IF #EOF.
+ END FILE.
ELSE.
+ KEYED DATA LIST FILE=EMPL, KEY=#NXTCASE, IN=#FOUND, NOTABLE
  /YRHIRED 1-2 SEX 3 JOBCLASS 4.
```

## KEYED DATA LIST

```

+ DO IF          #FOUND.
+   END CASE.          /* Return a case
+ ELSE.
+   PRINT / 'Oops. #NXTCASE=' #NXTCASE.
+ END IF.
END IF.
END INPUT PROGRAM.
EXECUTE.

```

- FILE HANDLE defines the handle for the data file to be read by the KEYED DATA LIST command. The record numbers for this example are generated by the transformation language; they are not based on data taken from another file.
- The INPUT PROGRAM and END INPUT PROGRAM commands begin and end the block of commands that build cases from the input file. Since the session generates cases, an input program is required.
- The first two COMPUTE statements determine the number of the next record to be selected. This is done in two steps. First, the integer portion is taken from the sum of 1 and a uniform pseudo-random number between 1 and 49. The result is a mean interval of 25. Second, the variable #NXTCASE is added to this number to generate the next record number. This record number, #NXTCASE, will be used for the key variable on the KEYED DATA LIST command. The third COMPUTE creates a logical scratch variable, #EOF, that has a value of 0 if the record number is less than or equal to 1000, or 1 if the value of the record number is greater than 1000.
- The DO IF-END IF structure controls the building of cases. If the record number is greater than 1000, #EOF equals 1, and the END FILE command tells the program to stop reading data and end the file.
- If the record number is less than or equal to 1000, the record is read via KEYED DATA LIST using the value of #NXTCASE. A case is generated if the record exists (#FOUND equals 1). If not, the program displays the record number and continues to the next case. The sample will have about 40 records.
- EXECUTE causes the transformations to be executed.
- This example illustrates the difference between DATA LIST, which always reads the next record in a file, and KEYED DATA LIST, which reads only specified records. The record numbers must be generated by another command or be contained in the active dataset.

**Reading a Keyed File**

\* Reading a keyed file: reading selected records.

```

GET FILE=STUDENTS/KEEP=AGE,SEX,COURSE.
FILE HANDLE COURSES/ file specifications.
STRING #KEY(A4).
COMPUTE #KEY = STRING(COURSE,N4). /* Create a string key
KEYED DATA LIST FILE=COURSES KEY=#KEY IN=#FOUND NOTABLE
/PERIOD 13 CREDITS 16.
SELECT IF #FOUND.
LIST.

```

- GET reads the STUDENTS file, which contains information on students, including a course identification for each student. The course identification will be used as the key for selecting one record from a file of courses.
- The FILE HANDLE command defines a file handle for the file of courses.



- The `STRING` and `COMPUTE` commands transform the course identification from numeric to string for use as a key. For keyed files, the key variable must be a string.
- `KEYED DATA LIST` uses the value of the newly created string variable `#KEY` as the key to search the course file. If a record that matches the value of `#KEY` is found, `#FOUND` is set to 1; otherwise, it is set to 0. Note that `KEYED DATA LIST` appears outside an input program in this example.
- If the course file contains the requested record, `#FOUND` equals 1. The variables `PERIOD` and `CREDITS` are added to the case and the case is selected via the `SELECT IF` command; otherwise, the case is dropped.
- `LIST` lists the values of the selected cases.
- This example shows how existing cases can be updated on the basis of information read from a keyed file.
- This task could also be accomplished by reading the entire course file with `DATA LIST` and combining it with the student file via the `MATCH FILES` command. The technique you should use depends on the percentage of the records in the course file that need to be accessed. If fewer than 10% of the course file records are read, `KEYED DATA LIST` is probably more efficient. As the percentage of the records that are read increases, reading the entire course file and using `MATCH` makes more sense.

## ***FILE Subcommand***

`FILE` specifies the handle for the direct-access or keyed data file. The file handle must have been defined on a previous `FILE HANDLE` command (or, in the case of the IBM OS environment, on a `DD` statement in the JCL).

## ***KEY Subcommand***

`KEY` specifies the variable whose value will be used as the key. This variable must already exist as the result of a prior `DATA LIST`, `KEYED DATA LIST`, `GET`, or transformation command.

- `KEY` is required. Its only specification is a single variable. The variable can be a permanent variable or a scratch variable.
- For direct-access files, the key variable must be numeric, and its value must be between 1 and the number of records in the file.
- For keyed files, the key variable must be string. If the keys are numbers, such as social security numbers, the `STRING` function can be used to convert the numbers to strings. For example, the following might be required to get the value of a numeric key into exactly the same format as used on the keyed file:

```
COMPUTE #KEY=STRING(123,IB4).
```

## ***IN Subcommand***

`IN` creates a numeric variable whose value indicates whether or not the specified record is found.

- `IN` is required. Its only specification is a single numeric variable. The variable can be a permanent variable or a scratch variable.
- The value of the variable is 1 if the record is successfully read or 0 if the record is not found. The `IN` variable can be used to select all cases that have been updated by `KEYED DATA LIST`.

**Example**

```
FILE HANDLE EMPL/ file specifications.
KEYED DATA LIST FILE=EMPL KEY=#NXTCASE IN=#FOUND
  /YRHIRED 1-2 SEX 3 JOBCLASS 4.
```

- `IN` creates the logical scratch variable `#FOUND`. The values of `#FOUND` will be 1 if the record indicated by the key value in `#NXTCASE` is found or 0 if the record does not exist.

**TABLE and NOTABLE Subcommands**

`TABLE` and `NOTABLE` determine whether the program displays a table that summarizes the variable definitions. `TABLE`, the default, displays the table. `NOTABLE` suppresses the table.

- `TABLE` and `NOTABLE` are optional and mutually exclusive.
- The only specification for `TABLE` or `NOTABLE` is the subcommand keyword. Neither subcommand has additional specifications.

**ENCODING Subcommand**

`ENCODING` specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is `UTF8`. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), `UTF8`, `UTF16`, `UTF16BE` (big endian), `UTF16LE` (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- In Unicode mode, the defined width of string variables is tripled for code page and UTF-16 text data files. Use [ALTER TYPE](#) to automatically adjust the defined width of string variables.

# KM

KM is available in the Advanced Statistics option.

```
KM varname [BY factor varname]
  /STATUS = varname [EVENT] (vallist) [LOST(vallist)]
  [/STRATA = varname]
  [/PLOT = {[SURVIVAL] [LOGSURV] [HAZARD] [OMS] }]
  [/ID = varname]
  [/PRINT = [TABLE**] [MEAN**] [NONE]]
  [/PERCENTILES = [( ){25, 50, 75 } ( )]]
  {value list }
  [/TEST = [LOGRANK**] [BRESLOW] [TARONE]]
  [/COMPARE = [{OVERALL**}] [{POOLED**}]]
  {PAIRWISE } {STRATA }
  [/TREND = [(METRIC)]]
  [/SAVE = tempvar[(newvar)], ...]
```

\*\*Default if the subcommand or keyword is omitted.

*Temporary variables created by Kaplan-Meier are:*

*SURVIVAL*

*HAZARD*

*SE*

*CUMEVENT*

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (2)
  /STRATA=LOCATION.
```

## **Overview**

KM (alias K-M) uses the Kaplan-Meier (product-limit) technique to describe and analyze the length of time to the occurrence of an event, often known as **survival time**. KM is similar to SURVIVAL in that it produces nonparametric estimates of the survival functions. However, instead of dividing the period of time under examination into arbitrary intervals, KM evaluates the survival function at the observed event times. For analysis of survival times with covariates, including time-dependent covariates, see the COXREG command.

### ***Options***

**KM Tables.** You can include one factor variable on the `KM` command. A KM table is produced for each level of the factor variable. You can also suppress the KM tables in the output with the `PRINT` subcommand.

**Survival Status.** You can specify the code(s) indicating that an event has occurred as well as code(s) for cases lost to follow-up using the `STATUS` subcommand.

**Plots.** You can plot the survival functions on a linear or log scale or plot the hazard function for each combination of factor and stratum with the `PLOT` subcommand.

**Test Statistics.** When a factor variable is specified, you can specify one or more tests of equality of survival distributions for the different levels of the factor using the `TEST` subcommand. You can also specify a trend metric for the requested tests with the `TREND` subcommand.

**Display ID and Percentiles.** You can specify an ID variable on the `ID` subcommand to identify each case. You can also request the display of percentiles in the output with the `PERCENTILES` subcommand.

**Comparisons.** When a factor variable is specified, you can use the `COMPARE` subcommand to compare the different levels of the factor, either pairwise or across all levels, and either pooled across all strata or within a stratum.

**Add New Variables to Active Dataset.** You can save new variables appended to the end of the active dataset with the `SAVE` subcommand.

### ***Basic Specification***

- The basic specification requires a survival variable and the `STATUS` subcommand naming a variable that indicates whether the event occurred.
- The basic specification prints one survival table followed by the mean and median survival time with standard errors and 95% confidence intervals.

### ***Subcommand Order***

- The survival variable and the factor variable (if there is one) must be specified first.
- Remaining subcommands can be specified in any order.

### ***Syntax Rules***

- Only one survival variable can be specified. To analyze multiple survival variables, use multiple `KM` commands.
- Only one factor variable can be specified following the `BY` keyword. If you have multiple factors, use the transformation language to create a single factor variable before invoking `KM`.
- Only one status variable can be listed on the `STATUS` subcommand. You must specify the value(s) indicating that the event occurred.
- Only one variable can be specified on the `STRATA` subcommand. If you have more than one stratum, use the transformation language to create a single variable to specify on the `STRATA` subcommand.

**Operations**

- KM deletes all cases that have negative values for the survival variable.
- KM estimates the survival function and associated statistics for each combination of factor and stratum.
- Three statistics can be computed to test the equality of survival functions across factor levels within a stratum or across all factor levels while controlling for strata. The statistics are the log rank (Mantel-Cox), generalized Wilcoxon (Breslow), and Tarone-Ware tests.
- When the `PLOTS` subcommand is specified, KM produces one plot of survival functions for each stratum, with all factor levels represented by different symbols or colors.

**Limitations**

- A maximum of 500 factor levels (symbols) can appear in a plot.

**Examples**

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (2)
  /STRATA=LOCATION.
```

- Survival analysis is used to examine the length of unemployment. The survival variable *LENGTH* contains the number of months a subject is unemployed. The factor variable *SEXRACE* combines sex and race factors.
- A value of 1 on the variable *EMPLOY* indicates the occurrence of the event (employment). All other observed cases are censored. A value of 2 on *EMPLOY* indicates cases lost to follow-up. Cases with other values for *EMPLOY* are known to have remained unemployed during the course of the study. KM separates the two types of censored cases in the KM table if *LOST* is specified.
- For each combination of *SEXRACE* and *LOCATION*, one KM table is produced, followed by the mean and median survival times with standard errors and confidence intervals.

**Using Kaplan-Meier to Test Competing Pain Relief Treatments**

```
KM
time BY treatment /STATUS=status(1)
  /PRINT TABLE MEAN
  /PERCENTILES
  /PLOT SURVIVAL
  /TEST LOGRANK BRESLOW TARONE
  /COMPARE OVERALL POOLED .
```

- Survival analysis is used to examine the time until the drug takes effect. The survival variable *time* is the time until effect. The factor variable *treatment* marks whether the experimental drug or the existing drug was administered.
- A value of 1 on the variable *status* indicates the occurrence of the event (drug takes effect). All other observed cases are censored.
- The `PRINT` subcommand produces a KM table for each value of *treatment*, followed by the mean and median survival times with standard errors and confidence intervals.

- The `PERCENTILES` subcommand, specified without keywords, requests the quartiles of the survival times for each value of *treatment*.
- The `PLOT` subcommand requests the cumulative survival distribution for each value of *treatment*.
- The `TEST` subcommand requests that the log rank, Breslow, and Tarone-Ware statistics be used to test the equality of the survival curves defined by the values of *treatment*. Per the `COMPARE` subcommand, all factor levels are compared in a single test.

## Survival and Factor Variables

You must identify the survival and factor variables for the analysis.

- The minimum specification is one, and only one, survival variable.
- Only one factor variable can be specified using the `BY` keyword. If you have more than one factor, create a new variable combining all factors. There is no limit to the factor levels.

### Example

```
DO IF SEX = 1.
+ COMPUTE SEXRACE = RACE.
ELSE.
+ COMPUTE SEXRACE = RACE + SEX.
END IF.
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (2).
```

- The two control variables, *SEX* and *RACE*, each with two values, 1 and 2, are combined into one factor variable, *SEXRACE*, with four values, 1 to 4.
- `KM` specifies *LENGTH* as the survival variable and *SEXRACE* as the factor variable.
- One `KM` table is produced for each factor level.

## STATUS Subcommand

To determine whether the terminal event has occurred for a particular observation, `KM` checks the value of a status variable. `STATUS` lists the status variable and the code(s) for the occurrence of the event. The code(s) for cases lost to follow-up can also be specified.

- Only one status variable can be specified. If multiple `STATUS` subcommands are specified, `KM` uses the last specification and displays a warning.
- The keyword `EVENT` is optional, but the value list in parentheses must be specified. Use `EVENT` for clarity's sake, especially when `LOST` is specified.
- The value list must be enclosed in parentheses. All cases with non-negative times that do not have a code within the range specified after `EVENT` are classified as **censored cases**—that is, cases for which the event has not yet occurred.
- The keyword `LOST` and the following value list are optional. `LOST` cannot be omitted if the value list for lost cases is specified.

- When `LOST` is specified, all cases with non-negative times that have a code within the specified value range are classified as lost to follow-up. Cases lost to follow-up are treated as censored in the analysis, and the statistics do not change, but the two types of censored cases are listed separately in the KM table.
- The value lists on `EVENT` or `LOST` can be one value, a list of values separated by blanks or commas, a range of values using the keyword `THRU`, or a combination.
- The status variable can be either numeric or string. If a string variable is specified, the `EVENT` or `LOST` values must be enclosed in apostrophes, and the keyword `THRU` cannot be used.

### Example

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (3,5 THRU 8).
```

- `STATUS` specifies that `EMPLOY` is the status variable.
- A value of 1 for `EMPLOY` means that the event (employment) occurred for the case.
- Values of 3 and 5 through 8 for `EMPLOY` mean that contact was lost with the case. The different values code different causes for the loss of contact.
- The summary table in the output includes columns for number lost and percentage lost, as well as for number censored and percentage censored.

## STRATA Subcommand

`STRATA` identifies a **stratification variable**—that is, a variable whose values are used to form subgroups (strata) within the categories of the factor variable. Analysis is done within each level of the strata variable for each factor level, and estimates are pooled over strata for an overall comparison of factor levels.

- The minimum specification is the subcommand keyword with one, and only one, variable name.
- If you have more than one strata variable, create a new variable to combine the levels on separate variables before invoking the `KM` command.
- There is no limit to the number of levels for the strata variable.

### Example

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (3,5 THRU 8)
  /STRATA=LOCATION.
```

- `STRATA` specifies `LOCATION` as the stratification variable. Analysis of the length of unemployment is done for each location within each sex and race subgroup.

## ***PLOT Subcommand***

PLOT plots the cumulative survival distribution on a linear or logarithmic scale or plots the cumulative hazard function. A separate plot with all factor levels is produced for each stratum. Each factor level is represented by a different symbol or color. Censored cases are indicated by markers.

- When PLOT is omitted, no plots are produced. The default is NONE.
- When PLOT is specified without a keyword, the default is SURVIVAL. A plot of survival functions for each stratum is produced.
- To request specific plots, specify, following the PLOT subcommand, any combination of the keywords defined below.
- Multiple keywords can be used on the PLOT subcommand, each requesting a different plot. The effect is cumulative.

<b>SURVIVAL</b>	<i>Plot the cumulative survival distribution on a linear scale. SURVIVAL is the default when PLOT is specified without a keyword.</i>
<b>LOGSURV</b>	<i>Plot the cumulative survival distribution on a logarithmic scale.</i>
<b>HAZARD</b>	<i>Plot the cumulative hazard function.</i>
<b>OMS</b>	<i>Plot the one-minus-survival function.</i>

### ***Example***

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (3,5 THRU 8)
  /STRATA=LOCATION
  /PLOT = SURVIVAL HAZARD.
```

- PLOT produces one plot of the cumulative survival distribution on a linear scale and one plot of the cumulative hazard rate for each value of *LOCATION*.

## ***ID Subcommand***

ID specifies a variable used for labeling cases. If the ID variable is a string, KM uses the string values as case identifiers in the KM table. If the ID variable is numeric, KM uses value labels or numeric values if value labels are not defined.

- ID is the first column of the KM table displayed for each combination of factor and stratum.
- If a string value or a value label exceeds 20 bytes in width, KM truncates the case identifier and displays a warning.



## **PRINT Subcommand**

By default, KM prints survival tables and the mean and median survival time with standard errors and confidence intervals if PRINT is omitted. If PRINT is specified, only the specified keyword is in effect. Use PRINT to suppress tables or the mean statistics.

<b>TABLE</b>	<i>Print the KM tables.</i> If PRINT is not specified, TABLE, together with MEAN, is the default. Specify TABLE on PRINT to suppress the mean statistics.
<b>MEAN</b>	<i>Print the mean statistics.</i> KM prints the mean and median survival time with standard errors and confidence intervals. If PRINT is not specified, MEAN, together with TABLE, is the default. Specify MEAN on PRINT to suppress the KM tables.
<b>NONE</b>	<i>Suppress both the KM tables and the mean statistics.</i> Only plots and comparisons are printed.

### **Example**

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (3,5 THRU 8)
  /STRATA=LOCATION
  /PLOT=SURVIVAL HAZARD
  /PRINT=NONE.
```

- PRINT=NONE suppresses both the KM tables and the mean statistics.

## **PERCENTILES Subcommand**

PERCENTILES displays percentiles for each combination of factor and stratum. Percentiles are not displayed without the PERCENTILES subcommand. If the subcommand is specified without a value list, the default is 25, 50, and 75 for quartile display. You can specify any values between 0 and 100.

## **TEST Subcommand**

TEST specifies the test statistic to use for testing the equality of survival distributions for the different levels of the factor.

- TEST is valid only when a factor variable is specified. If no factor variable is specified, KM issues a warning and TEST is not executed.
- If TEST is specified without a keyword, the default is LOGRANK. If a keyword is specified on TEST, only the specified test is performed.
- Each of the test statistics has a chi-square distribution with one degree of freedom.

<b>LOGRANK</b>	<i>Perform the log rank (Mantel-Cox) test.</i>
<b>BRESLOW</b>	<i>Perform the Breslow (generalized Wilcoxon) test.</i>
<b>TARONE</b>	<i>Perform the Tarone-Ware test.</i>

## COMPARE Subcommand

COMPARE compares the survival distributions for the different levels of the factor. Each of the keywords specifies a different method of comparison.

- COMPARE is valid only when a factor variable is specified. If no factor variable is specified, KM issues a warning and COMPARE is not executed.
- COMPARE uses whatever tests are specified on the TEST subcommand. If no TEST subcommand is specified, the log rank test is used.
- If COMPARE is not specified, the default is OVERALL and POOLED. All factor levels are compared across strata in a single test. The test statistics are displayed after the summary table at the end of output.
- Multiple COMPARE subcommands can be specified to request different comparisons.

<b>OVERALL</b>	<i>Compare all factor levels in a single test.</i> OVERALL, together with POOLED, is the default when COMPARE is not specified.
<b>PAIRWISE</b>	<i>Compare each pair of factor levels.</i> KM compares all distinct pairs of factor levels.
<b>POOLED</b>	<i>Pool the test statistics across all strata.</i> The test statistics are displayed after the summary table for all strata. POOLED, together with OVERALL, is the default when COMPARE is not specified.
<b>STRATA</b>	<i>Compare the factor levels for each stratum.</i> The test statistics are displayed for each stratum separately.

- If a factor variable has different levels across strata, you cannot request a pooled comparison. If you specify POOLED on COMPARE, KM displays a warning and ignores the request.

### Example

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (3,5 THRU 8)
  /STRATA=LOCATION
  /TEST = BRESLOW
  /COMPARE = PAIRWISE.
```

- TEST specifies the Breslow test.
- COMPARE uses the Breslow test statistic to compare all distinct pairs of *SEXRACE* values and pools the test results over all strata defined by *LOCATION*.
- Test statistics are displayed at the end of output for all strata.

## TREND Subcommand

TREND specifies that there is a trend across factor levels. This information is used when computing the tests for equality of survival functions specified on the TEST subcommand.

- The minimum specification is the subcommand keyword by itself. The default metric is chosen as follows:  
 If  $g$  is even,  
 $(-(g-1), \dots, -3, -1, 1, 3, \dots, (g-1))$   
 otherwise,

$$\left(-\frac{g-1}{2}, \dots, -1, 0, 1, \dots, \frac{g-1}{2}\right)$$

where  $g$  is the number of levels for the factor variable.

- If `TREND` is specified but `COMPARE` is not, `KM` performs the default log rank test with the trend metric for an `OVERALL POOLED` comparison.
- If the metric specified on `TREND` is longer than required by the factor levels, `KM` displays a warning and ignores extra values.

### Example

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (3,5 THRU 8)
  /STRATA=LOCATION
  /TREND.
```

- `TREND` is specified by itself. `KM` uses the default metric. Since `SEXRACE` has four levels, the default is  $(-3, -1, 1, 3)$ .
- Even though no `TEST` or `COMPARE` subcommand is specified, `KM` performs the default log rank test with the trend metric and does a default `OVERALL POOLED` comparison.

## SAVE Subcommand

`SAVE` saves the temporary variables created by `KM`. The following temporary variables can be saved:

<b>SURVIVAL</b>	<i>Survival function evaluated at current case.</i>
<b>SE</b>	<i>Standard error of the survival function.</i>
<b>HAZARD</b>	<i>Cumulative hazard function evaluated at current case.</i>
<b>CUMEVENT</b>	<i>Cumulative number of events.</i>

- To specify variable names for the new variables, assign the new names in parentheses following each temporary variable name.
- Assigned variable names must be unique in the active dataset. Scratch or system variable names cannot be used (that is, variable names cannot begin with `#` or `$`).
- If new variable names are not specified, `KM` generates default names. The default name is composed of the first three characters of the name of the temporary variable (two for `SE`), followed by an underscore and a number to make it unique.
- A temporary variable can be saved only once on the same `SAVE` subcommand.

### Example

```
KM LENGTH BY SEXRACE
  /STATUS=EMPLOY EVENT (1) LOST (3,5 THRU 8)
  /STRATA=LOCATION
  /SAVE SURVIVAL HAZARD.
```

- `KM` saves cumulative survival and cumulative hazard rates in two new variables, `SUR_1` and `HAZ_1`, provided that neither name exists in the active dataset. If one does, the numeric suffixes will be incremented to make a distinction.

# KNN

KNN is available in the Statistics Base option.

```
KNN [dependent variable [(MLEVEL = {S})]]
      {O}
      {N}
  [BY factor-list] [WITH covariate-list]

[/EXCEPT VARIABLES = varlist]

[/CASELABELS VARIABLE = varname]

[/FOCALCASES VARIABLE = varname]

[/RESCALE [COVARIATE = {ADJNORMALIZED**}]
      {NONE          } ]

[/PARTITION {TRAINING = {70** } HOLDOUT = {30** }}]
      {integer}      {integer}
      {VARIABLE = varname }

[/MODEL [METRIC = {EUCLID** }
      {CITYBLOCK}

      [NEIGHBORS = {FIXED**} [(K={3** }
      {integer}
      {AUTO } [(KMIN={3 } , KMAX={5 }
      {integer}      {integer}

      [FEATURES = {ALL**}
      {AUTO } [(FORCE = variable [variable ...])] ]

[/CRITERIA [NUMFEATURES = {AUTO** }
      {FIXED(integer)
      {ERRORRATIO(MINCHANGE={0.01 })}
      {value}

      [PREDICTED = {MEAN**}]
      {MEDIAN}

      [WEIGHTFEATURES = {NO**}]
      {YES }

[/CROSSVALIDATION {FOLDS = {10** }
      {integer}
      {VARIABLE = varname}

[/MISSING USERMISSING = {EXCLUDE**}]
      {INCLUDE }

[/VIEWMODEL [DISPLAY = {YES**}]
      {NO }

[/PRINT [CPS**] [NONE]]

[/SAVE [PREDFVAL[(varname)]]
      [PREDPROB[(rootname)]]
      [PARTITION[(varname)]]
      [FOLD[(varname)]]
      [MAXCAT({25** }
      {integer}

[/OUTFILE [MODEL = 'filename' ]
      [FOCALCASES = 'savfile' | 'dataset'].
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### Release History

Release 17.0

- Command introduced.

### Example

```
KNN DepVar BY A B C WITH X Y Z.
```

## Overview

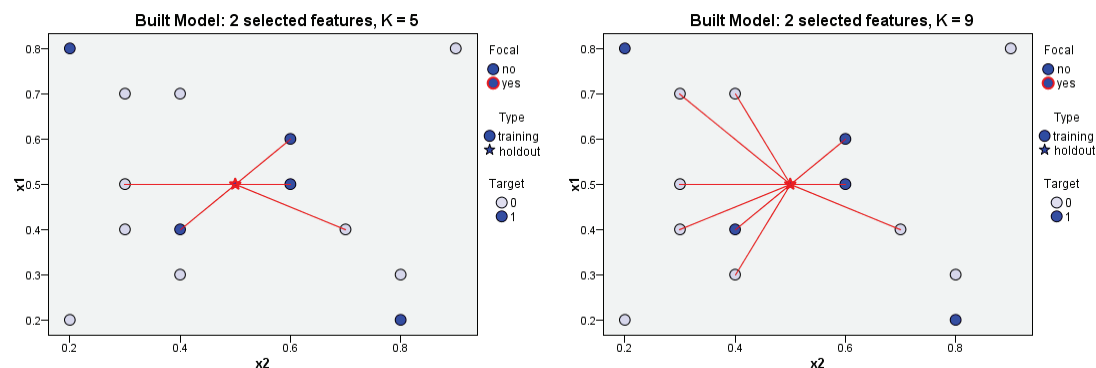
Nearest Neighbor Analysis is a method for classifying cases based on their similarity to other cases. In machine learning, it was developed as a way to recognize patterns of data without requiring an exact match to any stored patterns, or cases. Similar cases are near each other and dissimilar cases are distant from each other. Thus, the distance between two cases is a measure of their dissimilarity.

Cases that are near each other are said to be “neighbors.” When a new case (holdout) is presented, its distance from each of the cases in the model is computed. The classifications of the most similar cases – the nearest neighbors – are tallied and the new case is placed into the category that contains the greatest number of nearest neighbors.

You can specify the number of nearest neighbors to examine; this value is called  $k$ . The pictures show how a new case would be classified using two different values of  $k$ . When  $k = 5$ , the new case is placed in category 1 because a majority of the nearest neighbors belong to category 1. However, when  $k = 9$ , the new case is placed in category 0 because a majority of the nearest neighbors belong to category 0.

Figure 121-1

*The effects of changing  $k$  on classification*



Nearest neighbor analysis can also be used to compute values for a continuous target. In this situation, the average or median target value of the nearest neighbors is used to obtain the predicted value for the new case.

### ***Options***

**Prediction or classification.** The dependent variable may be scale, categorical, or a combination. If a dependent variable has scale measurement level, then the model predicts continuous values that approximate the “true” value of some continuous function of the input data. If a dependent variable is categorical, then the model is used to classify cases into the “best” category based on the input predictors.

**Rescaling.** KNN optionally rescales covariates; that is, predictors with scale measurement level, before training the model. Adjusted normalization is the rescaling method.

**Training and holdout partitions.** KNN optionally divides the data set into training and holdout partitions. The model is trained using the training partition. The holdout partition is completely excluded from the training process and is used for independent assessment of the final model.

**Missing Values.** The KNN procedure has an option for treating user-missing values of categorical variables as valid. User-missing values of scale variables are always treated as invalid. The procedure uses listwise deletion; that is, cases with invalid values for any variable are excluded from the model.

**Output.** KNN displays a case processing summary as pivot table output, and an interactive model view of other output. Tables in the model view include  $k$  nearest neighbors and distances for focal cases, classification of categorical response variables, and an error summary. Graphical output in the model view includes an automatic selection error log, feature importance chart, feature space chart, peers chart, and quadrant map. The procedure also optionally saves predicted values in the active dataset, PMML to an external file, and distances to focal cases to a new dataset or external file.

### ***Basic Specification***

The basic specification is the KNN command followed by zero or one dependent variable, the BY keyword and one or more factors, and the WITH keyword and one or more covariates.

By default, the KNN procedure normalizes covariates and selects a training sample before training the model. The model uses Euclidean distance to select the three nearest neighbors. User-missing values are excluded and default output is displayed.

If there are zero response variables, then the procedure finds the  $k$  nearest neighbors only – no classification or prediction is done.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### ***Syntax Rules***

- All subcommands are optional.
- Subcommands may be specified in any order.

- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not allowed.
- Any split variable defined on the `SPLIT FILE` command may not be used as a dependent variable, factor, covariate, or partition variable.

### ***Limitations***

Frequency weights specified on the `WEIGHT` command are ignored with a warning by the `KNN` procedure.

### ***Categorical Variables***

Although the `KNN` procedure accepts categorical variables as predictors or dependent variables, the user should be cautious when using a categorical variable with a very large number of categories.

The `KNN` procedure temporarily recodes categorical predictors using one-of- $c$  coding for the duration of the procedure. If there are  $c$  categories of a variable, then the variable is stored as  $c$  vectors, with the first category denoted  $(1,0,\dots,0)$ , the next category  $(0,1,0,\dots,0)$ , ..., and the final category  $(0,0,\dots,0,1)$ .

This coding scheme increases the dimensionality of the feature space. In particular, the total number of dimensions is the number of scale predictors plus the number of categories across all categorical predictors. As a result, this coding scheme can lead to slower training. If your nearest neighbors training is proceeding very slowly, you might try reducing the number of categories in your categorical predictors by combining similar categories or dropping cases that have extremely rare categories before running the `KNN` procedure.

All one-of- $c$  coding is based on the training data, even if a holdout sample is defined (see [PARTITION Subcommand](#)). Thus, if the holdout sample contains cases with predictor categories that are not present in the training data, then those cases are not scored. If the holdout sample contains cases with dependent variable categories that are not present in the training data, then those cases are scored.

### ***Replicating Results***

The `KNN` procedure uses random number generation during random assignment of partitions and cross-validation folds. To reproduce the same randomized results in the future, use the `SET` command to set the initialization value for the random number generator before each run of the `KNN` procedure, or use variables to define partitions and cross-validation folds.

## Examples

### Basic specification with default settings

```
KNN DepVar BY A B C WITH X Y Z.
```

- The KNN procedure treats *DepVar* as the dependent variable. Predictors *A*, *B*, and *C* are factors, and *X*, *Y*, and *Z* are covariates.
- By default, covariates are normalized before training. Also, the active dataset is partitioned into training and holdout data samples, with 70% going to the training data and 30% to the holdout data sample.
- The model uses Euclidean distance to select the three nearest neighbors.
- User-missing values are excluded and default output is displayed.

### No Target

```
*Nearest Neighbor Analysis.
KNN WITH price engine_s horsepow wheelbas widthlength curb_wgt fuel_cap mpg
/FOCALCASES VARIABLE=focal
/CASELABELS VARIABLE=model
/RESCALE COVARIATE=ADJNORMALIZED
/MODEL NEIGHBORS=FIXED(K=3) METRIC=EUCLID FEATURES=ALL
/CRITERIA WEIGHTFEATURES=NO
/PARTITION VARIABLE=partition
/PRINT CPS
/VIEWMODEL DISPLAY=YES
/MISSING USERMISSING=EXCLUDE.
```

- Since there is no dependent variable, the procedure finds the *k* nearest neighbors only – no classification or prediction is done. Predictors *price* through *mpg* are covariates.
- The FOCALCASES subcommand specifies *focal* as a focal case identifier. Cases with a positive value of *focal* are treated as focal cases.
- The CASELABELS subcommand specifies *model* as a case labeling variable. These values are used to label cases in the output.
- The PARTITION subcommand specifies that the variable *partition* should be used to assign each case in the active dataset to the training or holdout sample. Cases with a positive value on the variable are assigned to the training sample and cases with a non-positive value to the holdout sample.
- All other options are set to their default values.

### Automatic *k* Selection, Nominal Target

```
*Nearest Neighbor Analysis.
PRESERVE.
SET RNG=MT MTINDEX=20040509.
SHOW RNG.
KNN type (MLEVEL=0) WITH price engine_s horsepow wheelbas width length curb_wgt fuel_cap mp
/FOCALCASES VARIABLE=focal
/CASELABELS VARIABLE=model
/RESCALE COVARIATE=ADJNORMALIZED
/MODEL NEIGHBORS=AUTO(KMIN=3, KMAX=9) METRIC=EUCLID FEATURES=ALL
/CRITERIA WEIGHTFEATURES=YES
/PARTITION VARIABLE=partition
```



```

/CROSSVALIDATION FOLDS=10
/PRINT CPS
/VIEWMODEL DISPLAY=YES
/MISSING USERMISSING=EXCLUDE.
RESTORE.

```

- The `PRESERVE` and `RESTORE` commands ensure that the “current state” of the random number generator is retained and then restored after `KNN` is run.
- For this run of `KNN`, the random number generator is set to the Mersenne Twister with seed 20040509.
- The ordinal variable *type* is specified as a dependent variable with covariates *price* through *mpg* as predictors.
- The `FOCALCASES`, `CASELABELS`, and `PARTITION` subcommands are unchanged from the previous run.
- The `MODEL` subcommand specifies that the procedure should automatically select the “best” number of nearest neighbors (*k*) in the range 3 to 9, inclusive. Since feature selection is not requested, *k* selection is performed using cross-validation.
- The `CRITERIA` subcommand specifies that features should be weighted by importance when computing distances.
- The `CROSSVALIDATION` subcommand specifies that the procedure should subdivide the cases into 10 cross-validation folds for performing *k* selection.
- All other options are set to their default values.

### **Automatic *k* and Feature Selection, Scale Target**

```

*Nearest Neighbor Analysis.
KNN sales (MLEVEL=S) WITH price engine_s horsepower wheelbas width length curb_wgt fuel_cap r
/FOCALCASES VARIABLE=focal
/CASELABELS VARIABLE=model
/RESCALE COVARIATE=ADJNORMALIZED
/MODEL NEIGHBORS=AUTO (KMIN=3, KMAX=9) METRIC=EUCLID FEATURES=AUTO
/CRITERIA PREDICTED=MEAN WEIGHTFEATURES=YES NUMFEATURES=FIXED(5)
/PARTITION VARIABLE=partition
/SAVE PREDVAL
/PRINT CPS
/VIEWMODEL DISPLAY=YES
/MISSING USERMISSING=EXCLUDE.

```

- The scale variable *sales* is specified as a dependent variable with covariates *price* through *mpg* as predictors.
- The `FOCALCASES`, `CASELABELS`, and `PARTITION` subcommands are unchanged from the previous run.
- The `MODEL` subcommand specifies that the procedure should automatically select the “best” number of nearest neighbors (*k*) in the range 3 to 9, inclusive. Feature selection is also requested.
- The `CRITERIA` subcommand specifies that the final nearest neighbor model should have the 5 “best” features.

- The `SAVE` subcommand requests that the predicted value for `sales` be saved to the active dataset.
- All other options are set to their default values.

## ***Variable Lists***

The command line variable lists specify the dependent variable, any categorical predictors (also known as factors), and any scale predictors (also known as covariates).

### ***Dependent Variable***

- An optional dependent variable must be the first specification on the `KNN` command.
- The dependent variable may be followed by the measurement level specification, which contains, in parentheses, the `MLEVEL` keyword followed by an equals sign and then `S` for scale, `O` for ordinal, or `N` for nominal. `KNN` treats ordinal and nominal dependent variables equivalently as categorical.
- If a measurement level is specified, then it temporarily overrides a dependent variable's setting in the data dictionary.
- If no measurement level is specified, then `KNN` defaults to the dictionary setting.
- If a measurement level is not specified and no setting is recorded in the data dictionary, then a numeric variable is treated as scale and a string variable is treated as categorical.
- Dependent variables can be numeric or string.
- A string variable may be defined as ordinal or nominal only.

### ***Predictor Variables***

- The names of the factors, if any, must be preceded by the keyword `BY`.
- If keyword `BY` is specified with no factors, then a warning is issued and `BY` is ignored.
- The names of the covariates, if any, must be preceded by the keyword `WITH`.
- If keyword `WITH` is specified with no covariates, then a warning is issued and `WITH` is ignored.
- A dependent variable may not be specified within a factor or covariate list. If a dependent variable is specified within one of these lists, then an error is issued.
- All variables specified within a factor or covariate list must be unique. If duplicate variables are specified within a list, then the duplicates are ignored.
- If duplicate variables are specified across the factor and covariate lists, then an error is issued.
- Factor variables can be numeric or string.
- Covariates must be numeric.
- At least one predictor must be specified.

## ***EXCEPT Subcommand***

The `EXCEPT` subcommand lists any variables that the `KNN` procedure should exclude from the factor or covariate lists on the command line. This subcommand is useful if the factor or covariate lists contain a large number of variables – specified using the `TO` or `ALL` keyword, for example – but there are a few variables (for example, `Case ID`) that should be excluded.

The `EXCEPT` subcommand is introduced strictly for the purpose of simplifying syntax. Missing values on factors or covariates specified on `EXCEPT` do not affect whether a case is included in the analysis. For example, the following two `KNN` commands are equivalent. In both commands, listwise deletion is based on the dependent variable and factors *A*, *B*, and *C*.

```
KNN DepVar BY A B C.  
  
KNN DepVar BY A B C D  
  /EXCEPT VARIABLES=D.
```

- The `EXCEPT` subcommand ignores duplicate variables, and variables that are not specified on the command line's factor or covariate lists.
- There is no default variable list on the `EXCEPT` subcommand.

## ***CASELABELS Subcommand***

The `CASELABELS` subcommand allows you to specify a variable containing values used to label cases in the output.

- Cases are labeled using these values in the feature space chart, peers chart, and quadrant map.
- There is no default variable list on the `CASELABELS` subcommand.

## ***FOCALCASES Subcommand***

The `FOCALCASES` subcommand allows you to mark cases of particular interest. For example, a researcher wants to determine whether the test scores from one school district – the focal case – are comparable to those from similar school districts. He uses nearest neighbor analysis to find the school districts that are most similar with respect to a given set of features. Then he compares the test scores from the focal school district to those from the nearest neighbors.

Focal cases could also be used in clinical studies to select control cases that are similar to clinical cases. Focal cases are displayed in the *k* nearest neighbors and distances table, feature space chart, peers chart, and quadrant map. Information on focal cases is saved to the files specified on the `OUTFILE` subcommand.

- Cases with a positive value on the specified variable are treated as focal cases. It is invalid to specify a variable with no positive values.
- There is no default variable list on the `FOCALCASES` subcommand.

## **RESCALE Subcommand**

The `RESCALE` subcommand is used to rescale covariates.

All rescaling is performed based on the training data, even if a holdout sample is defined (see [PARTITION Subcommand](#)). That is, the minimum value or maximum value of a covariate are computed using only the training data. It is important that these covariates have similar distributions across the training and holdout samples. If the data are partitioned by specifying ratios on the `PARTITION` subcommand, then the `KNN` procedure attempts to ensure this similarity by random assignment. However, if you use the `PARTITION` subcommand `VARIABLE` keyword to assign cases to the training and holdout samples, then we recommend that you confirm the distributions are similar across samples before running the `KNN` procedure.

### **COVARIATE Keyword**

The `COVARIATE` keyword specifies the rescaling method to use for covariates specified following `WITH` on the command line. If no covariates are specified on the command line, then the `COVARIATE` keyword is ignored.

<b>ADJNORMAL- IZED</b>	<i>Adjusted version of subtract the minimum and divide by the range, <math>[2*(x-min)/(max-min)]-1</math>.</i>
<b>NONE</b>	<i>No rescaling of covariates.</i>

## **PARTITION Subcommand**

The `PARTITION` subcommand specifies the method of partitioning the active dataset into training and holdout samples. The training sample comprises the data records used to train the model. The holdout sample is an independent set of data records used to assess the final model.

- The partition can be defined by specifying the ratio of cases randomly assigned to each sample (training and holdout), or by a variable that assigns each case to the training or holdout sample.
- If the `PARTITION` subcommand is not specified, then the default partition randomly assigns 70% of the cases to the training sample and 30% to the holdout sample. If you want to specify a different random assignment, then you must specify new values for the `TRAINING` and `HOLDOUT` keywords. The value specified on each keyword gives the relative number of cases in the active dataset to assign to each sample. For example, `/PARTITION TRAINING = 50 HOLDOUT = 50` is equivalent to `/PARTITION TRAINING = 5 HOLDOUT = 5`; both subcommands randomly assign 50% of the cases to the training sample and 50% to the holdout sample.
- If you want to be able to reproduce results based on the `TRAINING` and `HOLDOUT` keywords later, use the `SET` command to set the initialization value for the random number generator before running the `KNN` procedure.
- Please see the discussion of the relationship between rescaling and partitioning in the `RESCALE` subcommand section above.

- All partitioning is performed after listwise deletion of any cases with invalid data for any variable used by the procedure. See the `MISSING` subcommand for details about valid and invalid data.
- It is invalid to specify both `TRAINING` and `VARIABLE`.

### ***TRAINING Keyword***

The `TRAINING` keyword specifies the relative number of cases in the active dataset to randomly assign to the training sample.

The value must be an integer greater than 0. The default (if the `PARTITION` subcommand is not specified) is 70.

### ***HOLDOUT Keyword***

The `HOLDOUT` keyword specifies the relative number of cases in the active dataset to randomly assign to the holdout sample.

The value must be an integer greater than or equal to 0. The default (if the `PARTITION` subcommand is not specified) is 30.

### ***VARIABLE Keyword***

The `VARIABLE` keyword specifies a variable that assigns each case in the active dataset to the training or holdout sample. Cases with a positive value on the variable are assigned to the training sample and cases with a non-positive value to the holdout sample. Cases with a system-missing value are excluded from the analysis. (Any user-missing values for the partition variable are always treated as valid.)

The variable may not be a dependent variable or any variable specified on the command line factor or covariate lists. The variable must be numeric.

## ***MODEL Subcommand***

The `MODEL` subcommand is used to specify the nearest neighbor “model”. By default, the procedure builds a model based on 3 nearest neighbors, using all features specified on the variables list and the Euclidean distance measure of “nearness”.

### ***METRIC Keyword***

The `METRIC` keyword allows you to specify the distance metric used to measure the similarity of cases.

- |                  |   |
|------------------|---|
| <b>EUCLID</b>    | <i>Euclidean distance.</i> This is the default specification for <code>METRIC</code> . The distance between two cases, $x$ and $y$ , is the square root of the sum, over all dimensions, of the squared differences between the values for the cases. |
| <b>CITYBLOCK</b> | <i>City-block or Manhattan distance.</i> The distance between two cases is the sum, over all dimensions, of the absolute differences between the values for the cases.  |

**NEIGHBORS Keyword**

The `NEIGHBORS` keyword indicates whether to use automatic selection of the number of nearest neighbors.

If no dependent variable is specified, then any specification other than `NEIGHBORS=FIXED` is ignored with a warning.

- FIXED**            *Use a fixed number of neighbors.* This is the default. The `FIXED` keyword may be followed by parentheses containing the `K` option, which specifies the number of neighbors. `K` must be a positive integer. The default value is 3.
- If `NEIGHBORS=FIXED` is specified, then any `CROSSVALIDATION` subcommand specifications are ignored with a warning.
- AUTO**            *Use automatic selection to determine the “best” number of neighbors.* The `AUTO` keyword may be followed by parentheses containing the `KMIN` and `KMAX` options, which specify the minimum and maximum number of neighbors, respectively, that automatic number of neighbors selection will consider in determining the “best” number of neighbors. It is invalid to specify only one option; you must specify both or neither. The options may be specified in any order and must be separated by a comma or space character. Both numbers must be integers greater than 0, with `KMIN` less than `KMAX`. The defaults are `KMIN=3`, `KMAX=5`.
- If `NEIGHBORS=AUTO` and `FEATURES=ALL` are specified, then  $V$ -fold cross-validation is used to select the “best” number of neighbors. The `CROSSVALIDATION` subcommand specifies the settings for  $V$ -fold cross-validation.

**FEATURES Keyword**

The `FEATURES` keyword indicates whether to use automatic selection of features (predictors).

If no dependent variable is specified, then any specification other than `FEATURES=ALL` is ignored with a warning.

- ALL**            *Use all predictors on the command line variable lists.* This is the default.
- AUTO**            *Use forward selection to determine the “best” feature set.* The `AUTO` keyword may be followed by parentheses containing the `FORCE` option, which specifies the starting set of predictors that must be included in the model. There is no default variable list on the `FORCE` option.
- If `FEATURES=AUTO` is specified, then any `CROSSVALIDATION` subcommand specifications are ignored with a warning.
- It is invalid for the variable list on the `FORCE` option to include all possible predictors; that is, there must be at least one predictor available for feature selection if `FEATURES=AUTO` is specified.

**Combined Neighbors and Features Selection**

When `NEIGHBORS=AUTO` and `FEATURES=AUTO`, the following method is used for combined neighbors and features selection:

1. For each  $k$ , use the forward selection method for feature selection.
2. Select the  $k$ , and accompanying feature set, with the lowest error rate or the lowest sum-of-squares error.

## CRITERIA Subcommand

The CRITERIA subcommand specifies computational and resource settings for the KNN procedure.

### NUMFEATURES Keyword

The NUMFEATURES keyword specifies how automatic feature selection should select the number of features.

This keyword is ignored if /MODEL FEATURES=AUTO is not in effect.

#### AUTO

*Select a fixed number of features, where the number of features is computed from a formula. The number of features the procedure selects is equal to  $\min(20, P) - J_{\text{Forced}}$ , where  $P$  is the total number of features and  $J_{\text{Forced}}$  is the number of forced features. This is the default.*

#### FIXED(integer)

*Select a fixed number of features, where the number of features is specified in advance. Specify a positive integer. It must be less than or equal to the number of unique predictors available for feature selection. This will be the number of predictors specified on the KNN command, minus any specified on the EXCEPT subcommand, minus any forced into the model.*

#### ERRORRATIO(MINCHANGE=value)

*Select features until the absolute change in the error ratio compared to the previous step is less than the criterion value. Specify a number greater than 0. The default value is 0.01.*

### PREDICTED Keyword

The PREDICTED keyword specifies the function used to compute the predicted value of scale response variables.

This keyword is ignored if no dependent variable is specified.

**MEAN**            *Compute predicted values based upon the mean value of the nearest neighbors. This is the default.*

**MEDIAN**        *Compute predicted values based upon the median value of the nearest neighbors.*

### WEIGHTFEATURES Keyword

The WEIGHTFEATURES keyword specifies whether to weight features by their normalized importance when computing distances.

**NO**                *Do not weight features by normalized importance. This is the default.*

**YES**              *Weight features by normalized importance. Feature importance for a predictor is calculated by the ratio of the error rate or sum-of-squares error of the model with the predictor removed from the model to the error rate or sum-of-squares error for the full model. Normalized importance is calculated by reweighting the feature importance values so that they sum to 1.*

*The specification WEIGHTFEATURES=YES is ignored with a warning if no dependent variable is specified.*

## **CROSSVALIDATION Subcommand**

The `CROSSVALIDATION` subcommand specifies settings for performing  $V$ -fold cross-validation to determine the “best” number of neighbors.

- $V$ -fold cross validation divides the data into  $V$  folds. Then, for a fixed  $k$ , it applies nearest neighbor analysis to make predictions on the  $v$ th fold (using the other  $V-1$  folds as the training sample) and evaluates the error. This process is successively applied to all possible choices of  $v$ . At the end of  $V$  folds, the computed errors are averaged. The above steps are repeated for various values of  $k$ . The value achieving the lowest average error is selected as the optimal value for  $k$ .
- If multiple values of  $k$  are tied on the lowest average error, then the smallest  $k$  among those that are tied is selected.
- Cross-validation is not used when `/MODEL NEIGHBORS=FIXED` or when `/MODEL FEATURES=AUTO`.
- It is invalid to specify both the `FOLDS` and `VARIABLE` keywords on the `CROSSVALIDATION` subcommand.

### **FOLDS Keyword**

The `FOLDS` keyword specifies the number of folds that should be used for cross-validation. The procedure randomly assigns cases to folds, numbered from 1 to the number of folds.

- Specify an integer greater than 1. The default is 10.
- For a given training set, the upper limit to the number of folds is the number of cases. If the value of `FOLDS` is greater than the number of cases in the training partition (and for any split, if `SPLIT FILE` is in effect), then the number of folds is set to the number of cases in the training partition (for that split).

### **VARIABLE Keyword**

The `VARIABLE` keyword specifies a variable that assigns each case in the active dataset to a fold from 1 to  $V$ .

The variable may not be a dependent variable or any variable specified on the command line factor or covariate lists. The variable must be numeric and take values from 1 to  $V$ . If any values in this range are missing on any splits (if `SPLIT FILE` is in effect) this will cause an error.

## **MISSING Subcommand**

The `MISSING` subcommand is used to control whether user-missing values for categorical variables are treated as valid values. By default, user-missing values for categorical variables are treated as invalid.

- The `MISSING` subcommand defines categorical variables as dependent variables with `MLEVEL` specification `O` or `N`, dependent variables with measurement level set at Ordinal or Nominal in the data dictionary and no `MLEVEL` specification, and factors. Scale variables are dependent



variables with MLEVEL specification S, dependent variables with Scale measurement level and no MLEVEL specification, and covariates.

- User-missing values for scale variables are always treated as invalid.
- System-missing values for any variables are always treated as invalid.

**USERMISSING=EXCLUDE**

*User-missing values for categorical variables are treated as invalid. This is the default.*

**USERMISSING=INCLUDE**

*User-missing values for categorical variables are treated as valid values.*

## **VIEWMODEL Subcommand**

The VIEWMODEL subcommand controls options for display of model-related output, including tables and charts. Tables in the model view include k nearest neighbors and distances for focal cases, classification of categorical response variables, and an error summary. Graphical output in the model view includes an automatic selection error log, feature importance chart, feature space chart, peers chart, and quadrant map.

### **DISPLAY Keyword**

The DISPLAY keyword specifies whether to display the model view in the output.

- YES**                    *Display the model view. This is the default.*  
**NO**                     *Do not display the model view.*

## **PRINT Subcommand**

The PRINT subcommand indicates the tabular output to display outside the Model Viewer. If PRINT is not specified, then the default tables are displayed. If PRINT is specified, then only the requested PRINT output is displayed.

### **CPS Keyword**

The CPS keyword displays the case processing summary table, which summarizes the number of cases included and excluded in the analysis, in total and by training and holdout samples. This table is shown by default.

### **NONE Keyword**

The NONE keyword suppresses all PRINT output. This keyword may not be specified with any other PRINT keywords.

## **SAVE Subcommand**

The SAVE subcommand writes optional temporary variables to the active dataset.

***PREDVAL Keyword***

This saves the predicted value for a scale dependent variable and the predicted category for a categorical dependent variable.

Specify a unique, valid variable name. The default variable name is *KNN\_PredictedValue*.

If there is no dependent variable, this keyword is ignored with a warning.

***PREDPROB Keyword***

If the dependent variable is categorical, then this keyword saves the predicted probabilities of the first *n* categories. The value of *n* is specified by the `MAXCAT` keyword.

- Specify a unique, valid variable name. The default root name is *KNN\_Probability*.
- If there is no dependent variable, this keyword is ignored with a warning.
- If the dependent variable is scale, this keyword is ignored with a warning.

***PARTITION Keyword***

This saves the value of the partition (training or holdout) to which the case was assigned.

- Specify a valid variable name. The default variable name is *KNN\_Partition*.
- This keyword is ignored with a warning if you specify a variable on the `PARTITION` subcommand.

***FOLD Keyword***

This saves the value of the fold to which the case was assigned.

- Specify a valid variable name. The default variable name is *KNN\_Fold*.
- This keyword is ignored if `/MODEL NEIGHBORS=FIXED` or `/MODEL FEATURES=AUTO`.
- This keyword is ignored with a warning if you specify a variable on the `CROSSVALIDATION` subcommand.

***MAXCAT Keyword***

The `MAXCAT` keyword specifies the number of categories for which predicted probabilities should be saved.

- Specify a positive integer in parentheses. The default is 25.
- This keyword is ignored if predicted probabilities are not saved.

***OUTFILE Subcommand***

The `OUTFILE` subcommand saves an XML-format file containing the nearest neighbor model. It also saves IBM® SPSS® Statistics-format data files containing distances from focal cases.

- File names must be specified in full. KNN does not supply extensions.
- The MODEL keyword is not honored if split-file processing is in effect (see the SPLIT FILE command). If this keyword is specified when split-file processing is on, then a warning is displayed.

**MODEL = 'file'**

*Writes the nearest neighbor model to an XML file.* Specify a valid file name. If the 'file' specification refers to an existing file, then the file is overwritten. The model is written in PMML with extensions.

**FOCALCASES = 'savfile' | 'dataset'**

*Writes the distances from focal cases to SPSS Statistics-format data files.* Specify a valid file or dataset name. If the specification refers to an existing file or dataset, then the file or dataset is overwritten.

For each focal case, a separate variable is created for each of the focal case's  $k$  nearest neighbors (from the training set) and the corresponding  $k$  nearest distances. The variable names are *KNN\_FocalCase\_CaseNumber*, *KNN\_NearestNeighbor\_CaseNumber\_1* through *KNN\_NearestNeighbor\_CaseNumber\_k* and *KNN\_NearestNeighbor\_Distance\_1* through *KNN\_NearestNeighbor\_CaseNumber\_k*.

# LEAVE

```
LEAVE varlist
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
COMPUTE TSALARY=TSALARY+SALARY.  
LEAVE TSALARY.  
FORMAT TSALARY (DOLLAR8)/ SALARY (DOLLAR7).  
EXECUTE.
```

## **Overview**

Normally, the program reinitializes variables each time it prepares to read a new case. `LEAVE` suppresses reinitialization and retains the current value of the specified variable or variables when the program reads the next case. It also sets the initial value received by a numeric variable to 0 instead of system-missing. `LEAVE` is frequently used with `COMPUTE` to create a variable to store an accumulating sum. `LEAVE` is also used to spread a variable's values across multiple cases when `VECTOR` is used within an input program to restructure a data file.

`LEAVE` cannot be used with scratch variables. [For more information, see the topic Scratch Variables on p. 51.](#)

## **Basic Specification**

The basic specification is the variable(s) whose values are not to be reinitialized as each new case is read.

## **Syntax Rules**

- Variables named on `LEAVE` must be new variables that do not already exist in the active dataset prior to the transformation block that defines them, but they must be defined in the transformation block prior to the `LEAVE` command that specifies them. [For more information, see the topic Examples on p. 1015.](#)
- Variables named on `LEAVE` cannot be scratch variables (but scratch variables can be used to obtain functionality equivalent to `LEAVE`). [For more information, see the topic Scratch Variables on p. 51.](#)
- Multiple variables can be named. The keyword `TO` can be used to refer to a list of consecutive variables.
- String and numeric variables can be specified on the same `LEAVE` command.

**Operations**

- Numeric variables named on LEAVE are initialized to 0 for the first case, and string variables are initialized to blanks. These variables are not reinitialized when new cases are read.

**Examples****Correct vs. Invalid Specifications for LEAVE**

```
DATA LIST LIST /Var1 Var2 Var3.
BEGIN DATA
1 2 3
4 5 6
7 8 9
END DATA.
```

```
*this is the correct form.
COMPUTE TotalVar1=TotalVar1+Var1.
LEAVE TotalVar1.
```

```
*this will change the value of Var2 but LEAVE will fail,
generating an error because Var2 already exists.
COMPUTE Var2=Var2+Var2.
LEAVE Var2.
```

```
*this will fail, generating an error because the LEAVE command
occurs before the command that defines the variable named on LEAVE.
LEAVE TotalVar3.
COMPUTE TotalVar3=TotalVar3+Var3.
```

```
LIST.
```

**Running Total**

```
COMPUTE TSALARY=TSALARY+SALARY.
LEAVE TSALARY.
FORMAT TSALARY (DOLLAR8)/ SALARY (DOLLAR7).
```

- These commands keep a running total of salaries across all cases. *SALARY* is the variable containing the employee's salary, and *TSALARY* is the new variable containing the cumulative salaries for all previous cases.
- For the first case, *TSALARY* is initialized to 0, and *TSALARY* equals *SALARY*. For the rest of the cases, *TSALARY* stores the cumulative totals for *SALARY*.
- LEAVE follows COMPUTE because *TSALARY* must first be defined before it can be specified on LEAVE.
- If LEAVE were not specified for this computation, *TSALARY* would be initialized to system-missing for all cases. *TSALARY* would remain system-missing because its value would be missing for every computation.

**Separate Sums for Each Category of a Grouping Variable**

```
SORT CASES DEPT.
IF DEPT NE LAG(DEPT,1) TSALARY=0. /*Initialize for new dept
COMPUTE TSALARY=TSALARY+SALARY. /*Sum salaries
LEAVE TSALARY. /*Prevent initialization each case
FORMAT TSALARY (DOLLAR8)/ SALARY (DOLLAR7).
```

- These commands accumulate a sum across cases for each department.
- `SORT` first sorts cases by the values of variable *DEPT*.
- `IF` specifies that if the value of *DEPT* for the current case is not equal to the value of *DEPT* for the previous case, *TSALARY* equals 0. Thus, *TSALARY* is reset to 0 each time the value of *DEPT* changes. (For the first case in the file, the logical expression on `IF` is missing. However, the desired effect is obtained because `LEAVE` initializes *TSALARY* to 0 for the first case, independent of the `IF` statement.)
- `LEAVE` prevents *TSALARY* from being initialized for cases within the same department.

# LINEAR

```
LINEAR
  /FIELDS
    TARGET = field
    INPUTS = fieldList
    [ANALYSIS_WEIGHT = field]

  [/BUILD_OPTIONS ]
    [OBJECTIVE = {STANDARD**}]
      {BOOSTING }
      {BAGGING }
      {LARGE }
    [USE_AUTO_DATA_PREPARATION = {TRUE**}]
      {FALSE }
    [CONFIDENCE_LEVEL = {95** }]
      {number}
    [MODEL_SELECTION = {FORWARDSTEPWISE**}]
      {BESTSUBSETS }
      {NONE }
    [CRITERIA_FORWARD_STEPWISE = {AICC** } ]
      {F }
      {ADJUSTEDRSQUARED}
      {ASE }
    [PROBABILITY_ENTRY = {0.05**}]
      {number}
    [PROBABILITY_REMOVAL = {0.1** } ]
      {number}
    [MAX_EFFECTS = {number}]
    [MAX_STEPS = {number}]
    [CRITERIA_BEST_SUBSETS = {AICC** } ]
      {ADJUSTEDRSQUARED}
      {ASE }
    [REPLICATE_RESULTS = {TRUE**}]
      {FALSE }
    [SEED = {54752075**}]
      {number }

  [/ENSEMBLES]
    [COMBINING_RULE_CONTINUOUS = {MEAN**}]
      {MEDIAN}
    [COMPONENT_MODELS_N = {10** } ]
      {integer}

  [/OUTFILE MODEL = 'file']

  [/SAVE PREDICTED_VALUES(varname)].
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 19

- Command introduced.

## **Example**

```
LINEAR
  /FIELDS
```

```
TARGET=y1  
INPUTS=x1 x2 .
```

## Overview

Linear models predict a continuous target based on linear relationships between the target and one or more predictors.

Linear models are relatively simple and give an easily interpreted mathematical formula for scoring. The properties of these models are well understood and are typically very fast to train.

### Options

**Ensemble models.** In addition to creating a “standard” linear model, you can use ensembles to improve model accuracy (boosting), improve model stability (bagging), or create a model for very large datasets.

**Model selection.** All fields can be entered into the model together, or forward stepwise or best subsets model selection can be used to choose the best fields.

**Output.** The procedure produces a model viewer object containing tables and charts for assessing the quality of the model; also you can save predicted values to the active dataset and the model to a file in PMML format.

### Basic Specification

The basic specification is the `LINEAR` command and `FIELDS` subcommand with `TARGET` and `INPUTS` keywords. There must be a single continuous target and at least one input.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### Syntax Rules

- All subcommands are optional.
- Subcommands may be specified in any order.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not allowed.
- Any split field defined on the `SPLIT FILE` command cannot be used on this command.
- Any field specified on the `WEIGHT` command cannot be used on this command.



**Limitations**

- If a categorical field has more than 1000 categories, the procedure does not run and no model is built.

**Examples****Automatic model selection**

```
*Automatic Linear Modeling.
LINEAR
  /FIELDS TARGET=claim_amount INPUTS=incident_date claim_type uninhabitable policy_date cov
    deductible townsize gender dob edcat job_start_date retire income marital reside occup
    primary_residence
  /BUILD_OPTIONS OBJECTIVE=STANDARD USE_AUTO_DATA_PREPARATION=TRUE CONFIDENCE_LEVEL=95
    MODEL_SELECTION=FORWARDSTEPWISE CRITERIA_FORWARD_STEPWISE=AICC REPLICATE_RESULTS=TRUE
  /ENSEMBLES COMBINING_RULE_CONTINUOUS=MEAN COMPONENT_MODELS_N=10.
```

- The procedure creates a linear model using *claim\_amount* as the target and *incident\_date* through *primary\_residence* as inputs (predictors).
- The `BUILD_OPTIONS` subcommand specifies forward stepwise as the model selection method and corrected Akaike information criterion (AICC) as the selection criterion.
- All other options are set to their default values.

**FIELDS Subcommand**

The `FIELDS` subcommand specifies the target, predictors, and optional analysis (regression) weight.

**TARGET.** The target field is what the model tries to predict. Specify a single continuous field.

**INPUTS.** The model uses values of the input fields to predict the target. Specify one or more fields. Continuous inputs are treated as covariates in the model, while categorical inputs are treated as factors.

- The universal keyword `TO` may be specified on the `INPUTS` list.

**ANALYSIS\_WEIGHT.** The model uses analysis (regression) weights to account for differences in variability across levels of the target field. Specify a single continuous field.

- This subcommand, and the `TARGET` and `INPUTS` keywords, are required.

**BUILD\_OPTIONS Subcommand**

The `BUILD_OPTIONS` subcommand specifies the criteria used to build the model.

**OBJECTIVE.** The objective determines which of the following model types is built.

- **STANDARD.** *Create a standard model.* The method builds a single model to predict the target using the predictors. Generally speaking, standard models are easier to interpret and can be faster to score than boosted, bagged, or large dataset ensembles.

- **BOOSTING.** *Enhance model accuracy (boosting).* The method builds an ensemble model using boosting, which generates a sequence of models to obtain more accurate predictions. Ensembles can take longer to build and to score than a standard model.
- **BAGGING.** *Enhance model stability (bagging).* The method builds an ensemble model using bagging (bootstrap aggregating), which generates multiple models to obtain more reliable predictions. Ensembles can take longer to build and to score than a standard model.
- **LARGE.** *Create a model for very large datasets (requires Server).* The method builds an ensemble model by splitting the dataset into separate data blocks. Choose this option if your dataset is too large to build any of the models above, or for incremental model building. This option can take less time to build, but can take longer to score than a standard model. This option requires Server connectivity and cannot be used when `SPLIT FILE` is active.

**USE\_AUTO\_DATA\_PREPARATION = TRUE | FALSE.** This option allows the procedure to transform the target and predictors in order to maximize the predictive power of the model. The original versions of transformed fields are excluded from the model. By default, the following automatic data preparation actions are performed. Note that this option is ignored and no automatic data preparation is performed if `OBJECTIVE=LARGE`.

- **Date and Time handling.** Each date predictor is transformed into new a continuous predictor containing the elapsed time since a reference date (1970-01-01). Each time predictor is transformed into a new continuous predictor containing the time elapsed since a reference time (00:00:00).
- **Adjust measurement level.** Continuous predictors with less than 5 distinct values are recast as ordinal predictors. Ordinal predictors with greater than 10 distinct values are recast as continuous predictors.
- **Outlier handling.** Values of continuous predictors that lie beyond a cutoff value (3 standard deviations from the mean) are set to the cutoff value.
- **Missing value handling.** Missing values of nominal predictors are replaced with the mode. Missing values of ordinal predictors are replaced with the median. Missing values of continuous predictors are replaced with the mean.
- **Supervised merging.** This makes a more parsimonious model by reducing the number of fields to be processed in association with the target. Similar categories are identified based upon the relationship between the input and the target. Categories that are not significantly different (that is, having a p-value greater than 0.1) are merged. If all categories are merged into one, the original and derived versions of the field are excluded from the model because they have no value as a predictor.

**CONFIDENCE\_LEVEL.** This is the level of confidence used to compute interval estimates of the model coefficients in the Coefficients view. Specify a value greater than 0 and less than 100. The default is 95.

**MODEL\_SELECTION.** Model selection methods determines how predictors are entered into the model.

- **FORWARDSTEPWISE.** This starts with no effects in the model and adds and removes effects one step at a time until no more can be added or removed according to the stepwise criteria. This is the default.

- **BESTSUBSETS.** This checks “all possible” models, or at least a larger subset of the possible models than forward stepwise, to choose the best according to the best subsets criterion. The model with the greatest value of the criterion is chosen as the best model. Note that Best subsets selection is more computationally intensive than forward stepwise selection. When best subsets is performed in conjunction with boosting, bagging, or very large datasets, it can take considerably longer to build than a standard model built using forward stepwise selection.
- **NONE.** Enters all available predictors into the model.

**CRITERIA\_FORWARD\_STEPWISE.** This is the statistic used to determine whether an effect should be added to or removed from the model when forward stepwise selection is used. If `MODEL_SELECTION = FORWARDSTEPWISE` is not specified, this keyword is ignored.

- **AICC.** Information Criterion (AICC) is based on the likelihood of the data given the model, and is adjusted to penalize overly complex models.
- **F.** The  $F$  statistic criterion is based on a statistical test of the improvement in model error.
- **ADJUSTEDRSQUARED.** Adjusted R-squared is based on the fit of the data, and is adjusted to penalize overly complex models.
- **ASE.** The average squared error (ASE) is an overfit prevention criterion based on the fit of the overfit prevention set. The overfit prevention set is a random subsample of approximately 30% of the original dataset that is not used to train the model.

If any criterion other than the  $F$  statistic is chosen, then at each step the effect that corresponds to the optimal change in the criterion is added to the model (the greatest increase for Adjusted R-squared, decrease for AICC and ASE). Any effects in the model that correspond to a decrease (increase for AICC and ASE) in the criterion are removed.

If the  $F$  statistic is chosen as the criterion, then at each step the effect that has the smallest p-value less than the specified `PROBABILITY_ENTRY` threshold is added to the model. Any effects in the model with a p-value greater than the specified `PROBABILITY_REMOVAL` threshold are removed.

**PROBABILITY\_ENTRY.** When forward stepwise selection is used with F Statistics as the criterion, this is the threshold for entering effects into the model. Specify a number greater than 0 and less than 1. The default is 0.05.

- The value of `PROBABILITY_ENTRY` must be less than the value of `PROBABILITY_REMOVAL`.

**PROBABILITY\_REMOVAL** When forward stepwise selection is used with F Statistics as the criterion, this is the threshold for removing effects from the model. Specify a number greater than 0 and less than 1. The default is 0.10.

**MAX\_EFFECTS = number.** Customize maximum number of effects in the final model when forward selection is used. By default, all available effects can be entered into the model. Alternatively, if the stepwise algorithm ends a step with the specified maximum number of effects, the algorithm stops with the current set of effects.

**MAX\_STEPS = number.** Customize the maximum number of steps when forward selection is used. The stepwise algorithm stops after a certain number of steps. By default, this is 3 times the number of available effects. Alternatively, specify a positive integer maximum number of steps.

**CRITERIA\_BEST\_SUBSETS.** This is the statistic used to choose the “best” model when best subsets selection is used. If `MODEL_SELECTION = FORWARDSTEPWISE` is not specified, this keyword is ignored.

- **AICC.** Information Criterion (AICC) is based on the likelihood of the data given the model, and is adjusted to penalize overly complex models.
- **ADJUSTEDRSQUARED.** Adjusted R-squared is based on the fit of the data, and is adjusted to penalize overly complex models.
- **ASE.** Overfit Prevention Criterion (ASE) is based on the fit of the overfit prevention set. The overfit prevention set is a random subsample of approximately 30% of the original dataset that is not used to train the model.

**REPLICATE\_RESULTS = TRUE\*\*|FALSE.** Setting a random seed allows you to replicate analyses. The random number generator is used to choose which records are in the overfit prevention set.

**SEED = 54752075\*\* | number.** When `REPLICATE_RESULTS=TRUE`, this is the value of the random seed. Specify an integer. The default is 54752075.

## ***ENSEMBLES Subcommand***

The `ENSEMBLES` subcommand determines the behavior of ensembling that occurs when boosting, bagging, or very large datasets are requested on the `OBJECTIVE` keyword. Options that do not apply to the selected objective are ignored.

**COMBINING\_RULE\_CONTINUOUS = MEAN\*\* | MEDIAN.** Ensemble predicted values for continuous targets can be combined using the mean or median of the predicted values from the base models.

Note that when the objective is to enhance model accuracy (`BAGGING`), the combining rule selections are ignored. Boosting always uses a weighted majority vote to score categorical targets and a weighted median to score continuous targets.

**COMPONENT\_MODELS\_N = 10\*\* | integer.** Specify the number of base models to build when the objective is to enhance model accuracy or stability; for bagging, this is the number of bootstrap samples. It should be a positive integer.

## ***SAVE Subcommand***

The `SAVE` subcommand writes optional temporary variables to the active dataset.

**PREDICTED\_VALUES(varname).** *Predicted value of the target.* The default variable name is *PredictedValue*.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand saves a file containing the scoring model.

**MODEL = 'file'**. Writes the model to an external *.zip* file. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#) Specify a unique, valid filename. If the file specification refers to an existing file, then the file is overwritten.

# LIST

```
LIST [VARIABLES={ALL** } ] [/FORMAT={WRAP**} [{UNNUMBERED**}] ]  
      {varlist}           {SINGLE}      {NUMBERED }  
  
  [/CASES=[FROM {1**} ] [TO {eof**}] [BY {1**}]]  
          {n }           {n }           {n }
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
LIST VARIABLES=V1 V2.
```

## Overview

LIST displays case values for variables in the active dataset. The output is similar to the output produced by the PRINT command. However, LIST is a procedure and reads data, whereas PRINT is a transformation and requires a procedure (or the EXECUTE command) to execute it.

### Options

**Selecting and Ordering Variables.** You can specify a list of variables to be listed using the VARIABLES subcommand.

**Format.** You can limit each case listing to a single line, and you can display the case number for each listed case with the FORMAT subcommand.

**Selecting Cases.** You can limit the listing to a particular sequence of cases using the CASES subcommand.

### Basic Specification

- The basic specification is simply LIST, which displays the values for all variables in the active dataset.
- By default, cases wrap to multiple lines if all the values do not fit within the page width (the page width is determined by the SET WIDTH command). Case numbers are not displayed for the listed cases.

### Subcommand Order

All subcommands are optional and can be named in any order.

### **Operations**

- If `VARIABLES` is not specified, variables are listed in the order in which they appear in the active dataset.
- `LIST` does not display values for scratch or system variables.
- `LIST` uses print formats contained in the dictionary of the active dataset. Alternative formats cannot be specified on `LIST`. See `FORMATS` or `PRINT FORMATS` for information on changing print formats.
- `LIST` output uses the width specified on `SET`.
- If a numeric value is longer than its defined width, the program first attempts to list the value by removing punctuation characters, then uses scientific notation, and finally prints asterisks.
- If a long string variable cannot be listed within the output width, it is truncated.
- Values of the variables listed for a case are always separated by at least one blank.
- System-missing values are displayed as a period for numeric variables and a blank for string variables.
- If cases fit on one line, the column width for each variable is determined by the length of the variable name or the format, whichever is greater. If the variable names do not fit on one line, they are printed vertically.
- If cases do not fit on one line within the output width specified on `SET`, they are wrapped. `LIST` displays a table illustrating the location of the variables in the output and prints the name of the first variable in each line at the beginning of the line.
- Each execution of `LIST` begins at the top of a new page. If `SPLIT FILE` is in effect, each split also begins at the top of a new page.

## **Examples**

### ***LIST with No Subcommands***

```
LIST.
```

- `LIST` by itself requests a display of the values for all variables in the active dataset.

### ***Controlling Listed Cases with CASES Subcommand***

```
LIST VARIABLES=V1 V2 /CASES=FROM 10 TO 100 BY 2.
```

- `LIST` produces a list of every second case for variables `V1` and `V2`, starting with case 10 and stopping at case 100.

## **VARIABLES Subcommand**

`VARIABLES` specifies the variables to be listed.

- The variables must already exist, and they cannot be scratch or system variables.
- If `VARIABLES` is used, only the specified variables are listed.
- Variables are listed in the order in which they are named on `VARIABLES`.

- If a variable is named more than once, it is listed more than once.
- The keyword `ALL` (the default) can be used to request all variables. `ALL` can also be used with a variable list (see example below).

**ALL** *List all user-defined variables. Variables are listed in the order in which they appear in the active dataset. This is the default if `VARIABLES` is omitted.*

### **Example**

```
LIST VARIABLES=V15 V31 ALL.
```

- `VARIABLES` is used to list values for `V15` and `V31` before all other variables. The keyword `ALL` then lists all variables, including `V15` and `V31`, in the order in which they appear in the active dataset. Values for `V15` and `V31` are therefore listed twice.

## **FORMAT Subcommand**

`FORMAT` controls whether cases wrap if they cannot fit on a single line and whether the case number is displayed for each listed case. The default display uses more than one line per case (if necessary) and does not number cases.

- The minimum specification is a single keyword.
- `WRAP` and `SINGLE` are alternatives, as are `NUMBERED` and `UNNUMBERED`. Only one of each pair can be specified.
- If `SPLIT FILE` is in effect for `NUMBERED`, case numbering restarts at each split. To get sequential numbering regardless of splits, create a variable and set it equal to the system variable `$CASENUM` and then name this variable as the first variable on the `VARIABLES` subcommand. An appropriate format should be specified for the new variable before it is used on `LIST`.

**WRAP** *Wrap cases if they do not fit on a single line. Page width is determined by the `SET WIDTH` command. This is the default.*

**SINGLE** *Limit each case to one line. Only variables that fit on a single line are displayed.*

**UNNUMBERED** *Do not include the sequence number of each case. This is the default.*

**NUMBERED** *Include the sequence number of each case. The sequence number is displayed to the left of the listed values.*

## **CASES Subcommand**

`CASES` limits the number of cases listed. By default, all cases in the active dataset are listed.

- Any or all of the keywords below can be used. Defaults that are not changed remain in effect.
- If `LIST` is preceded by a `SAMPLE` or `SELECT IF` command, case selections specified by `CASES` are taken from those cases that were selected by `SAMPLE` or `SELECT IF`.



- If `SPLIT FILE` is in effect, case selections specified by `CASES` are restarted for each split.

**FROM n**                      *Number of the first case to be listed.* The default is 1.

**TO n**                         *Number of the last case to be listed.* The default is the end of the active dataset.  
`CASES 100` is interpreted as `CASES TO 100`.

**BY n**                         *Increment used to choose cases for listing.* The default is 1.

### ***Example***

```
LIST CASES BY 3 /FORMAT=NUMBERED.
```

- Every third case is listed for all variables in the active dataset. The listing begins with the first case and includes every third case up to the end of the file.
- `FORMAT` displays the case number of each listed case.

### ***Example***

```
LIST CASES FROM 10 TO 20.
```

- Cases from case 10 through case 20 are listed for all variables in the active dataset.

# LOGISTIC REGRESSION

LOGISTIC REGRESSION is available in the Regression option.

```
LOGISTIC REGRESSION VARIABLES = dependent var
    [WITH independent varlist [BY var [BY var] ... ]]

[/CATEGORICAL = var1, var2, ... ]

[/CONTRAST (categorical var) = [{INDICATOR [(refcat)]    }]]
    {DEVIATION [(refcat)]    }
    {SIMPLE [(refcat)]      }
    {DIFFERENCE             }
    {HELMERT                }
    {REPEATED               }
    {POLYNOMIAL[({1,2,3...})] }
    {metric                  }
    {SPECIAL (matrix)       }

[/METHOD = {ENTER**        }   [{ALL    }]]
    {BSTEP [{COND}]        }   {varlist}
    {LR    }
    {WALD }
    {FSTEP [{COND}]        }
    {LR    }
    {WALD }

[/SELECT = {ALL**          }]]
    {varname relation value}

[/{NOORIGIN**}]
    {ORIGIN    }

[/ID = [variable]]

[/PRINT = [DEFAULT**] [SUMMARY] [CORR] [ALL] [ITER [({1})]] [GOODFIT]]
    {n}
    [CI(level)]

[/CRITERIA = [BCON ({0.001**})] [ITERATE({20**})] [LCON({0**  })]
    {value } {n } {value }
    [PIN({0.05**})] [POUT({0.10**})] [EPS({.00000001**})]]
    {value } {value } {value }
    [CUT[({0.5**  })]
    {value }

[/CLASSPLOT]

[/MISSING = {EXCLUDE **}]
    {INCLUDE    }

[/CASEWISE = [tempvarlist] [OUTLIER({2  })]]
    {value}

[/SAVE = tempvar[(newname)] tempvar[(newname)]...]

[/OUTFILE = [{MODEL    } (filename)]
    {PARAMETER}

[/EXTERNAL]
```

\*\*Default if the subcommand or keyword is omitted.

Temporary variables that are created by LOGISTIC REGRESSION are as follows:

<i>PRED</i>	<i>LEVER</i>	<i>COOK</i>
<i>PGROUP</i>	<i>LRESID</i>	<i>DFBETA</i>
<i>RESID</i>	<i>SRESID</i>	
<i>DEV</i>	<i>ZRESID</i>	

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- `OUTFILE` subcommand introduced.

Release 14.0

- Modification to the method of recoding string variables. [For more information, see the topic Overview on p. 1029.](#)

### **Example**

```
LOGISTIC REGRESSION VARIABLES = PROMOTED WITH AGE, JOBTIME, JOBRATE.
```

## **Overview**

LOGISTIC REGRESSION regresses a dichotomous dependent variable on a set of independent variables. Categorical independent variables are replaced by sets of contrast variables, each set entering and leaving the model in a single step.

### **Options**

**Processing of Independent Variables.** You can specify which independent variables are categorical in nature on the `CATEGORICAL` subcommand. You can control treatment of categorical independent variables by the `CONTRAST` subcommand. Seven methods are available for entering independent variables into the model. You can specify any one of them on the `METHOD` subcommand. You can also use the keyword `BY` between variable names to enter interaction terms.

**Selecting Cases.** You can use the `SELECT` subcommand to define subsets of cases to be used in estimating a model.

**Regression through the Origin.** You can use the `ORIGIN` subcommand to exclude a constant term from a model.

**Specifying Termination and Model-Building Criteria.** You can further control computations when building the model by specifying criteria on the `CRITERIA` subcommand.

**Adding New Variables to the Active Dataset.** You can save the residuals, predicted values, and diagnostics that are generated by LOGISTIC REGRESSION in the active dataset.

**Output.** You can use the `PRINT` subcommand to print optional output, use the `CASEWISE` subcommand to request analysis of residuals, and use the `ID` subcommand to specify a variable whose values or value labels identify cases in output. You can request plots of the actual values and predicted values for each case with the `CLASSPLOT` subcommand.

### **Basic Specification**

- The minimum specification is the `VARIABLES` subcommand with one dichotomous dependent variable. You must specify a list of independent variables either following the keyword `WITH` on the `VARIABLES` subcommand or on a `METHOD` subcommand.
- The default output includes goodness-of-fit tests for the model ( $-2$  log-likelihood, goodness-of-fit statistic, Cox and Snell  $R^2$ , and Nagelkerke  $R^2$ ) and a classification table for the predicted and observed group memberships. The regression coefficient, standard error of the regression coefficient, Wald statistic and its significance level, and a multiple correlation coefficient adjusted for the number of parameters (Atkinson, 1980) are displayed for each variable in the equation.

### **Subcommand Order**

- Subcommands can be named in any order. If the `VARIABLES` subcommand is not specified first, a slash (/) must precede it.
- The ordering of `METHOD` subcommands determines the order in which models are estimated. Different sequences may result in different models.

### **Syntax Rules**

- Only one dependent variable can be specified for each `LOGISTIC REGRESSION`.
- Any number of independent variables may be listed. The dependent variable may not appear on this list.
- The independent variable list is required if any of the `METHOD` subcommands are used without a variable list or if the `METHOD` subcommand is not used. The keyword `TO` cannot be used on any variable list.
- If you specify the keyword `WITH` on the `VARIABLES` subcommand, all independent variables must be listed.
- If the keyword `WITH` is used on the `VARIABLES` subcommand, interaction terms do not have to be specified on the variable list, but the individual variables that make up the interactions must be listed.
- Multiple `METHOD` subcommands are allowed.
- The minimum truncation for this command is `LOGI REG`.

### **Operations**

- Independent variables that are specified on the `CATEGORICAL` subcommand are replaced by sets of contrast variables. In stepwise analyses, the set of contrast variables associated with a categorical variable is entered or removed from the model as a single step.
- Independent variables are screened to detect and eliminate redundancies.

- If the linearly dependent variable is one of a set of contrast variables, the set will be reduced by the redundant variable or variables. A warning will be issued, and the reduced set will be used.
- For the forward stepwise method, redundancy checking is done when a variable is to be entered into the model.
- When backward stepwise or direct-entry methods are requested, all variables for each METHOD subcommand are checked for redundancy before that analysis begins.

### **Compatibility**

Prior to version 14.0, the order of recoded string values was dependent on the order of values in the data file. For example, when recoding the dependent variable, the first string value encountered was recoded to 0, and the second string value encountered was recoded to 1. Beginning with version 14.0, the procedure recodes string variables so that the order of recoded values is the alphanumeric order of the string values. Thus, the procedure may recode string variables differently than in previous versions.

### **Limitations**

- The dependent variable must be dichotomous for each split-file group. Specifying a dependent variable with more or less than two nonmissing values per split-file group will result in an error.

## **Examples**

```
LOGISTIC REGRESSION VAR=default
  /SELECT validate EQ 1
  /METHOD=FSTEP(LR) age ed employ address income debtinc creddebt othdebt
  /CONTRAST (ed)=Indicator
  /SAVE PRED COOK SRESID
  /CLASSPLOT
  /PRINT=GOODFIT
  /CRITERIA PIN(.05) POUT(.10) ITERATE(20) CUT(.5) .
```

- The procedure fits a logistic regression model for *default*.
- The SELECT subcommand specifies that only cases with a value of 1 on the variable *validate* will be used to build the model. Other cases will be scored and can be used to validate the final model.
- The variables specified on the METHOD subcommand will be entered into the model using forward stepwise regression based on the significance of the likelihood ratio statistic.
- The CONTRAST subcommand specifies that *ed* is a categorical variable and should be processed using indicator coding.
- The SAVE subcommand requests that the casewise predicted values, Cook's influence statistics, and studentized residuals be saved as new variables to the dataset.
- The CLASSPLOT subcommand requests a classification plot of the observed and predicted values.
- The PRINT subcommand requests the Hosmer-Lemeshow goodness-of-fit statistic.
- All other options are set to their default values.

## VARIABLES Subcommand

VARIABLES specifies the dependent variable and, optionally, all independent variables in the model. The dependent variable appears first on the list and is separated from the independent variables by the keyword WITH.

- One VARIABLES subcommand is allowed for each Logistic Regression procedure.
- The dependent variable must be dichotomous—that is, it must have exactly two values other than system-missing and user-missing values for each split-file group.
- The dependent variable may be a string variable if its two values can be differentiated by their first eight characters.
- You can indicate an interaction term on the variable list by using the keyword BY to separate the individual variables.
- If all METHOD subcommands are accompanied by independent variable lists, the keyword WITH and the list of independent variables may be omitted.
- If the keyword WITH is used, *all* independent variables must be specified. For interaction terms, only the individual variable names that make up the interaction (for example, X1, X2) need to be specified. Specifying the actual interaction term (for example, X1 BY X2) on the VARIABLES subcommand is optional if you specify it on a METHOD subcommand.

### Example

```
LOGISTIC REGRESSION VARIABLES = PROMOTED WITH AGE,JOBTIME,JOBRATE,  
    AGE BY JOBTIME.
```

- *PROMOTED* is specified as the dependent variable.
- *AGE*, *JOBTIME*, *JOBRATE*, and the interaction *AGE* by *JOBTIME* are specified as the independent variables.
- Because no METHOD is specified, all three single independent variables and the interaction term are entered into the model.
- LOGISTIC REGRESSION produces the default output.

## CATEGORICAL Subcommand

CATEGORICAL identifies independent variables that are nominal or ordinal. Variables that are declared to be categorical are automatically transformed to a set of contrast variables as specified on the CONTRAST subcommand. If a variable that is coded as 0 – 1 is declared as categorical, its coding scheme is given indicator contrasts by default.

- Independent variables that are not specified on CATEGORICAL are assumed to be at least interval level, except for string variables.
- Any variable that is specified on CATEGORICAL is ignored if it does not appear either after WITH on the VARIABLES subcommand or on any METHOD subcommand.
- Variables that are specified on CATEGORICAL are replaced by sets of contrast variables. If the categorical variable has  $n$  distinct values, there will be  $n-1$  contrast variables generated. The set of contrast variables associated with a categorical variable is entered or removed from the model as a step.

- If any one of the variables in an interaction term is specified on CATEGORICAL, the interaction term is replaced by contrast variables.
- All string variables are categorical. Only the first eight characters of each value of a string variable are used in distinguishing between values. Thus, if two values of a string variable are identical for the first eight characters, the values are treated as though they were the same.

### Example

```
LOGISTIC REGRESSION VARIABLES = PASS WITH GPA, GRE, MAT, CLASS, TEACHER
/CATEGORICAL = CLASS,TEACHER.
```

- The dichotomous dependent variable *PASS* is regressed on the interval-level independent variables *GPA*, *GRE*, and *MAT* and the categorical variables *CLASS* and *TEACHER*.

## CONTRAST Subcommand

CONTRAST specifies the type of contrast that is used for categorical independent variables. The interpretation of the regression coefficients for categorical variables depends on the contrasts that are used. The default is INDICATOR. The categorical independent variable is specified in parentheses following CONTRAST. The closing parenthesis is followed by one of the contrast-type keywords.

- If the categorical variable has  $n$  values, there will be  $n-1$  rows in the contrast matrix. Each contrast matrix is treated as a set of independent variables in the analysis.
- Only one categorical independent variable can be specified per CONTRAST subcommand, but multiple CONTRAST subcommands can be specified.

The following contrast types are available (Finn, 1974), (Kirk, 1982).

<b>INDICATOR(refcat)</b>	<i>Indicator variables.</i> Contrasts indicate the presence or absence of category membership. By default, refcat is the last category (represented in the contrast matrix as a row of zeros). To omit a category (other than the last category), specify the sequence number of the omitted category (which is not necessarily the same as its value) in parentheses after the keyword INDICATOR.
<b>DEVIATION(refcat)</b>	<i>Deviations from the overall effect.</i> The effect for each category of the independent variable (except one category) is compared to the overall effect. Refcat is the category for which parameter estimates are not displayed (they must be calculated from the others). By default, refcat is the last category. To omit a category (other than the last category), specify the sequence number of the omitted category (which is not necessarily the same as its value) in parentheses after the keyword DEVIATION.
<b>SIMPLE(refcat)</b>	<i>Each category of the independent variable (except the last category) is compared to the last category.</i> To use a category other than the last as the omitted reference category, specify its sequence number (which is not necessarily the same as its value) in parentheses following the keyword SIMPLE.
<b>DIFFERENCE</b>	<i>Difference or reverse Helmert contrasts.</i> The effects for each category of the independent variable (except the first category) are compared to the mean effects of the previous categories.

<b>HELMERT</b>	<i>Helmert contrasts.</i> The effects for each category of the independent variable (except the last category) are compared to the mean effects of subsequent categories.
<b>POLYNOMIAL(metric)</b>	<i>Polynomial contrasts.</i> The first degree of freedom contains the linear effect across the categories of the independent variable, the second degree of freedom contains the quadratic effect, and so on. By default, the categories are assumed to be equally spaced; unequal spacing can be specified by entering a metric consisting of one integer for each category of the independent variable in parentheses after the keyword <code>POLYNOMIAL</code> . For example, <code>CONTRAST(STIMULUS)=POLYNOMIAL(1,2,4)</code> indicates that the three levels of <code>STIMULUS</code> are actually in the proportion 1:2:4. The default metric is always (1,2, ..., k), where k categories are involved. Only the relative differences between the terms of the metric matter: (1,2,4) is the same metric as (2,3,5) or (20,30,50) because the difference between the second and third numbers is twice the difference between the first and second numbers in each instance.
<b>REPEATED</b>	<i>Comparison of adjacent categories.</i> Each category of the independent variable (except the last category) is compared to the next category.
<b>SPECIAL(matrix)</b>	<i>A user-defined contrast.</i> After this keyword, a matrix is entered in parentheses with $k-1$ rows and $k$ columns (where $k$ is the number of categories of the independent variable). The rows of the contrast matrix contain the special contrasts indicating the desired comparisons between categories. If the special contrasts are linear combinations of each other, <code>LOGISTIC REGRESSION</code> reports the linear dependency and stops processing. If $k$ rows are entered, the first row is discarded and only the last $k-1$ rows are used as the contrast matrix in the analysis.

**Example**

```
LOGISTIC REGRESSION VARIABLES = PASS WITH GRE, CLASS
/CATEGORICAL = CLASS
/CONTRAST(CLASS)=HELMERT.
```

- A logistic regression analysis of the dependent variable *PASS* is performed on the interval independent variable *GRE* and the categorical independent variable *CLASS*.
- *PASS* is a dichotomous variable representing course pass/fail status and *CLASS* identifies whether a student is in one of three classrooms. A `HELMERT` contrast is requested.

**Example**

```
LOGISTIC REGRESSION VARIABLES = PASS WITH GRE, CLASS
/CATEGORICAL = CLASS
/CONTRAST(CLASS)=SPECIAL(2 -1 -1
                        0 1 -1).
```

- In this example, the contrasts are specified with the keyword `SPECIAL`.

**METHOD Subcommand**

`METHOD` indicates how the independent variables enter the model. The specification is the `METHOD` subcommand followed by a single method keyword. The keyword `METHOD` can be omitted. Optionally, specify the independent variables and interactions for which the method is to be used. Use the keyword `BY` between variable names of an interaction term.



- If no variable list is specified, or if the keyword `ALL` is used, all of the independent variables following the keyword `WITH` on the `VARIABLES` subcommand are eligible for inclusion in the model.
- If no `METHOD` subcommand is specified, the default method is `ENTER`.
- Variables that are specified on `CATEGORICAL` are replaced by sets of contrast variables. The set of contrast variables associated with a categorical variable is entered or removed from the model as a single step.
- Any number of `METHOD` subcommands can appear in a Logistic Regression procedure. `METHOD` subcommands are processed in the order in which they are specified. Each method starts with the results from the previous method. If `BSTEP` is used, all remaining eligible variables are entered at the first step. All variables are then eligible for entry and removal unless they have been excluded from the `METHOD` variable list.
- The beginning model for the first `METHOD` subcommand is either the constant variable (by default or if `NOORIGIN` is specified) or an empty model (if `ORIGIN` is specified).

The available `METHOD` keywords are as follows:

<b>ENTER</b>	<i>Forced entry.</i> All variables are entered in a single step. This setting is the default if the <code>METHOD</code> subcommand is omitted.
<b>FSTEP</b>	<i>Forward stepwise.</i> The variables (or interaction terms) that are specified on <code>FSTEP</code> are tested for entry into the model one by one, based on the significance level of the score statistic. The variable with the smallest significance less than <code>PIN</code> is entered into the model. After each entry, variables that are already in the model are tested for possible removal, based on the significance of the conditional statistic, the Wald statistic, or the likelihood-ratio criterion. The variable with the largest probability greater than the specified <code>POUT</code> value is removed, and the model is reestimated. Variables in the model are then evaluated again for removal. When no more variables satisfy the removal criterion, covariates that are not in the model are evaluated for entry. Model building stops when no more variables meet entry or removal criteria or when the current model is the same as a previous model.
<b>BSTEP</b>	<i>Backward stepwise.</i> As a first step, the variables (or interaction terms) that are specified on <code>BSTEP</code> are entered into the model together and are tested for removal one by one. Stepwise removal and entry then follow the same process as described for <code>FSTEP</code> until no more variables meet entry or removal criteria or when the current model is the same as a previous model.

The statistic that is used in the test for removal can be specified by an additional keyword in parentheses following `FSTEP` or `BSTEP`. If `FSTEP` or `BSTEP` is specified by itself, the default is `COND`.

<b>COND</b>	<i>Conditional statistic.</i> This setting is the default if <code>FSTEP</code> or <code>BSTEP</code> is specified by itself.
<b>WALD</b>	<i>Wald statistic.</i> The removal of a variable from the model is based on the significance of the Wald statistic.
<b>LR</b>	<i>Likelihood ratio.</i> The removal of a variable from the model is based on the significance of the change in the log-likelihood. If <code>LR</code> is specified, the model must be reestimated without each of the variables in the model. This process can substantially increase computational time. However, the likelihood-ratio statistic is the best criterion for deciding which variables are to be removed.

**Example**

```
LOGISTIC REGRESSION
  VARIABLES = PROMOTED WITH AGE JOBTIME JOBRATE RACE SEX AGENCY
  /CATEGORICAL RACE SEX AGENCY
  /METHOD ENTER AGE JOBTIME
  /METHOD BSTEP (LR) RACE SEX JOBRATE AGENCY.
```

- *AGE*, *JOBTIME*, *JOBRATE*, *RACE*, *SEX*, and *AGENCY* are specified as independent variables. *RACE*, *SEX*, and *AGENCY* are specified as categorical independent variables.
- The first `METHOD` subcommand enters *AGE* and *JOBTIME* into the model.
- Variables in the model at the termination of the first `METHOD` subcommand are included in the model at the beginning of the second `METHOD` subcommand.
- The second `METHOD` subcommand adds the variables *RACE*, *SEX*, *JOBRATE*, and *AGENCY* to the previous model.
- Backward stepwise logistic regression analysis is then done with only the variables on the `BSTEP` variable list tested for removal by using the `LR` statistic.
- The procedure continues until all variables from the `BSTEP` variable list have been removed or the removal of a variable will not result in a decrease in the log-likelihood with a probability larger than `POUT`.

**SELECT Subcommand**

By default, all cases in the active dataset are considered for inclusion in `LOGISTIC REGRESSION`. Use the optional `SELECT` subcommand to include a subset of cases in the analysis.

- The specification is either a logical expression or keyword `ALL`. `ALL` is the default. Variables that are named on `VARIABLES`, `CATEGORICAL`, or `METHOD` subcommands cannot appear on `SELECT`.
- In the logical expression on `SELECT`, the relation can be `EQ`, `NE`, `LT`, `LE`, `GT`, or `GE`. The variable must be numeric, and the value can be any number.
- Only cases for which the logical expression on `SELECT` is true are included in calculations. All other cases, including those cases with missing values for the variable that is named on `SELECT`, are unselected.
- Diagnostic statistics and classification statistics are reported for both selected and unselected cases.
- Cases that are deleted from the active dataset with the `SELECT IF` or `SAMPLE` command are not included among either the selected or unselected cases.

**Example**

```
LOGISTIC REGRESSION VARIABLES=GRADE WITH GPA, TUCE, PSI
  /SELECT SEX EQ 1 /CASEWISE=RESID.
```

- Only cases with the value 1 for *SEX* are included in the logistic regression analysis.
- Residual values that are generated by `CASEWISE` are displayed for both selected and unselected cases.

## **ORIGIN and NOORIGIN Subcommands**

ORIGIN and NOORIGIN control whether the constant is included. NOORIGIN (the default) includes a constant term (intercept) in all equations. ORIGIN suppresses the constant term and requests regression through the origin. (NOCONST can be used as an alias for ORIGIN.)

- The only specification is either ORIGIN or NOORIGIN.
- ORIGIN or NOORIGIN can be specified only once per Logistic Regression procedure, and it affects all METHOD subcommands.

### **Example**

```
LOGISTIC REGRESSION VARIABLES=PASS WITH GPA,GRE,MAT /ORIGIN.
```

- ORIGIN suppresses the automatic generation of a constant term.

## **ID Subcommand**

ID specifies a variable whose values or value labels identify the casewise listing. By default, cases are labeled by their case number.

- The only specification is the name of a single variable that exists in the active dataset.
- Only the first eight characters of the variable's value labels are used to label cases. If the variable has no value labels, the values are used.
- Only the first eight characters of a string variable are used to label cases.

## **PRINT Subcommand**

PRINT controls the display of optional output. If PRINT is omitted, DEFAULT output (defined below) is displayed.

- The minimum specification is PRINT followed by a single keyword.
- If PRINT is used, only the requested output is displayed.

<b>DEFAULT</b>	<i>Goodness-of-fit tests for the model, classification tables, and statistics for the variables in and not in the equation at each step.</i> Tables and statistics are displayed for each split file and METHOD subcommand.
<b>SUMMARY</b>	<i>Summary information.</i> This output is the same output as DEFAULT, except that the output for each step is not displayed.
<b>CORR</b>	<i>Correlation matrix of parameter estimates for the variables in the model.</i>
<b>ITER(value)</b>	<i>Iterations at which parameter estimates are to be displayed.</i> The value in parentheses controls the spacing of iteration reports. If the value is $n$ , the parameter estimates are displayed for every $n$ th iteration, starting at 0. If a value is not supplied, intermediate estimates are displayed at each iteration.
<b>GOODFIT</b>	<i>Hosmer-Lemeshow goodness-of-fit statistic</i> (Hosmer and Lemeshow, 2000).
<b>CI(level)</b>	<i>Confidence interval for <math>\exp(B)</math>.</i> The value in parentheses must be an integer between 1 and 99.
<b>ALL</b>	<i>All available output.</i>

**Example**

```
LOGISTIC REGRESSION VARIABLES=PASS WITH GPA, GRE, MAT
/METHOD FSTEP
/PRINT CORR SUMMARY ITER(2).
```

- A forward stepwise logistic regression analysis of *PASS* on *GPA*, *GRE*, and *MAT* is specified.
- The `PRINT` subcommand requests the display of the correlation matrix of parameter estimates for the variables in the model (`CORR`), classification tables and statistics for the variables in and not in the equation for the final model (`SUMMARY`), and parameter estimates at every second iteration (`ITER(2)`).

**CRITERIA Subcommand**

`CRITERIA` controls the statistical criteria that are used in building the logistic regression models. The way in which these criteria are used depends on the method that is specified on the `METHOD` subcommand. The default criteria are noted in the description of each keyword below. Iterations will stop if the criterion for `BCON`, `LCON`, or `ITERATE` is satisfied.

<b>BCON(value)</b>	<i>Change in parameter estimates to terminate iteration.</i> Iteration terminates when the parameters change by less than the specified value. The default is 0.001. To eliminate this criterion, specify a value of 0.
<b>ITERATE</b>	<i>Maximum number of iterations.</i> The default is 20.
<b>LCON(value)</b>	<i>Percentage change in the log-likelihood ratio for termination of iterations.</i> If the log-likelihood decreases by less than the specified value, iteration terminates. The default is 0, which is equivalent to not using this criterion.
<b>PIN(value)</b>	<i>Probability of score statistic for variable entry.</i> The default is 0.05. The larger the specified probability, the easier it is for a variable to enter the model.
<b>POUT(value)</b>	<i>Probability of conditional, Wald, or LR statistic to remove a variable.</i> The default is 0.1. The larger the specified probability, the easier it is for a variable to remain in the model.
<b>EPS(value)</b>	<i>Epsilon value used for redundancy checking.</i> The specified value must be less than or equal to 0.05 and greater than or equal to $10^{-12}$ . The default is $10^{-8}$ . Larger values make it harder for variables to pass the redundancy check—that is, they are more likely to be removed from the analysis.
<b>CUT(value)</b>	<i>Cutoff value for classification.</i> A case is assigned to a group when the predicted event probability is greater than or equal to the cutoff value. The cutoff value affects the value of the dichotomous derived variable in the classification table, the predicted group ( <code>PGROUP</code> on <code>CASEWISE</code> ), and the classification plot ( <code>CLASSPLOT</code> ). The default cutoff value is 0.5. You can specify a value between 0 and 1 ( $0 < \text{value} < 1$ ).

**Example**

```
LOGISTIC REGRESSION VARIABLES = PROMOTED WITH AGE JOBTIME RACE
/CATEGORICAL RACE
/METHOD BSTEP
/CRITERIA BCON(0.01) PIN(0.01) POUT(0.05).
```

- A backward stepwise logistic regression analysis is performed for the dependent variable *PROMOTED* and the independent variables *AGE*, *JOBTIME*, and *RACE*.

- **CRITERIA** alters four of the statistical criteria that control the building of a model.
- **BCON** specifies that if the change in the absolute value of all of the parameter estimates is less than 0.01, the iterative estimation process should stop. Larger values lower the number of required iterations. Notice that the **ITER** and **LCON** criteria remain unchanged and that if either of them is met before **BCON**, iterations will terminate. (**LCON** can be set to 0 if only **BCON** and **ITER** are to be used.)
- **POUT** requires that the probability of the statistic that is used to test whether a variable should remain in the model be smaller than 0.05. This requirement is more stringent than the default value of 0.1.
- **PIN** requires that the probability of the score statistic that is used to test whether a variable should be included be smaller than 0.01. This requirement makes it more difficult for variables to be included in the model than the default value of 0.05.

## **CLASSPLOT Subcommand**

**CLASSPLOT** generates a classification plot of the actual and predicted values of the dichotomous dependent variable at each step.

- Keyword **CLASSPLOT** is the only specification.
- If **CLASSPLOT** is not specified, plots are not generated.

### **Example**

```
LOGISTIC REGRESSION VARIABLES = PROMOTED WITH JOBTIME RACE
/CATEGORICAL RACE
/CLASSPLOT.
```

- A logistic regression model is constructed for the dichotomous dependent variable *PROMOTED* and the independent variables *JOBTIME* and *RACE*.
- **CLASSPLOT** produces a classification plot for the dependent variable *PROMOTED*. The vertical axis of the plot is the frequency of the variable *PROMOTED*. The horizontal axis is the predicted probability of membership in the second of the two levels of *PROMOTED*.

## **CASEWISE Subcommand**

**CASEWISE** produces a casewise listing of the values of the temporary variables that are created by **LOGISTIC REGRESSION**.

The following keywords are available for specifying temporary variables (see Fox, 1984). When **CASEWISE** is specified by itself, the default is to list *PRED*, *PGROUP*, *RESID*, and *ZRESID*. If a list of variable names is given, only those named temporary variables are displayed.

<b>PRED</b>	<i>Predicted probability.</i> For each case, the predicted probability of having the second of the two values of the dichotomous dependent variable.
<b>PGROUP</b>	<i>Predicted group.</i> The group to which a case is assigned based on the predicted probability.
<b>RESID</b>	<i>Difference between observed and predicted probabilities.</i>
<b>DEV</b>	<i>Deviance values.</i> For each case, a log-likelihood-ratio statistic, which measures how well the model fits the case, is computed.

<b>LRESID</b>	<i>Logit residual.</i> Residual divided by the product of <i>PRED</i> and $1-PRED$ .
<b>SRESID</b>	<i>Studentized residual.</i>
<b>ZRESID</b>	<i>Normalized residual.</i> Residual divided by the square root of the product of <i>PRED</i> and $1-PRED$ .
<b>LEVER</b>	<i>Leverage value.</i> A measure of the relative influence of each observation on the model's fit.
<b>COOK</b>	<i>Analog of Cook's influence statistic.</i>
<b>DFBETA</b>	<i>Difference in beta.</i> The difference in the estimated coefficients for each independent variable if the case is omitted.

The following keyword is available for restricting the cases to be displayed, based on the absolute value of *SRESID*:

<b>OUTLIER (value)</b>	<i>Cases with absolute values of SRESID greater than or equal to the specified value are displayed.</i> If <b>OUTLIER</b> is specified with no value, the default is 2.
------------------------	---

### Example

```
LOGISTIC REGRESSION VARIABLES = PROMOTED WITH JOBTIME SEX RACE
/CATEGORICAL SEX RACE
/METHOD ENTER
/CASEWISE SRESID LEVER DFBETA.
```

- **CASEWISE** produces a casewise listing of the temporary variables *SRESID*, *LEVER*, and *DFBETA*.
- There will be one *DFBETA* value for each parameter in the model. The continuous variable *JOBTIME*, the two-level categorical variable *SEX*, and the constant each require one parameter, while the four-level categorical variable *RACE* requires three parameters. Thus, six values of *DFBETA* will be produced for each case.

## MISSING Subcommand

**LOGISTIC REGRESSION** excludes all cases with missing values on any of the independent variables. For a case with a missing value on the dependent variable, predicted values are calculated if it has nonmissing values on all independent variables. The **MISSING** subcommand controls the processing of user-missing values. If the subcommand is not specified, the default is **EXCLUDE**.

<b>EXCLUDE</b>	<i>Delete cases with user-missing values as well as system-missing values.</i> This setting is the default.
<b>INCLUDE</b>	<i>Include user-missing values in the analysis.</i>

## OUTFILE Subcommand

The **OUTFILE** subcommand allows you to specify files to which output is written.

- Only one **OUTFILE** subcommand is allowed. If you specify more than one subcommand, only the last subcommand is executed.

- You must specify at least one keyword and a valid filename in parentheses. There is no default.
- `MODEL` cannot be used if split-file processing is on (`SPLIT FILE` command) or if more than one dependent variable is specified (`DEPENDENT` subcommand).

**MODEL(filename)**      *Write parameter estimates and their covariances to an XML file.* Specify the filename in full. `LOGISTIC REGRESSION` does not supply an extension. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#)

**PARAMETER(filename)**      *Write parameter estimates only to an XML file.* Specify the filename in full. `LOGISTIC REGRESSION` does not supply an extension. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#)

## **SAVE Subcommand**

`SAVE` saves the temporary variables that are created by `LOGISTIC REGRESSION`. To specify variable names for the new variables, assign the new names in parentheses following each temporary variable name. If new variable names are not specified, `LOGISTIC REGRESSION` generates default names.

- Assigned variable names must be unique in the active dataset. Scratch or system variable names (that is, names that begin with # or \$) cannot be used.
- A temporary variable can be saved only once on the same `SAVE` subcommand.

### **Example**

```
LOGISTIC REGRESSION VARIABLES = PROMOTED WITH JOBTIME AGE
  /SAVE PRED (PREDPRO) DFBETA (DF) .
```

- A logistic regression analysis of *PROMOTED* on the independent variables *JOBTIME* and *AGE* is performed.
- `SAVE` adds four variables to the active dataset: one variable named *PREDPRO*, containing the predicted value from the specified model for each case, and three variables named *DF0*, *DF1*, and *DF2*, containing, respectively, the *DFBETA* values for each case of the constant, the independent variable *JOBTIME*, and the independent variable *AGE*.

## **EXTERNAL Subcommand**

`EXTERNAL` indicates that the data for each split-file group should be held in an external scratch file during processing. This process can help conserve memory resources when running complex analyses or analyses with large data sets.

- The keyword `EXTERNAL` is the only specification.
- Specifying `EXTERNAL` may result in slightly longer processing time.
- If `EXTERNAL` is not specified, all data are held internally, and no scratch file is written.

## **References**

- Agresti, A. 2002. *Categorical Data Analysis*, 2nd ed. New York: John Wiley and Sons.
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- Fox, J. 1984. *Linear statistical models and related methods: With applications to social research*. New York: John Wiley and Sons.
- Hosmer, D. W., and S. Lemeshow. 2000. *Applied Logistic Regression*, 2nd ed. New York: John Wiley and Sons.
- Kirk, R. E. 1982. *Experimental design*, 2nd ed. Monterey, California: Brooks/Cole.
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# LOGLINEAR

LOGLINEAR is available in the Advanced Statistics option.

The syntax for LOGLINEAR is available only in a syntax window, not from the dialog box interface. See [GENLOG](#) for information on loglinear models available from the dialog box interface.

```
LOGLINEAR varlist(min,max)...[BY] varlist(min,max)
        [WITH covariate varlist]
[/CWEIGHT={varname } ] [/CWEIGHT=(matrix)...]
        { (matrix) }
[/GRESID={varlist } ] [/GRESID=(matrix)...]
        { (matrix) }
[/CONTRAST (varname)={DEVIATION [(refcat)]          } [/CONTRAST...]]
        {DIFFERENCE                               }
        {HELMERT                                   }
        {SIMPLE [(refcat)]                         }
        {REPEATED                                  }
        {POLYNOMIAL [({1,2,3,...})]}
        {          {metric } }
        {[BASIS] SPECIAL(matrix) }
[/CRITERIA=[CONVERGE({0.001**})] [ITERATE({20**})] [DELTA({0.5**})]]
        {n          }          {n          }          {n          }
        [DEFAULT]]
[/PRINT={ [FREQ**] [RESID**] [DESIGN] [ESTIM] [COR] }]
        {DEFAULT          }
        {ALL              }
        {NONE             }
[/PLOT={NONE** } ]
        {DEFAULT }
        {RESID  }
        {NORMPROB}
[/MISSING=[ {EXCLUDE**}]]
        {INCLUDE  }
[/DESIGN=effect[(n)] effect[(n)]... effect BY effect...] [/DESIGN...]
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
LOGLINEAR JOBSAT (1,2) ZODIAC (1,12) /DESIGN=JOBSAT.
```

## Overview

LOGLINEAR is a general procedure for model fitting, hypothesis testing, and parameter estimation for any model that has categorical variables as its major components. As such, LOGLINEAR subsumes a variety of related techniques, including general models of multiway contingency tables, logit models, logistic regression on categorical variables, and quasi-independence models.

LOGLINEAR models cell frequencies using the multinomial response model and produces maximum likelihood estimates of parameters by means of the Newton-Raphson algorithm (Haberman, 1978). HILOGLINEAR, which uses an iterative proportional-fitting algorithm, is more efficient for hierarchical models, but it cannot produce parameter estimates for unsaturated models, does not permit specification of contrasts for parameters, and does not display a correlation matrix of the parameter estimates.

### **Comparison of the GENLOG and LOGLINEAR Commands**

The General Loglinear Analysis and Logit Loglinear Analysis dialog boxes are both associated with the GENLOG command. In previous releases, these dialog boxes were associated with the LOGLINEAR command. The LOGLINEAR command is now available only as a syntax command. The differences are described below.

#### **Distribution Assumptions**

- GENLOG can handle both Poisson and multinomial distribution assumptions for observed cell counts.
- LOGLINEAR assumes only multinomial distribution.

#### **Approach**

- GENLOG uses a regression approach to parameterize a categorical variable in a design matrix.
- LOGLINEAR uses contrasts to reparameterize a categorical variable. The major disadvantage of the reparameterization approach is in the interpretation of the results when there is a redundancy in the corresponding design matrix. Also, the reparameterization approach may result in incorrect degrees of freedom for an incomplete table, leading to incorrect analysis results.

#### **Contrasts and Generalized Log-Odds Ratios (GLOR)**

- GENLOG doesn't provide contrasts to reparameterize the categories of a factor. However, it offers generalized log-odds ratios (GLOR) for cell combinations. Often, comparisons among categories of factors can be derived from GLOR.
- LOGLINEAR offers contrasts to reparameterize the categories of a factor.

#### **Deviance Residual**

- GENLOG calculates and displays the deviance residual and its normal probability plot in addition to the other residuals.
- LOGLINEAR does not calculate the deviance residual.

#### **Factor-by-Covariate Design**

- When there is a factor-by-covariate term in the design, GENLOG generates one regression coefficient of the covariate for each combination of factor values. The estimates of these regression coefficients are calculated and displayed.
- LOGLINEAR estimates and displays the contrasts of these regression coefficients.

### Partition Effect

- In GENLOG, the term *partition effect* refers to the category of a factor.
- In LOGLINEAR, the term *partition effect* refers to a particular contrast.

### Options

**Model Specification.** You can specify the model or models to be fit using the DESIGN subcommand.

**Cell Weights.** You can specify cell weights, such as structural zeros, for the model with the CWEIGHT subcommand.

**Output Display.** You can control the output display with the PRINT subcommand.

**Optional Plots.** You can produce plots of adjusted residuals against observed and expected counts, normal plots, and detrended normal plots with the PLOT subcommand.

**Linear Combinations.** You can calculate linear combinations of observed cell frequencies, expected cell frequencies, and adjusted residuals using the GRESID subcommand.

**Contrasts.** You can indicate the type of contrast desired for a factor using the CONTRAST subcommand.

**Criteria for Algorithm.** You can control the values of algorithm-tuning parameters with the CRITERIA subcommand.

### Basic Specification

The basic specification is two or more variables that define the crosstabulation. The minimum and maximum values for each variable must be specified in parentheses after the variable name.

By default, LOGLINEAR estimates the saturated model for a multidimensional table. Output includes the factors or effects, their levels, and any labels; observed and expected frequencies and percentages for each factor and code; residuals, standardized residuals, and adjusted residuals; two goodness-of-fit statistics (the likelihood-ratio chi-square and Pearson's chi-square); and estimates of the parameters with accompanying  $z$  values and 95% confidence intervals.

### Limitations

- A maximum of 10 independent (factor) variables
- A maximum of 200 covariates

### Subcommand Order

- The variables specification must come first.
- The subcommands that affect a specific model must be placed before the DESIGN subcommand specifying the model.

- All subcommands can be used more than once and, with the exception of the `DESIGN` subcommand, are carried from model to model unless explicitly overridden.
- If the last subcommand is not `DESIGN`, `LOGLINEAR` generates a saturated model in addition to the explicitly requested model(s).

## Examples

### Example: Main Effects General Loglinear Model

```
LOGLINEAR JOBSAT (1,2) ZODIAC (1,12) /DESIGN=JOBSAT, ZODIAC.
```

- The variable list specifies two categorical variables, *JOBSAT* and *ZODIAC*. *JOBSAT* has values 1 and 2. *ZODIAC* has values 1 through 12.
- `DESIGN` specifies a model with main effects only.

### Example: Saturated General Loglinear Model

```
LOGLINEAR DPREF (2,3) RACE CAMP (1,2).
```

- *DPREF* is a categorical variable with values 2 and 3. *RACE* and *CAMP* are categorical variables with values 1 and 2.
- This is a general loglinear model because no `BY` keyword appears. The design defaults to a saturated model that includes all main effects and interaction effects.

### Example: Logit Loglinear Model

```
LOGLINEAR GSLEVEL (4,8) BY EDUC (1,4) SEX (1,2)
/DESIGN=GSLEVEL, GSLEVEL BY EDUC, GSLEVEL BY SEX.
```

- *GSLEVEL* is a categorical variable with values 4 through 8. *EDUC* is a categorical variable with values 1 through 4. *SEX* has values 1 and 2.
- The keyword `BY` on the variable list specifies a logit model in which *GSLEVEL* is the dependent variable and *EDUC* and *SEX* are the independent variables.
- `DESIGN` specifies a model that can test for the absence of a joint effect of *SEX* and *EDUC* on *GSLEVEL*.

## Variable List

The variable list specifies the variables to be included in the model. `LOGLINEAR` analyzes two classes of variables: categorical and continuous. Categorical variables are used to define the cells of the table. Continuous variables are used as cell covariates. Continuous variables can be specified only after the keyword `WITH` following the list of categorical variables.

- The list of categorical variables must be specified first. Categorical variables must be numeric and integer.
- A range must be defined for each categorical variable by specifying, in parentheses after each variable name, the minimum and maximum values for that variable. Separate the two values with at least one space or a comma.

- To specify the same range for a list of variables, specify the list of variables followed by a single range. The range applies to all variables on the list.
- To specify a logit model, use the keyword `BY` (see [Logit Model on p. 1047](#)). A variable list without the keyword `BY` generates a general loglinear model.
- Cases with values outside the specified range are excluded from the analysis. Non-integer values within the range are truncated for the purpose of building the table.

### **Logit Model**

- To segregate the independent (factor) variables from the dependent variables in a logit model, use the keyword `BY`. The categorical variables preceding `BY` are the dependent variables; the categorical variables following `BY` are the independent variables.
- A total of 10 categorical variables can be specified. In most cases, one of them is dependent.
- A `DESIGN` subcommand should be used to request the desired logit model.
- `LOGLINEAR` displays an analysis of dispersion and two measures of association: entropy and concentration. These measures are discussed elsewhere (Haberman, 1982) and can be used to quantify the magnitude of association among the variables. Both are proportional reduction in error measures. The entropy statistic is analogous to Theil's entropy measure, while the concentration statistic is analogous to Goodman and Kruskal's tau-*b*. Both statistics measure the strength of association between the dependent variable and the predictor variable set.

### **Cell Covariates**

- Continuous variables can be used as covariates. When used, the covariates must be specified after the keyword `WITH` following the list of categorical variables. Ranges are not specified for the continuous variables.
- A variable cannot be named as both a categorical variable and a cell covariate.
- To enter cell covariates into a model, the covariates must be specified on the `DESIGN` subcommand.
- Cell covariates are not applied on a case-by-case basis. The mean covariate value for a cell in the contingency table is applied to that cell.

#### **Example**

```
LOGLINEAR DPREF(2,3) RACE CAMP (1,2) WITH CONSTANT
  /DESIGN=DPREF RACE CAMP CONSTANT.
```

- The variable `CONSTANT` is a continuous variable specified as a cell covariate. Cell covariates must be specified after the keyword `WITH` following the variable list. No range is defined for cell covariates.
- To include the cell covariate in the model, the variable `CONSTANT` is specified on `DESIGN`.

### **CWEIGHT Subcommand**

`CWEIGHT` specifies cell weights, such as structural zeros, for a model. By default, cell weights are equal to 1.

- The specification is either one numeric variable or a matrix of weights enclosed in parentheses.
- If a matrix of weights is specified, the matrix must contain the same number of elements as the product of the levels of the categorical variables. An asterisk can be used to signify repetitions of the same value.
- If weights are specified for a multiple-factor model, the index value of the rightmost factor increments the most rapidly.
- If a numeric variable is specified, only one `CWEIGHT` subcommand can be used on `LOGLINEAR`.
- To use multiple cell weights on the same `LOGLINEAR` command, specify all weights in matrix format. Each matrix must be specified on a separate `CWEIGHT` subcommand, and each `CWEIGHT` specification remains in effect until explicitly overridden by another `CWEIGHT` subcommand.
- `CWEIGHT` can be used to impose structural, or *a priori*, zeros on the model. This feature is useful in the analysis of symmetric tables.

### Example

```
COMPUTE CWT=1.
IF (HUSED EQ WIFED) CWT=0.
LOGLINEAR HUSED WIFED(1,4) WITH DISTANCE
  /CWEIGHT=CWT
  /DESIGN=HUSED WIFED DISTANCE.
```

- `COMPUTE` initially assigns `CWT` the value 1 for all cases.
- `IF` assigns `CWT` the value 0 when `HUSED` equals `WIFED`.
- `CWEIGHT` imposes structural zeros on the diagonal of the symmetric crosstabulation. Because a variable name is specified, only one `CWEIGHT` can be used.

### Example

```
LOGLINEAR HUSED WIFED(1,4) WITH DISTANCE
  /CWEIGHT=(0, 4*1, 0, 4*1, 0, 4*1, 0)
  /DESIGN=HUSED WIFED DISTANCE
  /CWEIGHT=(16*1)
  /DESIGN=HUSED WIFED DISTANCE.
```

- The first `CWEIGHT` matrix specifies the same values as variable `CWT` provided in the first example. The specified matrix is as follows:

```
0 1 1 1
1 0 1 1
1 1 0 1
1 1 1 0
```

- The same matrix can be specified in full as (0 1 1 1 1 0 1 1 1 1 0 1 1 1 1 0).
- By using the matrix format on `CWEIGHT` rather than a variable name, a different `CWEIGHT` subcommand can be used for the second model.

## ***GRESID Subcommand***

**GRESID** (generalized residual) calculates linear combinations of observed cell frequencies, expected cell frequencies, and adjusted residuals.

- The specification is either a numeric variable or a matrix whose contents are coefficients of the desired linear combinations.
- If a matrix of coefficients is specified, the matrix must contain the same number of elements as the number of cells implied by the variables specification. An asterisk can be used to signify repetitions of the same value.
- Each **GRESID** subcommand specifies a single linear combination. Each matrix or variable must be specified on a separate **GRESID** subcommand. All **GRESID** subcommands specified are displayed for each design.

### ***Example***

```
LOGLINEAR MONTH(1,18) WITH Z
  /GRESID=(6*1,12*0)
  /GRESID=(6*0,6*1,6*0)
  /GRESID=(12*0,6*1)
  /DESIGN=Z.
```

- The first **GRESID** subcommand combines the first six months into a single effect. The second **GRESID** subcommand combines the second six months, and the third **GRESID** subcommand combines the last six months.
- For each effect, **LOGLINEAR** displays the observed and expected counts, the residual, and the adjusted residual.

## ***CONTRAST Subcommand***

**CONTRAST** indicates the type of contrast desired for a factor, where a factor is any categorical dependent or independent variable. The default contrast is **DEVIATION** for each factor.

- The specification is **CONTRAST**, which is followed by a variable name in parentheses and the contrast-type keyword.
- To specify a contrast for more than one factor, use a separate **CONTRAST** subcommand for each specified factor. Only one contrast can be in effect for each factor on each **DESIGN**.
- A contrast specification remains in effect for subsequent designs until explicitly overridden by another **CONTRAST** subcommand.
- The design matrix used for the contrasts can be displayed by specifying the keyword **DESIGN** on the **PRINT** subcommand. However, this matrix is the basis matrix that is used to determine contrasts; it is not the contrast matrix itself.
- **CONTRAST** can be used for a multinomial logit model, in which the dependent variable has more than two categories.
- **CONTRAST** can be used for fitting linear logit models. The keyword **BASIS** is not appropriate for such models.
- In a logit model, **CONTRAST** is used to transform the independent variable into a metric variable. Again, the keyword **BASIS** is not appropriate.

The following contrast types are available:

<b>DEVIATION(refcat)</b>	<i>Deviations from the overall effect.</i> <code>DEVIATION</code> is the default contrast if the <code>CONTRAST</code> subcommand is not used. Refcat is the category for which parameter estimates are not displayed (they are the negative of the sum of the others). By default, refcat is the last category of the variable.
<b>DIFFERENCE</b>	<i>Levels of a factor with the average effect of previous levels of a factor.</i> Also known as reverse Helmert contrasts.
<b>HELMERT</b>	<i>Levels of a factor with the average effect of subsequent levels of a factor.</i>
<b>SIMPLE(refcat)</b>	<i>Each level of a factor to the reference level.</i> By default, <code>LOGLINEAR</code> uses the last category of the factor variable as the reference category. Optionally, any level can be specified as the reference category enclosed in parentheses after the keyword <code>SIMPLE</code> . The sequence of the level, not the actual value, must be specified.
<b>REPEATED</b>	<i>Adjacent comparisons across levels of a factor.</i>
<b>POLYNOMIAL(metric)</b>	<i>Orthogonal polynomial contrasts.</i> The default is equal spacing. Optionally, the coefficients of the linear polynomial can be specified in parentheses, indicating the spacing between levels of the treatment measured by the given factor.
<b>[BASIS]SPECIAL(matrix)</b>	<i>User-defined contrast.</i> As many elements as the number of categories squared must be specified. If <code>BASIS</code> is specified before <code>SPECIAL</code> , a basis matrix is generated for the special contrast, which makes the coefficients of the contrast equal to the special matrix. Otherwise, the matrix specified is transposed and then used as the basis matrix to determine coefficients for the contrast matrix.

### Example

```
LOGLINEAR  A(1,4) BY B(1,4)
/CONTRAST(B)=POLYNOMIAL
/DESIGN=A A BY B(1)
/CONTRAST(B)=SIMPLE
/DESIGN=A A BY B(1).
```

- The first `CONTRAST` subcommand requests polynomial contrasts of *B* for the first design.
- The second `CONTRAST` subcommand requests the simple contrast of *B*, with the last category (value 4) used as the reference category for the second `DESIGN` subcommand.

### Example

```
* Multinomial logit model
LOGLINEAR  PREF(1,5) BY RACE ORIGIN CAMP(1,2)
/CONTRAST(PREF)=SPECIAL(5*1, 1 1 1 1 -4, 3 -1 -1 -1 0,
0 1 1 -2 0, 0 1 -1 0 0).
```

- `LOGLINEAR` builds special contrasts among the five categories of the dependent variable *PREF*, which measures preference for training camps among Army recruits. For *PREF*, 1=stay, 2=move to north, 3=move to south, 4=move to unnamed camp, and 5=undecided.
- The four contrasts are: (1) move or stay versus undecided, (2) stay versus move, (3) named camp versus unnamed, and (4) northern camp versus southern. Because these contrasts are orthogonal, `SPECIAL` and `BASIS SPECIAL` produce equivalent results.



**Example**

```
* Contrasts for a linear logit model

LOGLINEAR RESPONSE(1,2) BY YEAR(0,20)
/PRINT=DEFAULT ESTIM
/CONTRAST(YEAR)=SPECIAL(21*1, -10, -9, -8, -7, -6, -5, -4,
                        -3, -2, -1, 0, 1, 2, 3, 4, 5, 6, 7,
                        8, 9, 10, 399*1)
/DESIGN=RESPONSE RESPONSE BY YEAR(1).
```

- *YEAR* measures years of education and ranges from 0 through 20. Therefore, allowing for the constant effect, *YEAR* has 20 estimable parameters associated with it.
- The *SPECIAL* contrast specifies the constant—that is, 21\*1—and the linear effect of *YEAR*—that is, -10 to 10. The other 399 1's fill out the 21\*21 matrix.

**Example**

```
* Contrasts for a logistic regression model

LOGLINEAR RESPONSE(1,2) BY TIME(1,4)
/CONTRAST(TIME) = SPECIAL(4*1, 7 14 27 51, 8*1)
/PRINT=ALL /PLOT=DEFAULT
/DESIGN=RESPONSE, TIME(1) BY RESPONSE.
```

- *CONTRAST* is used to transform the independent variable into a metric variable.
- *TIME* represents elapsed time in days. Therefore, the weights in the contrast represent the metric of the passage of time.

**CRITERIA Subcommand**

*CRITERIA* specifies the values of some constants in the Newton-Raphson algorithm. Defaults or specifications remain in effect until overridden with another *CRITERIA* subcommand.

<b>CONVERGE(n)</b>	<i>Convergence criterion.</i> Specify a value for the convergence criterion. The default is 0.001.
<b>ITERATE(n)</b>	<i>Maximum number of iterations.</i> Specify the maximum number of iterations for the algorithm. The default number is 20.
<b>DELTA(n)</b>	<i>Cell delta value.</i> The value of delta is added to each cell frequency for the first iteration. For saturated models, it remains in the cell. The default value is 0.5. <i>LOGLINEAR</i> does not display parameter estimates or correlation matrices of parameter estimates if any sampling zero cells exist in the expected table after delta is added. Parameter estimates and correlation matrices can be displayed in the presence of structural zeros.
<b>DEFAULT</b>	<i>Default values are used.</i> <i>DEFAULT</i> can be used to reset the parameters to the default.

**Example**

```
LOGLINEAR DPREF(2,3) BY RACE ORIGIN CAMP(1,2)
/CRITERIA=ITERATION(50) CONVERGE(.0001).
```

- *ITERATION* increases the maximum number of iterations to 50.
- *CONVERGE* lowers the convergence criterion to 0.0001.

## **PRINT Subcommand**

PRINT requests statistics that are not produced by default.

- By default, LOGLINEAR displays the frequency table and residuals. The parameter estimates of the model are also displayed if DESIGN is not used.
- Multiple PRINT subcommands are permitted. The specifications are cumulative.

The following keywords can be used on PRINT:

<b>FREQ</b>	<i>Observed and expected cell frequencies and percentages.</i> This is displayed by default.
<b>RESID</b>	<i>Raw, standardized, and adjusted residuals.</i> This is displayed by default.
<b>DESIGN</b>	<i>The design matrix of the model, showing the basis matrix corresponding to the contrasts used.</i>
<b>ESTIM</b>	<i>The parameter estimates of the model.</i> If you do not specify a design on the DESIGN subcommand, LOGLINEAR generates a saturated model and displays the parameter estimates for the saturated model. LOGLINEAR does not display parameter estimates or correlation matrices of parameter estimates if any sampling zero cells exist in the expected table after delta is added. Parameter estimates and a correlation matrix are displayed when structural zeros are present.
<b>COR</b>	<i>The correlation matrix of the parameter estimates.</i> Alias COV.
<b>ALL</b>	<i>All available output.</i>
<b>DEFAULT</b>	<i>FREQ and RESID.</i> ESTIM is also displayed by default if the DESIGN subcommand is not used.
<b>NONE</b>	<i>The design information and goodness-of-fit statistics only.</i> This option overrides all other specifications on the PRINT subcommand. The NONE option applies only to the PRINT subcommand.

### **Example**

```
LOGLINEAR A(1,2) B(1,2)
  /PRINT=ESTIM
  /DESIGN=A,B,A BY B
  /PRINT=ALL
  /DESIGN=A,B.
```

- The first design is the saturated model. The parameter estimates are displayed with ESTIM specified on PRINT.
- The second design is the main-effects model, which tests the hypothesis of no interaction. The second PRINT subcommand displays all available display output for this model.

## **PLOT Subcommand**

PLOT produces optional plots. No plots are displayed if PLOT is not specified or is specified without any keyword. Multiple PLOT subcommands can be used. The specifications are cumulative.

<b>RESID</b>	<i>Plots of adjusted residuals against observed and expected counts.</i>
<b>NORMPROB</b>	<i>Normal and detrended normal plots of the adjusted residuals.</i>
<b>NONE</b>	<i>No plots.</i>
<b>DEFAULT</b>	<i>RESID and NORMPROB.</i> Alias ALL.

**Example**

```
LOGLINEAR RESPONSE(1,2) BY TIME(1,4)
  /CONTRAST(TIME)=SPECIAL(4*1, 7 14 27 51, 8*1)
  /PLOT=DEFAULT
  /DESIGN=RESPONSE TIME(1) BY RESPONSE
  /PLOT=NONE
  /DESIGN.
```

- RESID and NORMPROB plots are displayed for the first design.
- No plots are displayed for the second design.

**MISSING Subcommand**

MISSING controls missing values. By default, LOGLINEAR excludes all cases with system- or user-missing values on any variable. You can specify INCLUDE to include user-missing values. If INCLUDE is specified, user-missing values must also be included in the value range specification.

- EXCLUDE**            *Delete cases with user-missing values.* This is the default if the subcommand is omitted. You can also specify the keyword DEFAULT.
- INCLUDE**            *Include user-missing values.* Only cases with system-missing values are deleted.

**Example**

```
MISSING VALUES A(0).
LOGLINEAR A(0,2) B(1,2) /MISSING=INCLUDE
  /DESIGN=B.
```

- Even though 0 was specified as missing, it is treated as a nonmissing category of *A* in this analysis.

**DESIGN Subcommand**

DESIGN specifies the model or models to be fit. If DESIGN is omitted or used with no specifications, the saturated model is produced. The saturated model fits all main effects and all interaction effects.

- To specify more than one model, use more than one DESIGN subcommand. Each DESIGN specifies one model.
- To obtain main-effects models, name all the variables listed on the variables specification.
- To obtain interactions, use the keyword BY to specify each interaction, as in A BY B and C BY D. To obtain the single-degree-of-freedom partition of a specified contrast, specify the partition in parentheses following the factor (see the example below).
- To include cell covariates in the model, first identify them on the variable list by naming them after the keyword WITH, and then specify the variable names on DESIGN.
- To specify an equiprobability model, name a cell covariate that is actually a constant of 1.

**Example**

```
* Testing the linear effect of the dependent variable
```

```
COMPUTE X=MONTH.
LOGLINEAR MONTH (1,12) WITH X
  /DESIGN X.
```

- The variable specification identifies *MONTH* as a categorical variable with values 1 through 12. The keyword *WITH* identifies *X* as a covariate.
- *DESIGN* tests the linear effect of *MONTH*.

**Example**

\* Specifying main effects models

```
LOGLINEAR A(1,4) B(1,5)
  /DESIGN=A
  /DESIGN=A,B.
```

- The first design tests the homogeneity of category probabilities for *B*; it fits the marginal frequencies on *A*, but assumes that membership in any of the categories of *B* is equiprobable.
- The second design tests the independence of *A* and *B*. It fits the marginals on both *A* and *B*.

**Example**

\* Specifying interactions

```
LOGLINEAR A(1,4) B(1,5) C(1,3)
  /DESIGN=A,B,C, A BY B.
```

- This design consists of the *A* main effect, the *B* main effect, the *C* main effect, and the interaction of *A* and *B*.

**Example**

\* Single-degree-of-freedom partitions

```
LOGLINEAR A(1,4) BY B(1,5)
  /CONTRAST(B)=POLYNOMIAL
  /DESIGN=A,A BY B(1).
```

- The value 1 following *B* refers to the first partition of *B*, which is the linear effect of *B*; this follows from the contrast specified on the *CONTRAST* subcommand.

**Example**

\* Specifying cell covariates

```
LOGLINEAR HUSED WIFED(1,4) WITH DISTANCE
  /DESIGN=HUSED WIFED DISTANCE.
```

- The continuous variable *DISTANCE* is identified as a cell covariate by specifying it after *WITH* on the variable list. The cell covariate is then included in the model by naming it on *DESIGN*.

**Example**

\* Equiprobability model

```
COMPUTE X=1.
```

```
LOGLINEAR MONTH(1,18) WITH X  
/DESIGN=X.
```

- This model tests whether the frequencies in the 18-cell table are equal by using a cell covariate that is a constant of 1.

# LOOP-END LOOP

```
LOOP [varname=n TO m [BY {1**}]] [IF [(logical expression)]]
      {n }
transformation commands
END LOOP [IF [(logical expression)]]
```

**\*\*Default if the subcommand is omitted.**

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## Examples

```
SET MXLOOPS=10. /*Maximum number of loops allowed
LOOP. /*Loop with no limit other than MXLOOPS
COMPUTE X=X+1.
END LOOP.
```

```
LOOP #I=1 TO 5. /*Loop five times
COMPUTE X=X+1.
END LOOP.
```

## Overview

The LOOP-END LOOP structure performs repeated transformations specified by the commands within the loop until they reach a specified cutoff. The cutoff can be specified by an indexing clause on the LOOP command, an IF clause on the END LOOP command, or a BREAK command within the loop structure (see BREAK). In addition, the maximum number of iterations within a loop can be specified on the MXLOOPS subcommand on SET. The default MXLOOPS is 40.

The IF clause on the LOOP command can be used to perform repeated transformations on a subset of cases. The effect is similar to nesting the LOOP-END LOOP structure within a DO IF-END IF structure, but using IF on LOOP is simpler and more efficient. You have to use the DO IF-END IF structure, however, if you want to perform different transformations on different subsets of cases. You can also use IF on LOOP to specify the cutoff, especially when the cutoff may be reached before the first iteration.

LOOP and END LOOP are usually used within an input program or with the VECTOR command. Since the loop structure repeats transformations on a single case or on a single input record containing information on multiple cases, it allows you to read complex data files or to generate data for a active dataset. For more information, see INPUT PROGRAM-END INPUT PROGRAM and VECTOR.

The loop structure repeats transformations on single cases across variables. It is different from the DO REPEAT-END REPEAT structure, which replicates transformations on a specified set of variables. When both can be used to accomplish a task, such as selectively transforming data for some cases on some variables, LOOP and END LOOP are generally more efficient and

more flexible, but `DO REPEAT` allows selection of nonadjacent variables and use of replacement values with different intervals.

### **Options**

**Missing Values.** You can prevent cases with missing values for any of the variables used in the loop structure from entering the loop. [For more information, see the topic Missing Values on p. 1064.](#)

**Creating Data.** A loop structure within an input program can be used to generate data. [For more information, see the topic Creating Data on p. 1065.](#)

**Defining Complex File Structures.** A loop structure within an input program can be used to define complex files that cannot be handled by standard file definition facilities.

### **Basic Specification**

The basic specification is `LOOP` followed by at least one transformation command. The structure must end with the `END LOOP` command. Commands within the loop are executed until the cutoff is reached.

### **Syntax Rules**

- If `LOOP` and `END LOOP` are specified before an active dataset exists, they must be specified within an input program.
- If both an indexing and an `IF` clause are used on `LOOP`, the indexing clause must be first.
- Loop structures can be nested within other loop structures or within `DO IF` structures, and vice versa.

### **Operations**

- The `LOOP` command defines the beginning of a loop structure and the `END LOOP` command defines its end. The `LOOP` command returns control to `LOOP` unless the cutoff has been reached. When the cutoff has been reached, control passes to the command immediately following `END LOOP`.
- Commands that take effect immediately without reading the data or executing pending commands (e.g., `VALUE LABELS`, `MISSING VALUES`, `VARIABLE LABEL`), commands that don't require an active dataset (e.g., `DISPLAY`, `HOST`, `INSERT`, `OMS`, `SET`) and the `GET` command are processed unconditionally and immediately each time they are encountered in a loop structure. The specified condition is ignored for these commands.
- Commands that require a data pass, such as statistical or charting procedures, cannot be included in a loop structure and will cause the `LOOP` command to fail.
- An indexing clause (e.g., `LOOP #i=1 to 1000`) will override the `SET MXLOOPS` limit, but a loop with an `IF` condition will terminate if the `MXLOOPS` limit is reached before the condition is satisfied.

## Examples

### Example

```
SET MXLOOPS=10.
LOOP. /*Loop with no limit other than MXLOOPS
COMPUTE X=X+1.
END LOOP.
```

- This and the following examples assume that an active dataset and all of the variables mentioned in the loop exist.
- The `SET MXLOOPS` command limits the number of times the loop is executed to 10. The function of `MXLOOPS` is to prevent infinite loops when there is no indexing clause.
- Within the loop structure, each iteration increments *X* by 1. After 10 iterations, the value of *X* for all cases is increased by 10, and, as specified on the `SET` command, the loop is terminated.

### Example

```
*Assume MXLOOPS set to default value of 40.
COMPUTE newvar1=0.
LOOP IF newvar1<100.
COMPUTE newvar1=newvar1+1.
END LOOP.

PRESERVE.
SET MXLOOPS 500.
COMPUTE newvar2=0.
LOOP IF newvar2<100.
COMPUTE newvar2=newvar2+1.
END LOOP.
RESTORE.

COMPUTE newvar3=0.
LOOP #i=1 to 1000.
COMPUTE newvar3=newvar3+1.
END LOOP.
EXECUTE.
```

- In the first loop, the value of *newvar1* will reach 40, at which point the loop will terminate because the `MXLOOPS` limit has been exceeded.
- In the second loop, the value of `MXLOOPS` is increased to 500, and the loop will continue to iterate until the value of *newvar2* reaches 100, at which point the `IF` condition is reached and the loop terminates.
- In the third loop, the indexing clause overrides the `MXLOOPS` setting, and the loop will iterate 1,000 times.

## IF Keyword

The keyword `IF` and a logical expression can be specified on `LOOP` or on `END LOOP` to control iterations through the loop.

- The specification on `IF` is a logical expression enclosed in parentheses.



**Example**

```
LOOP.
COMPUTE X=X+1.
END LOOP IF (X EQ 5). /*Loop until X is 5
```

- Iterations continue until the logical expression on END LOOP is true, which for every case is when  $X$  equals 5. Each case does not go through the same number of iterations.
- This corresponds to the programming notion of DO UNTIL. The loop is always executed at least once.

**Example**

```
LOOP IF (X LT 5). /*Loop while X is less than 5
COMPUTE X=X+1.
END LOOP.
```

- The IF clause is evaluated each trip through the structure, so looping stops once  $X$  equals 5.
- This corresponds to the programming notion of DO WHILE. The loop may not be executed at all.

**Example**

```
LOOP IF (Y GT 10). /*Loop only for cases with Y GT 10
COMPUTE X=X+1.
END LOOP IF (X EQ 5). /*Loop until X IS 5
```

- The IF clause on LOOP allows transformations to be performed on a subset of cases.  $X$  is increased by 5 only for cases with values greater than 10 for  $Y$ .  $X$  is not changed for all other cases.

## Indexing Clause

The indexing clause limits the number of iterations for a loop by specifying the number of times the program should execute commands within the loop structure. The indexing clause is specified on the LOOP command and includes an indexing variable followed by initial and terminal values.

- The program sets the *indexing variable* to the *initial value* and increases it by the specified increment each time the loop is executed for a case. When the indexing variable reaches the specified *terminal value*, the loop is terminated for that case.
- By default, the program increases the indexing variable by 1 for each iteration. The keyword BY overrides this increment.
- The indexing variable can have any valid variable name. Unless you specify a scratch variable, the indexing variable is treated as a permanent variable and is saved in the active dataset. If the indexing variable is assigned the same name as an existing variable, the values of the existing variable are altered by the LOOP structure as it is executed, and the original values are lost. [For more information, see the topic Creating Data on p. 1065.](#)
- The indexing clause overrides the maximum number of loops specified by SET MXLOOPS.
- The initial and terminal values of the indexing clause can be numeric expressions. Noninteger and negative expressions are allowed.

- If the expression for the initial value is greater than the terminal value, the loop is not executed. For example, #J=X TO Y is a zero-trip loop if X is 0 and Y is -1.
- If the expressions for the initial and terminal values are equal, the loop is executed once. #J=0 TO Y is a one-trip loop when Y is 0.
- If the loop is exited via BREAK or a conditional clause on the END LOOP statement, the iteration variable is not updated. If the LOOP statement contains both an indexing clause and a conditional clause, the indexing clause is executed first, and the iteration variable is updated regardless of which clause causes the loop to terminate.

**Example**

```
LOOP #I=1 TO 5. /*LOOP FIVE TIMES
COMPUTE X=X+1.
END LOOP.
```

- The scratch variable #I (the indexing variable) is set to the initial value of 1 and increased by 1 each time the loop is executed for a case. When #I increases beyond the terminal value 5, no further loops are executed. Thus, the value of X will be increased by 5 for every case.

**Example**

```
LOOP #I=1 TO 5 IF (Y GT 10). /*Loop to X=5 only if Y GT 10
COMPUTE X=X+1.
END LOOP.
```

- Both an indexing clause and an IF clause are specified on LOOP. X is increased by 5 for all cases where Y is greater than 10.

**Example**

```
LOOP #I=1 TO Y. /*Loop to the value of Y
COMPUTE X=X+1.
END LOOP.
```

- The number of iterations for a case depends on the value of the variable Y for that case. For a case with value 0 for the variable Y, the loop is not executed and X is unchanged. For a case with value 1 for the variable Y, the loop is executed once and X is increased by 1.

**Example**

```
* Factorial routine.

DATA LIST FREE / X.
BEGIN DATA
1 2 3 4 5 6 7
END DATA.

COMPUTE FACTOR=1.
LOOP #I=1 TO X.
COMPUTE FACTOR=FACTOR * #I.
END LOOP.
LIST.
```

- The loop structure computes FACTOR as the factorial value of X.

**Example**

\* Example of nested loops:  
compute every possible combination of values for each variable.

```
INPUT PROGRAM.
-LOOP #I=1 TO 4. /* LOOP TO NUMBER OF VALUES FOR I
- LOOP #J=1 TO 3. /* LOOP TO NUMBER OF VALUES FOR J
- LOOP #K=1 TO 4. /* LOOP TO NUMBER OF VALUES FOR K
- COMPUTE I=#I.
- COMPUTE J=#J.
- COMPUTE K=#K.
- END CASE.
- END LOOP.
- END LOOP.
- END LOOP.
END FILE.
END INPUT PROGRAM.
LIST.
```

- The first loop iterates four times. The first iteration sets the indexing variable *#I* equal to 1 and then passes control to the second loop. *#I* remains 1 until the second loop has completed all of its iterations.
- The second loop is executed 12 times, three times for each value of *#I*. The first iteration sets the indexing variable *#J* equal to 1 and then passes control to the third loop. *#J* remains 1 until the third loop has completed all of its iterations.
- The third loop results in 48 iterations ( $4 \times 3 \times 4$ ). The first iteration sets *#K* equal to 1. The COMPUTE statements set the variables *I*, *J*, and *K* each to 1, and END CASE creates a case. The third loop iterates a second time, setting *#K* equal to 2. Variables *I*, *J*, and *K* are then computed with values 1, 1, 2, respectively, and a second case is created. The third and fourth iterations of the third loop produce cases with *I*, *J*, and *K*, equal to 1, 1, 3 and 1, 1, 4, respectively. After the fourth iteration within the third loop, control passes back to the second loop.
- The second loop is executed again. *#I* remains 1, while *#J* increases to 2, and control returns to the third loop. The third loop completes its iterations, resulting in four more cases with *I* equal to 1, *J* to 2, and *K* increasing from 1 to 4. The second loop is executed a third time, resulting in cases with *I*=1, *J*=3, and *K* increasing from 1 to 4. Once the second loop has completed three iterations, control passes back to the first loop, and the entire cycle is repeated for the next increment of *#I*.
- Once the first loop completes four iterations, control passes out of the looping structures to END FILE. END FILE defines the resulting cases as a data file, the input program terminates, and the LIST command is executed.
- This example does not require a LEAVE command because the iteration variables are scratch variables. If the iteration variables were *I*, *J*, and *K*, LEAVE would be required because the variables would be reinitialized after each END CASE command.

**Example**

\* Modifying the loop iteration variable.

```
INPUT PROGRAM.
PRINT SPACE 2.
LOOP A = 1 TO 3. /*Simple iteration
+ PRINT /'A WITHIN LOOP: ' A(F1).
+ COMPUTE A = 0.
```

## LOOP-END LOOP

```

END LOOP.
PRINT          /'A AFTER LOOP:  ' A(F1).

NUMERIC      #B.
LOOP        B = 1 TO 3. /*Iteration + UNTIL
+ PRINT    /'B WITHIN LOOP:  ' B(F1).
+ COMPUTE  B = 0.
+ COMPUTE  #B = #B+1.
END LOOP    IF #B = 3.
PRINT      /'B AFTER LOOP:  ' B(F1).

NUMERIC      #C.
LOOP        C = 1 TO 3 IF #C NE 3. /*Iteration + WHILE
+ PRINT    /'C WITHIN LOOP:  ' C(F1).
+ COMPUTE  C = 0.
+ COMPUTE  #C = #C+1.
END LOOP.
PRINT      /'C AFTER LOOP:  ' C(F1).

NUMERIC      #D.
LOOP        D = 1 TO 3. /*Iteration + BREAK
+ PRINT    /'D WITHIN LOOP:  ' D(F1).
+ COMPUTE  D = 0.
+ COMPUTE  #D = #D+1.
+ DO IF    #D = 3.
+ BREAK.
+ END IF.
END LOOP.
PRINT      /'D AFTER LOOP:  ' D(F1).

LOOP        E = 3 TO 1. /*Zero-trip iteration
+ PRINT    /'E WITHIN LOOP:  ' E(F1).
+ COMPUTE  E = 0.
END LOOP.
PRINT      /'E AFTER LOOP:  ' E(F1).
END FILE.
END INPUT PROGRAM.
EXECUTE.

```

- If a loop is exited via BREAK or a conditional clause on the END LOOP statement, the iteration variable is not updated.
- If the LOOP statement contains both an indexing clause and a conditional clause, the indexing clause is executed first, and the actual iteration variable will be updated regardless of which clause causes termination of the loop.

The output from this example is shown below.

**Figure 127-1**  
*Modifying the loop iteration value*

```

A WITHIN LOOP: 1
A WITHIN LOOP: 2
A WITHIN LOOP: 3
A AFTER LOOP:  4
B WITHIN LOOP: 1
B WITHIN LOOP: 2
B WITHIN LOOP: 3
B AFTER LOOP:  0
C WITHIN LOOP: 1
C WITHIN LOOP: 2
C WITHIN LOOP: 3
C AFTER LOOP:  4
D WITHIN LOOP: 1
D WITHIN LOOP: 2
D WITHIN LOOP: 3

```

```
D AFTER LOOP: 0
E AFTER LOOP: 3
```

## BY Keyword

By default, the program increases the indexing variable by 1 for each iteration. The keyword BY overrides this increment.

- The *increment value* can be a numeric expression and can therefore be non-integer or negative. Zero causes a warning and results in a zero-trip loop.
- If the initial value is greater than the terminal value and the increment is positive, the loop is never entered. #I=1 TO 0 BY 2 results in a zero-trip loop.
- If the initial value is less than the terminal value and the increment is negative, the loop is never entered. #I=1 TO 2 BY -1 also results in a zero-trip loop.
- Order is unimportant: 2 BY 2 TO 10 is equivalent to 2 TO 10 BY 2.

### Example

```
LOOP #I=2 TO 10 BY 2. /*Loop five times by 2'S
COMPUTE X=X+1.
END LOOP.
```

- The scratch variable *#I* starts at 2 and increases by 2 for each of five iterations until it equals 10 for the last iteration.

### Example

```
LOOP #I=1 TO Y BY Z. /*Loop to Y incrementing by Z
COMPUTE X=X+1.
END LOOP.
```

- The loop is executed once for a case with *Y* equal to 2 and *Z* equal to 2 but twice for a case with *Y* equal to 3 and *Z* equal to 2.

### Example

\* Repeating data using LOOP.

```
INPUT PROGRAM.
DATA LIST      NOTABLE/ ORDER 1-4(N) #BKINFO 6-71(A).
LEAVE ORDER.
LOOP          #I = 1 TO 66 BY 6 IF SUBSTR(#BKINFO,#I,6) <> ' '.
+ REREAD      COLUMN = #I+5.
+ DATA LIST  NOTABLE/ ISBN 1-3(N) QUANTITY 4-5.
+ END CASE.
END LOOP.
END INPUT PROGRAM.
SORT CASES    BY ISBN ORDER.
BEGIN DATA
1045 182 2 155 1 134 1 153 5
1046 155 3 153 5 163 1
1047 161 5 182 2 163 4 186 6
1048 186 2
1049 155 2 163 2 153 2 074 1 161 1
END DATA.
```

---

**LOOP-END LOOP**

```
DO IF          $CASENUM = 1.
+ PRINT EJECT  /'Order' 1  'ISBN' 7  'Quantity' 13.
END IF.
PRINT         /ORDER 2-5 (N) ISBN 8-10 (N) QUANTITY 13-17.
EXECUTE.
```

- This example uses LOOP to simulate a REPEATING DATA command.
- DATA LIST specifies the scratch variable #BKINFO as a string variable (format A) to allow blanks in the data.
- LOOP is executed if the SUBSTR function returns anything other than a blank or null value. SUBSTR returns a six-character substring of #BKINFO, beginning with the character in the position specified by the value of the indexing variable #I. As specified on the indexing clause, #I begins with a value of 1 and is increased by 6 for each iteration of LOOP, up to a maximum #I value of 61 ( $1 + 10 \times 6 = 61$ ). The next iteration would exceed the maximum #I value ( $1 + 11 \times 6 = 67$ ).

## Missing Values

- If the program encounters a case with a missing value for the initial, terminal, or increment value or expression, or if the conditional expression on the LOOP command returns missing, a zero-trip loop results and control is passed to the first command after the END LOOP command.
- If a case has a missing value for the conditional expression on an END LOOP command, the loop is terminated after the first iteration.
- To prevent cases with missing values for any variable used in the loop structure from entering the loop, use the IF clause on the LOOP command (see third example below).

### Example

```
LOOP #I=1 TO Z IF (Y GT 10). /*Loop to X=Z for cases with Y GT 10
COMPUTE X=X+1.
END LOOP.
```

- The value of *X* remains unchanged for cases with a missing value for *Y* or a missing value for *Z* (or if *Z* is less than 1).

### Example

```
MISSING VALUES X(5).
LOOP.
COMPUTE X=X+1.
END LOOP IF (X GE 10). /*Loop until X is at least 10 or missing
```

- Looping is terminated when the value of *X* is 5 because 5 is defined as missing for *X*.

### Example

```
LOOP IF NOT MISSING(Y). /*Loop only when Y isn't missing
COMPUTE X=X+Y.
END LOOP IF (X GE 10). /*Loop until X is at least 10
```

- The variable *X* is unchanged for cases with a missing value for *Y*, since the loop is never entered.

## Creating Data

A loop structure and an `END CASE` command within an input program can be used to create data without any data input. The `END FILE` command must be used outside the loop (but within the input program) to terminate processing.

### Example

```
INPUT PROGRAM.
LOOP #I=1 TO 20.
COMPUTE AMOUNT=RND(UNIFORM(5000))/100.
END CASE.
END LOOP.
END FILE.
END INPUT PROGRAM.

PRINT FORMATS AMOUNT (DOLLAR6.2).
PRINT /AMOUNT.
EXECUTE.
```

- This example creates 20 cases with a single variable, *AMOUNT*. *AMOUNT* is a uniformly distributed number between 0 and 5,000, rounded to an integer and divided by 100 to provide a variable in dollars and cents.
- The `END FILE` command is required to terminate processing once the loop structure is complete.

### Scratch vs. Permanent Index Variables

Permanent variables are reinitialized to system-missing for each case, but scratch variables are initialized to 0 and are not reinitialized for each case; instead they retain their previous values. For loops that don't span cases, this is not an important factor, but in an input program a nested loop can persist across cases. In such instances, the index counter is not affected because it is cached and restored after execution of the loop, but the results of commands within the loop that use the value of the index variables will be different for scratch and permanent index variables.

```
*using scratch index variables.
INPUT PROGRAM.
LOOP #i=1 to 3.
- LOOP #j=#i to 4.
- COMPUTE var1=#i.
- COMPUTE var2=#j.
- END CASE.
- END LOOP.
END LOOP.
END FILE.
END INPUT PROGRAM.
LIST.

*using non-scratch index variables.
INPUT PROGRAM.
LOOP i=1 to 3.
- LOOP j=i to 4.
- COMPUTE var1=i.
- COMPUTE var2=j.
- END CASE.
- END LOOP.
END LOOP.
END FILE.
```

```
END INPUT PROGRAM.
LIST.
```

**Figure 127-2***Case listing from loop using scratch variables*

var1	var2
1.00	1.00
1.00	2.00
1.00	3.00
1.00	4.00
2.00	2.00
2.00	3.00
2.00	4.00
3.00	3.00
3.00	4.00

**Figure 127-3***Case listing from loop using permanent variables*

i	j	var1	var2
1.00	1.00	1.00	1.00
.	2.00	.	2.00
.	3.00	.	3.00
.	4.00	.	4.00
2.00	2.00	2.00	2.00
.	3.00	.	3.00
.	4.00	.	4.00
3.00	3.00	3.00	3.00
.	4.00	.	4.00

- Aside from the fact that the use of permanent variables as index variables results in the creation (or replacement) of those variables in the active dataset, note that many of the value for *var1* (and *i*) are missing in the results from the input program that uses permanent variables as index variables, while none are missing from the one that uses scratch variables.
- In both cases, the inner loop ends with the `END CASE` command, which causes the permanent variable *i* to be reinitialized to system-missing for subsequent iterations of the inner loop, but the scratch variable *#i* retains its previous value.
- When control passes to the outer loop again, the cached index value is used to set and increment *i*; so it has a non-missing value for the first iteration of the inner loop, but once `END CASE` is encountered, it becomes missing again until control passes to the outer loop again.

For more information, see the topic [Scratch Variables](#) on p. 51.



# MANOVA

MANOVA is available in the Advanced Statistics option.

```
MANOVA dependent varlist [BY factor list (min,max)[factor list...]
                        [WITH covariate list]]

[/WSFACTORS=varname (levels) [varname... ]

[/WSDESIGN]*

[/TRANSFORM [(dependent varlist [/dependent varlist])]=
             [ORTHONORM] [{CONTRAST}] {DEVIATION (refcat) } ]
             {BASIS } {DIFFERENCE
                       {HELMERT
                       {SIMPLE (refcat)
                       {REPEATED
                       {POLYNOMIAL [{(1,2,3...)}]}
                       {metric } }
             {SPECIAL (matrix) }

[/MEASURE=newname newname...]

[/RENAME={newname} {newname}...]
        { * } { * }

[/ERROR={WITHIN } ]
        {RESIDUAL
        {WITHIN + RESIDUAL}
        {n }

[/CONTRAST (factorname)={DEVIATION** [(refcat)] } ] +
        {POLYNOMIAL**[{(1,2,3...)}]}
        {metric }
        {SIMPLE [(refcat)]
        {DIFFERENCE
        {HELMERT
        {REPEATED
        {SPECIAL (matrix)

[/PARTITION (factorname) [=({1,1... })]]
        {n1,n2...}

[/METHOD={UNIQUE** } ] [{CONSTANT**}] [{QR** }]]
        {SEQUENTIAL} {NOCONSTANT} {CHOLESKY}

[/{PRINT }= [CELLINFO [{MEANS} [SSCP] [COV] [COR] [ALL]]]]
        {NOPRINT} [HOMOGENEITY [{ALL} [BARTLETT] [COCHRAN] [BOXM]]]]
        [DESIGN [{OVERALL} [ONEWAY] [DECOMP] [BIAS] [SOLUTION]
                 [REDUNDANCY] [COLLINEARITY] [ALL]]]]
        [PARAMETERS [{ESTIM} [ORTHO][COR][NEGSUM][EFSIZE][OPTIMAL][ALL]]]]
        [SIGNIF [(SINGLEDF)
                 [(MULTIV**)] [(EIGEN)] [(DIMENR)]
                 [(UNIV**)] [(HYPOTH)][(STEPDOWN)] [(BRIEF)]
                 [{(AVERF**)}] [(HF)] [(GG)] [(EFSIZE)]
                 {(AVONLY) }
        [ERROR[(STDDEV)][(COR)][(COV)][(SSCP)]]

[/OMEANS =[VARIABLES(varlist)] [TABLES (factor name ) ] ]
        {factor BY factor}
        {CONSTANT }

[/PMEANS =[VARIABLES(varlist)] [TABLES (factor name )]] [PLOT]] ]
        {factor BY factor}
        {CONSTANT }

[/RESIDUALS=[CASEWISE] [PLOT] ]

[/POWER=[T({.05**})] [F({.05**})] [{APPROXIMATE}]]]
```

```

      {a      }      {a      }      {EXACT      }
[/CINTERVAL={INDIVIDUAL}][(.95) ]
      {JOINT      }      {a      }
      UNIVARIATE ({SCHEFFE})]
      {BONFER      }
      MULTIVARIATE ({ROY      })] ]
      {PILLAI      }
      {BONFER      }
      {HOTELLING}
      {WILKS      }
[/PCOMPS [COR] [COV] [ROTATE(rotttype)]
      [NCOMP(n)] [MINEIGEN(eigencut)] [ALL] ]

[/PLOT={BOXPLOTS} [CELLPLOTS] [NORMAL] [ALL] ]

[/DISCRIM [RAW] [STAN] [ESTIM] [COR] [ALL]
      [ROTATE(rotttype)] [ALPHA(.25**)]]
      {a      }

[/MISSING={LISTWISE**} [{EXCLUDE**}] ]
      {INCLUDE      }

[/MATRIX={IN({file})} [OUT({file})] ]
      {[*]      }      {[*]      }

[/ANALYSIS [{UNCONDITIONAL**}]=([dependent varlist
      {CONDITIONAL      } [WITH covariate varlist]
      [/dependent varlist...][])[WITH varlist] ]

[/DESIGN={factor [(n)] } [BY factor[(n)]] [WITHIN factor[(n)]] [WITHIN... ]
      {POOL(varlist)}

      [+ {factor [(n)] }... ]
      {POOL(varlist)}

      [= n] {AGAINST} {WITHIN      }
      {VS      } {RESIDUAL}
      {WR      }
      {n      }

      [{factor [(n)] } ... ]
      {POOL(varlist)}

      [MWITHIN factor(n)]
      [MUPLUS]
      [CONSTANT [=n] ]

```

\* WSDSIGN uses the same specification as DESIGN, with only within-subjects factors.

† DEVIATION is the default for between-subjects factors, while POLYNOMIAL is the default for within-subjects factors.

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### Example 1

```

* Analysis of Variance
MANOVA RESULT BY TREATMNT(1,4) GROUP(1,2).

```

### Example 2

```

* Analysis of Covariance

```

MANOVA RESULT BY TREATMNT(1,4) GROUP(1,2) WITH RAINFALL.

### Example 3

\* Repeated Measures Analysis

```
MANOVA SCORE1 TO SCORE4 BY CLASS(1,2)
  /WSFACTORS=MONTH(4) .
```

### Example 4

\* Parallelism Test with Crossed Factors

```
MANOVA YIELD BY PLOT(1,4) TYPEFERT(1,3) WITH FERT
  /ANALYSIS YIELD
  /DESIGN FERT, PLOT, TYPEFERT, PLOT BY TYPEFERT,
  FERT BY PLOT + FERT BY TYPEFERT
  + FERT BY PLOT BY TYPEFERT.
```

## Overview

MANOVA (multivariate analysis of variance) is a generalized procedure for analysis of variance and covariance. MANOVA is a powerful procedure and can be used for both univariate and multivariate designs. MANOVA allows you to perform the following tasks:

- Specify nesting of effects.
- Specify individual error terms for effects in mixed-model analyses.
- Estimate covariate-by-factor interactions to test the assumption of homogeneity of regressions.
- Obtain parameter estimates for a variety of contrast types, including irregularly spaced polynomial contrasts with multiple factors.
- Test user-specified special contrasts with multiple factors.
- Partition effects in models.
- Pool effects in models.

## MANOVA and General Linear Model (GLM)

MANOVA is available only in syntax. GLM (general linear model), the other generalized procedure for analysis of variance and covariance, is available both in syntax and via the dialog boxes. The major distinction between GLM and MANOVA in terms of statistical design and functionality is that GLM uses a non-full-rank, or overparameterized, indicator variable approach to parameterization of linear models (instead of the full-rank reparameterization approach that is used in MANOVA). GLM uses a generalized inverse approach and uses the aliasing of redundant parameters to zero to allow greater flexibility in handling a variety of data situations, particularly situations involving empty cells. For features that are provided by GLM but unavailable in MANOVA, refer to [General Linear Model \(GLM\) and MANOVA on p. 885](#).

To simplify the presentation, MANOVA reference material is divided into three sections: *univariate* designs with one dependent variable; *multivariate* designs with several interrelated dependent variables; and *repeated measures* designs in which the dependent variables represent the same types of measurements, taken at more than one time.

The full syntax diagram for *MANOVA* is presented here. The sections that follow include partial syntax diagrams that show the subcommands and specifications that are discussed in that section. Individually, those diagrams are incomplete. Subcommands that are listed for univariate designs are available for any analysis, and subcommands that are listed for multivariate designs can be used in any multivariate analysis, including repeated measures.

*MANOVA* was designed and programmed by Philip Burns of Northwestern University.

# MANOVA: Univariate

MANOVA is available in the Advanced Statistics option.

```
MANOVA dependent var [BY factor list (min,max)][factor list...]  
[WITH covariate list]  
[/ERROR={WITHIN  
  {RESIDUAL }  
  {WITHIN + RESIDUAL}  
  {n }  
[/CONTRAST (factor name)={DEVIATION** [(refcat)] }]  
  {POLYNOMIAL [(1,2,3...)]}  
  { {metric } }  
  {SIMPLE [(refcat)] }  
  {DIFFERENCE }  
  {HELMERT }  
  {REPEATED }  
  {SPECIAL (matrix) }  
[/PARTITION (factor name)=[(1,1... )]]  
  {n1,n2... }  
[/METHOD={UNIQUE** }] [{CONSTANT**}] [{QR** }]  
  {SEQUENTIAL} {NOCONSTANT} {CHOLESKY}  
[/PRINT ]= [CELLINFO [(MEANS) [SSCP] [COV] [COR] [ALL]]]  
  {NOPRINT} [HOMOGENEITY [(ALL) [BARTLETT] [COCHRAN]]]  
  [DESIGN [(OVERALL] [ONEWAY] [DECOMP] [BIAS] [SOLUTION]  
  [REDUNDANCY] [COLLINEARITY]]]  
  [PARAMETERS [(ESTIM) [ORTHO] [COR] [NEGSUM] [EFSIZE] [OPTIMAL] [ALL]]]  
  [SIGNIF[(SINGLEDF)]]  
  [ERROR[(STDDEV)] ]  
[/OMEANS =[VARIABLES (varlist)] [TABLES ({factor name } ) ]  
  {factor BY factor}  
  {CONSTANT }  
[/PMEANS =[TABLES ({factor name } ) ] [PLOT]] ]  
  {factor BY factor}  
  {CONSTANT }  
[/RESIDUALS=[CASEWISE] [PLOT] ]  
[/POWER=[T(.05**)] [F(.05**)] [{APPROXIMATE}] ]  
  {a } {a } {EXACT }  
[/CINTERVAL=[INDIVIDUAL] [(0.95) ]] [UNIVARIATE ({SCHEFFE}) ]  
  {JOINT } {a } {BONFER }  
[/PLOT=[BOXPLOTS] [CELLPLOTS] [NORMAL] [ALL] ]  
[/MISSING=[LISTWISE**] [{EXCLUDE**}] ]  
  {INCLUDE }  
[/MATRIX=[IN({file})] [OUT({file})]  
  {[*] } {[*] }  
[/ANALYSIS=dependent var [WITH covariate list]]  
[/DESIGN={factor [(n) ]} [BY factor[(n)]] [WITHIN factor[(n)]] [WITHIN...]  
  {POOL(varlist)}  
  [+ {factor [(n) ]} ... ]  
  {POOL(varlist)}  
  [[= n] {AGAINST} {WITHIN }  
  {VS } {RESIDUAL}  
  {WR }  
  {n }  
  [{factor [(n) ]} ... ]  
  {POOL(varlist)}  
  [MUPLUS]  
  [MWITHIN factor(n)]  
  [CONSTANT [=n] ]
```

\*\* Default if the subcommand or keyword is omitted.

## Example

```
MANOVA YIELD BY SEED(1,4) FERT(1,3)  
/DESIGN.
```

## Overview

This section describes the use of `MANOVA` for univariate analyses. However, the subcommands that are described here can be used in any type of analysis with `MANOVA`. For additional subcommands that are used in multivariate analysis, see *MANOVA: Multivariate*. For additional subcommands that are used in repeated measures analysis, see *MANOVA: Repeated Measures*. For basic specification, syntax rules, and limitations of the `MANOVA` procedures, see *MANOVA*.

### Options

**Design Specification.** You can use the `DESIGN` subcommand to specify which terms to include in the design. This ability allows you to estimate a model other than the default full factorial model, incorporate factor-by-covariate interactions, indicate nesting of effects, and indicate specific error terms for each effect in mixed models. You can specify a different continuous variable as a dependent variable or work with a subset of the continuous variables with the `ANALYSIS` subcommand.

**Contrast Types.** You can specify contrasts other than the default deviation contrasts on the `CONTRAST` subcommand. You can also subdivide the degrees of freedom associated with a factor (using the `PARTITION` subcommand) and test the significance of a specific contrast or group of contrasts.

**Optional Output.** You can choose from a variety of optional output on the `PRINT` subcommand or suppress output using the `NOPRINT` subcommand. Output that is appropriate to univariate designs includes cell means, design or other matrices, parameter estimates, and tests for homogeneity of variance across cells. Using the `OMEANS`, `PMEANS`, `RESIDUAL`, and `PLOT` subcommands, you can also request tables of observed and/or predicted means, casewise values and residuals for your model, and various plots that are useful in checking assumptions. In addition, you can use the `POWER` subcommand to request observed power values (based on fixed-effect assumptions), and you can use the `CINTERVAL` subcommand to request simultaneous confidence intervals for each parameter estimate and regression coefficient.

**Matrix Materials.** You can write matrices of intermediate results to a matrix data file, and you can read such matrices in performing further analyses by using the `MATRIX` subcommand.

### Basic Specification

- The basic specification is a variable list that identifies the dependent variable, the factors (if any), and the covariates (if any).
- By default, `MANOVA` uses a full factorial model, which includes all main effects and all possible interactions among factors. Estimation is performed by using the cell-means model and `UNIQUE` (regression-type) sums of squares, adjusting each effect for all other effects in the model. Parameters are estimated by using `DEVIATION` contrasts to determine whether their categories differ significantly from the mean.

### Subcommand Order

- The variable list must be specified first.

- Subcommands that are applicable to a specific design must be specified before that `DESIGN` subcommand. Otherwise, subcommands can be used in any order.

### **Syntax Rules**

- For many analyses, the `MANOVA` variable list and the `DESIGN` subcommand are the only specifications that are needed. If a full factorial design is desired, `DESIGN` can be omitted.
- All other subcommands apply only to designs that follow. If you do not enter a `DESIGN` subcommand, or if the last subcommand is not `DESIGN`, `MANOVA` uses a full factorial model.
- Unless replaced, `MANOVA` subcommands (other than `DESIGN`) remain in effect for all subsequent models.
- `MISSING` can be specified only once.
- The following words are reserved as keywords or internal commands in the `MANOVA` procedure: `AGAINST`, `CONSPLUS`, `CONSTANT`, `CONTIN`, `MUPLUS`, `MWITHIN`, `POOL`, `R`, `RESIDUAL`, `RW`, `VERSUS`, `VS`, `W`, `WITHIN`, and `WR`. Variable names that duplicate these words should be changed before you invoke `MANOVA`.
- If you enter one of the multivariate specifications in a univariate analysis, `MANOVA` will ignore it.

### **Limitations**

- A maximum of 20 factors is in place.
- A maximum of 200 dependent variables is in place.
- Memory requirements depend primarily on the number of cells in the design. For the default full factorial model, the number of cells equals the product of the number of levels or categories in each factor.

## **Example**

```
MANOVA YIELD BY SEED(1,4) FERT(1,3) WITH RAINFALL
/PRINT=CELLINFO(MEANS) PARAMETERS(ESTIM)
/DESIGN.
```

- `YIELD` is the dependent variable; `SEED` (with values 1, 2, 3, and 4) and `FERT` (with values 1, 2, and 3) are factors; `RAINFALL` is a covariate.
- The `PRINT` subcommand requests the means of the dependent variable for each cell and the default deviation parameter estimates.
- The `DESIGN` subcommand requests the default design, a full factorial model. This subcommand could have been omitted or could have been specified in full as:  
`/DESIGN = SEED, FERT, SEED BY FERT.`

## **MANOVA Variable List**

The variable list specifies all variables that will be used in any subsequent analyses.

- The dependent variable must be the first specification on `MANOVA`.

- By default, MANOVA treats a list of dependent variables as jointly dependent, implying a multivariate design. However, you can use the ANALYSIS subcommand to change the role of a variable or its inclusion status in the analysis.
- The names of the factors follow the dependent variable. Use the keyword BY to separate the factors from the dependent variable.
- Factors must have adjacent integer values, and you must supply the minimum and maximum values in parentheses after the factor name(s).
- If several factors have the same value range, you can specify a list of factors followed by a single value range in parentheses.
- Certain one-cell designs, such as univariate and multivariate regression analysis, canonical correlation, and one-sample Hotelling's  $T^2$ , do not require a factor specification. To perform these analyses, omit the keyword BY and the factor list.
- Enter the covariates, if any, following the factors and their ranges. Use the keyword WITH to separate covariates from factors (if any) and the dependent variable.

**Example**

```
MANOVA DEPENDNT BY FACTOR1 (1,3) FACTOR2, FACTOR3 (1,2).
```

- In this example, three factors are specified.
- *FACTOR1* has values 1, 2, and 3, while *FACTOR2* and *FACTOR3* have values 1 and 2.
- A default full factorial model is used for the analysis.

**Example**

```
MANOVA Y BY A(1,3) WITH X
/DESIGN.
```

- In this example, the *A* effect is tested after adjusting for the effect of the covariate *X*. It is a test of equality of adjusted *A* means.
- The test of the covariate *X* is adjusted for *A*. The test is a test of the pooled within-groups regression of *Y* on *X*.

**ERROR Subcommand**

ERROR allows you to specify or change the error term that is used to test all effects for which you do not explicitly specify an error term on the DESIGN subcommand. ERROR affects all terms in all subsequent designs, except terms for which you explicitly provide an error term.

**WITHIN**

*Terms in the model are tested against the within-cell sum of squares.* This specification can be abbreviated to *w*. This setting is the default unless there is no variance within cells or a continuous variable is named on the DESIGN subcommand.

**RESIDUAL**

*Terms in the model are tested against the residual sum of squares.* This specification can be abbreviated to *r*. This specification includes all terms not named on the DESIGN subcommand.



<b>WITHIN+RESIDUAL</b>	<i>Terms are tested against the pooled within-cells and residual sum of squares. This specification can be abbreviated to WR or RW. This setting is the default for designs in which a continuous variable appears on the DESIGN subcommand.</i>
<b>error number</b>	<i>Terms are tested against a numbered error term. The error term must be defined on each DESIGN subcommand. For a discussion of error terms, see <a href="#">DESIGN Keyword on p. 1081</a>.</i>

- If you specify `ERROR=WITHIN+RESIDUAL` and one of the components does not exist, MANOVA uses the other component alone.
- If you specify your own error term by number and a design does not have an error term with the specified number, MANOVA does not carry out significance tests. MANOVA will, however, display hypothesis sums of squares and, if requested, parameter estimates.

### Example

```
MANOVA DEP BY A(1,2) B(1,4)
  /ERROR = 1
  /DESIGN = A, B, A BY B = 1 VS WITHIN
  /DESIGN = A, B.
```

- `ERROR` defines error term 1 as the default error term.
- In the first design, *A* by *B* is defined as error term 1 and is therefore used to test the *A* and *B* effects. The *A* by *B* effect itself is explicitly tested against the within-cells error.
- In the second design, no term is defined as error term 1, so no significance tests are carried out. Hypothesis sums of squares are displayed for *A* and *B*.

## CONTRAST Subcommand

`CONTRAST` specifies the type of contrast that is desired among the levels of a factor. For a factor with  $k$  levels or values, the contrast type determines the meaning of its  $k-1$  degrees of freedom. If the subcommand is omitted or is specified with no keyword, the default is `DEVIATION` for between-subjects factors.

- Specify the factor name in parentheses following the subcommand `CONTRAST`.
- You can specify only one factor per `CONTRAST` subcommand, but you can enter multiple `CONTRAST` subcommands.
- After closing the parentheses, enter an equals sign followed by one of the contrast keywords.
- To obtain  $F$  tests for individual degrees of freedom for the specified contrast, enter the factor name followed by a number in parentheses on the `DESIGN` subcommand. The number refers to a partition of the factor's degrees of freedom. If you do not use the `PARTITION` subcommand, each degree of freedom is a distinct partition.

The following contrast types are available:

- DEVIATION**      *Deviations from the grand mean.* This setting is the default for between-subjects factors. Each level of the factor (except one level) is compared to the grand mean. One category (by default, the last category) must be omitted so that the effects will be independent of one another. To omit a category other than the last category, specify the number of the omitted category (which is not necessarily the same as its value) in parentheses after the keyword `DEVIATION`. An example is as follows:  
`MANOVA A BY B(2,4) /CONTRAST(B)=DEVIATION(1).`  
 The specified contrast omits the first category, in which *B* has the value 2. Deviation contrasts are not orthogonal.
- POLYNOMIAL**      *Polynomial contrasts.* This setting is the default for within-subjects factors. The first degree of freedom contains the linear effect across the levels of the factor, the second degree of freedom contains the quadratic effect, and so on. In a balanced design, polynomial contrasts are orthogonal. By default, the levels are assumed to be equally spaced; you can specify unequal spacing by entering a metric—consisting of one integer for each level of the factor—in parentheses after the keyword `POLYNOMIAL`. An example is as follows:  
`MANOVA RESPONSE BY STIMULUS (4,6) /CONTRAST(STIMULUS) = POLYNOMIAL(1,2,4).`  
 The specified contrast indicates that the three levels of *STIMULUS* are actually in the proportion 1:2:4. The default metric is always (1,2,...,k), where *k* levels are involved. Only the relative differences between the terms of the metric matter. (1,2,4) is the same metric as (2,3,5) or (20,30,50) because, in each instance, the difference between the second and third numbers is twice the difference between the first and second numbers.
- DIFFERENCE**      *Difference or reverse Helmert contrasts.* Each level of the factor (except the first level) is compared to the mean of the previous levels. In a balanced design, difference contrasts are orthogonal.
- HELMERT**      *Helmert contrasts.* Each level of the factor (except the last level) is compared to the mean of subsequent levels. In a balanced design, Helmert contrasts are orthogonal.
- SIMPLE**      *Contrast where each level of the factor (except the last level) is compared to the last level.* To use a category (other than the last category) as the omitted reference category, specify its number (which is not necessarily the same as its value) in parentheses following the keyword `SIMPLE`. An example is as follows:  
`MANOVA A BY B(2,4) /CONTRAST(B)=SIMPLE(1).`  
 The specified contrast compares the other levels to the first level of *B*, in which *B* has the value 2. Simple contrasts are not orthogonal.
- REPEATED**      *Comparison of adjacent levels.* Each level of the factor (except the last level) is compared to the next level. Repeated contrasts are not orthogonal.
- SPECIAL**      *A user-defined contrast.* After this keyword, enter a square matrix in parentheses with as many rows and columns as there are levels in the factor. The first row represents the mean effect of the factor and is generally a vector of 1's. The row represents a set of weights indicating how to collapse over the categories of this factor in estimating parameters for other factors. The other rows of the contrast matrix contain the special contrasts indicating the desired comparisons between levels of the factor. If the special contrasts are linear combinations of each other, MANOVA reports the linear dependency and stops processing.

**Orthogonal contrasts** are particularly useful. In a balanced design, contrasts are orthogonal if the sum of the coefficients in each contrast row is 0 and if, for any pair of contrast rows, the products of corresponding coefficients sum to 0. `DIFFERENCE`, `HELMERT`, and `POLYNOMIAL` contrasts always meet these criteria in balanced designs.

### Example

```
MANOVA DEP BY FAC(1,5)
  /CONTRAST(FAC)=DIFFERENCE
  /DESIGN=FAC(1) FAC(2) FAC(3) FAC(4) .
```

- The factor *FAC* has five categories and therefore four degrees of freedom.
- `CONTRAST` requests `DIFFERENCE` contrasts, which compare each level (except the first level) with the mean of the previous levels.
- Each of the four degrees of freedom is tested individually on the `DESIGN` subcommand.

## ***PARTITION Subcommand***

`PARTITION` subdivides the degrees of freedom that are associated with a factor. This process permits you to test the significance of the effect of a specific contrast or group of contrasts of the factor instead of the overall effect of all contrasts of the factor. The default is a single degree of freedom for each partition.

- Specify the factor name in parentheses following the `PARTITION` subcommand.
- Specify an integer list in parentheses after the optional equals sign to indicate the degrees of freedom for each partition.
- Each value in the partition list must be a positive integer, and the sum of the values cannot exceed the degrees of freedom for the factor.
- The degrees of freedom that are available for a factor are one less than the number of levels of the factor.
- The meaning of each degree of freedom depends on the contrast type for the factor. For example, with deviation contrasts (the default for between-subjects factors), each degree of freedom represents the deviation of the dependent variable in one level of the factor from its grand mean over all levels. With polynomial contrasts, the degrees of freedom represent the linear effect, the quadratic effect, and so on.
- If your list does not account for all the degrees of freedom, `MANOVA` adds one final partition containing the remaining degrees of freedom.
- You can use a repetition factor of the form  $n^*$  to specify a series of partitions with the same number of degrees of freedom.
- To specify a model that tests only the effect of a specific partition of a factor in your design, include the number of the partition in parentheses on the `DESIGN` subcommand (see the example below).
- If you want the default single degree-of-freedom partition, you can omit the `PARTITION` subcommand and simply enter the appropriate term on the `DESIGN` subcommand.

**Example**

```
MANOVA OUTCOME BY TREATMNT(1,12)
  /PARTITION(TREATMNT) = (3*2,4)
  /DESIGN TREATMNT(2) .
```

- The factor *TREATMNT* has 12 categories (hence, 11 degrees of freedom).
- *PARTITION* divides the effect of *TREATMNT* into four partitions, containing, respectively, 2, 2, 2, and 4 degrees of freedom. A fifth partition is formed to contain the remaining 1 degree of freedom.
- *DESIGN* specifies a model in which only the second partition of *TREATMNT* is tested. This partition contains the third and fourth degrees of freedom.
- Because the default contrast type for between-subjects factors is *DEVIATION*, this second partition represents the deviation of the third and fourth levels of *TREATMNT* from the grand mean.

**METHOD Subcommand**

*METHOD* controls the computational aspects of the MANOVA analysis. You can specify one of two different methods for partitioning the sums of squares. The default is *UNIQUE*.

<b>UNIQUE</b>	<i>Regression approach.</i> Each term is corrected for every other term in the model. With this approach, sums of squares for various components of the model do not add up to the total sum of squares unless the design is balanced. This is the default if the <i>METHOD</i> subcommand is omitted or if neither of the two keywords is specified.
<b>SEQUENTIAL</b>	<i>Hierarchical decomposition of the sums of squares.</i> Each term is adjusted only for the terms that precede it on the <i>DESIGN</i> subcommand. This decomposition is an orthogonal decomposition, and the sums of squares in the model add up to the total sum of squares.

You can control how parameters are to be estimated by specifying one of the following two keywords that are available on *MANOVA*. The default is *QR*.

<b>QR</b>	<i>Use modified Givens rotations.</i> <i>QR</i> bypasses the normal equations and the inaccuracies that can result from creating the cross-products matrix, and it generally results in extremely accurate parameter estimates. This setting is the default if the <i>METHOD</i> subcommand is omitted or if neither of the two keywords is specified.
<b>CHOLESKY</b>	<i>Use Cholesky decomposition of the cross-products matrix.</i> This method is useful for large data sets with covariates entered on the <i>DESIGN</i> subcommand.

You can also control whether a constant term is included in all models. Two keywords are available on *METHOD*. The default is *CONSTANT*.

<b>CONSTANT</b>	<i>All models include a constant (grand mean) term, even if none is explicitly specified on the DESIGN subcommand.</i> This setting is the default if neither of the two keywords is specified.
<b>NOCONSTANT</b>	<i>Exclude constant terms from models that do not include the keyword CONSTANT on the DESIGN subcommand.</i>

**Example**

```
MANOVA DEP BY A B C (1,4)
  /METHOD=NOCONSTANT
  /DESIGN=A, B, C
  /METHOD=CONSTANT SEQUENTIAL
  /DESIGN.
```

- For the first design, a main-effects model, the METHOD subcommand requests the model to be fitted with no constant.
- The second design requests a full factorial model to be fitted with a constant and with a sequential decomposition of sums of squares.

**PRINT and NOPRINT Subcommands**

PRINT and NOPRINT control the display of optional output.

- Specifications on PRINT remain in effect for all subsequent designs.
- Some PRINT output, such as CELLINFO, applies to the entire MANOVA procedure and is displayed only once.
- You can turn off optional output that you have requested on PRINT by entering a NOPRINT subcommand with the specifications that were originally used on the PRINT subcommand.
- Additional output can be obtained on the PCOMPS, DISCRIM, OMEANS, PMEANS, PLOT, and RESIDUALS subcommands.
- Some optional output greatly increases the processing time. Request only the output that you want to see.

The following specifications are appropriate for univariate MANOVA analyses. For information about PRINT specifications that are appropriate for multivariate models, see [PRINT and NOPRINT Subcommands on p. 1103](#). For information about PRINT specifications that are appropriate for repeated measures models, see [PRINT Subcommand on p. 1118](#).

<b>CELLINFO</b>	<i>Basic information about each cell in the design.</i>
<b>PARAMETERS</b>	<i>Parameter estimates.</i>
<b>HOMOGENEITY</b>	<i>Tests of homogeneity of variance.</i>
<b>DESIGN</b>	<i>Design information.</i>
<b>ERROR</b>	<i>Error standard deviations.</i>

**CELLINFO Keyword**

You can request any of the following cell information by specifying the appropriate keyword(s) in parentheses after CELLINFO. The default is MEANS.

<b>MEANS</b>	<i>Cell means, standard deviations, and counts for the dependent variable and covariates. Confidence intervals for the cell means are displayed if you have set a wide width. This setting is the default when CELLINFO is requested with no further specification.</i>
<b>SSCP</b>	<i>Within-cell sum-of-squares and cross-products matrices for the dependent variable and covariates.</i>

<b>COV</b>	<i>Within-cell variance-covariance matrices for the dependent variable and covariates.</i>
<b>COR</b>	<i>Within-cell correlation matrices, with standard deviations on the diagonal, for the dependent variable and covariates.</i>
<b>ALL</b>	<i>MEANS, SSCP, COV, and COR.</i>

- Output from `CELLINFO` is displayed once before the analysis of any particular design. Specify `CELLINFO` only once.
- When you specify `SSCP`, `COV`, or `COR`, the cells are numbered for identification, beginning with cell 1.
- The levels vary most rapidly for the factor named last on the `MANOVA` variables specification.
- Empty cells are neither displayed nor numbered.
- At the beginning of `MANOVA` output, a table is displayed, showing the levels of each factor corresponding to each cell number.

### **Example**

```
MANOVA DEP BY A(1,4) B(1,2) WITH COV
/PRINT=CELLINFO(MEANS COV)
/DESIGN.
```

- For each combination of levels of *A* and *B*, `MANOVA` displays separately the means and standard deviations of *DEP* and *COV*. Beginning with cell 1, `MANOVA` will then display the variance-covariance matrix of *DEP* and *COV* within each non-empty cell.
- A table of cell numbers will be displayed to show the factor levels corresponding to each cell.
- The keyword `COV`, as a parameter of `CELLINFO`, is not confused with the variable *COV*.

## **PARAMETERS Keyword**

The keyword `PARAMETERS` displays information about the estimated size of the effects in the model. You can specify any of the following keywords in parentheses on `PARAMETERS`. The default is `ESTIM`.

<b>ESTIM</b>	<i>The estimated parameters themselves, along with their standard errors, t tests, and confidence intervals. Only nonredundant parameters are displayed. This setting is the default if <code>PARAMETERS</code> is requested without further specification.</i>
<b>NEGSUM</b>	<i>The negative of the sum of parameters for each effect. For <code>DEVIATION</code> main effects, this value equals the parameter for the omitted (redundant) contrast. <code>NEGSUM</code> is displayed, along with the parameter estimates.</i>
<b>ORTHO</b>	<i>The orthogonal estimates of parameters that are used to produce the sums of squares.</i>
<b>COR</b>	<i>Covariance factors and correlations among the parameter estimates.</i>
<b>EFSIZE</b>	<i>The effect size values.</i>
<b>OPTIMAL</b>	<i>Optimal Scheffé contrast coefficients.</i>
<b>ALL</b>	<i>ESTIM, NEGSUM, ORTHO, COR, EFSIZE, and OPTIMAL.</i>

## ***SIGNIF Keyword***

`SIGNIF` requests special significance tests, most of which apply to multivariate designs (see [SIGNIF Keyword on p. 1103](#)). The following specification is useful in univariate applications of MANOVA:

**SINGLEDF**            *Significance tests for each single degree of freedom making up each effect for analysis-of-variance tables.*

- When non-orthogonal contrasts are requested or when the design is unbalanced, the `SINGLEDF` effects will differ from single degree-of-freedom partitions. `SINGLEDEF` effects are orthogonal within an effect; single degree-of-freedom partitions are not orthogonal within an effect.

### ***Example***

```
MANOVA DEP BY FAC(1,5)
/CONTRAST(FAC)=POLY
/PRINT=SIGNIF(SINGLEDF)
/DESIGN.
```

- `POLYNOMIAL` contrasts are applied to `FAC`, testing the linear, quadratic, cubic, and quartic components of its five levels. `POLYNOMIAL` contrasts are orthogonal in balanced designs.
- The `SINGLEDF` specification on `SIGNIF` requests significance tests for each of these four components.

## ***HOMOGENEITY Keyword***

`HOMOGENEITY` requests tests for the homogeneity of variance of the dependent variable across the cells of the design. You can specify one or more of the following specifications in parentheses. If `HOMOGENEITY` is requested without further specification, the default is `ALL`.

**BARTLETT**            *Bartlett-Box F test.*  
**COCHRAN**            *Cochran's C.*  
**ALL**                    *Both BARTLETT and COCHRAN. This setting is the default.*

## ***DESIGN Keyword***

You can enter one or more of the following specifications in parentheses following the keyword `DESIGN`. If `DESIGN` is requested without further specification, the default is `OVERALL`.

The `DECOMP` and `BIAS` matrices can provide valuable information about the confounding of the effects and the estimability of the chosen contrasts. If two effects are confounded, the entry corresponding to them in the `BIAS` matrix will be nonzero; if the effects are orthogonal, the entry

will be zero. This feature is particularly useful in designs with unpatterned empty cells. For further discussion of the matrices, see Bock (1985).

<b>OVERALL</b>	<i>The overall reduced-model design matrix (not the contrast matrix). This setting is the default.</i>
<b>ONEWAY</b>	<i>The one-way basis matrix (not the contrast matrix) for each factor.</i>
<b>DECOMP</b>	<i>The upper triangular QR/CHOLESKY decomposition of the design.</i>
<b>BIAS</b>	<i>Contamination coefficients displaying the bias that is present in the design.</i>
<b>SOLUTION</b>	<i>Coefficients of the linear combinations of the cell means that are used in significance testing.</i>
<b>REDUNDANCY</b>	<i>Exact linear combinations of parameters that form a redundancy. This keyword displays a table only if QR (the default) is the estimation method.</i>
<b>COLLINEARITY</b>	<i>Collinearity diagnostics for design matrices. These diagnostics include the singular values of the normalized design matrix (which are the same as those values of the normalized decomposition matrix), condition indexes corresponding to each singular value, and the proportion of variance of the corresponding parameter accounted for by each principal component. For greatest accuracy, use the QR method of estimation whenever you request collinearity diagnostics.</i>
<b>ALL</b>	<i>All available options.</i>

## **ERROR Keyword**

Generally, the keyword `ERROR` on `PRINT` produces error matrices. In univariate analyses, the only valid specification for `ERROR` is `STDDEV`, which is the default if `ERROR` is specified by itself.

<b>STDDEV</b>	<i>The error standard deviation. Normally, this deviation is the within-cells standard deviation of the dependent variable. If you specify multiple error terms on <code>DESIGN</code>, this specification will display the standard deviation for each term.</i>
---------------	---

## **OMEANS Subcommand**

`OMEANS` (observed means) displays tables of the means of continuous variables for levels or combinations of levels of the factors.

- Use the keywords `VARIABLES` and `TABLES` to indicate which observed means you want to display.
- With no specifications, the `OMEANS` subcommand is equivalent to requesting `CELLINFO (MEANS)` on `PRINT`.
- `OMEANS` displays confidence intervals for the cell means if you have set the width to 132.



- Output from OMEANS is displayed once before the analysis of any particular design. This subcommand should be specified only once.

**VARIABLES**      *Continuous variables for which you want means.* Specify the variables in parentheses after the keyword VARIABLES. You can request means for the dependent variable or any covariates. If you omit the VARIABLES keyword, observed means are displayed for the dependent variable and all covariates. If you enter the keyword VARIABLES, you must also enter the keyword TABLES (discussed below).

**TABLES**          *Factors for which you want the observed means displayed.* In parentheses, list the factors, or combinations of factors, separated with BY. Observed means are displayed for each level, or combination of levels, of the factors that are named (see the example below). Both weighted means and unweighted means (where all cells are weighted equally, regardless of the number of cases that they contain) are displayed. If you enter the keyword CONSTANT, the grand mean is displayed.

### Example

```
MANOVA DEP BY A(1,3) B(1,2)
  /OMEANS=TABLES(A,B)
  /DESIGN.
```

- Because there is no VARIABLES specification on the OMEANS subcommand, observed means are displayed for all continuous variables. DEP is the only dependent variable here, and there are no covariates.
- The TABLES specification on the OMEANS subcommand requests tables of observed means for each of the three categories of A (collapsing over B) and for both categories of B (collapsing over A).
- MANOVA displays weighted means, in which all cases count equally, and displays unweighted means, in which all cells count equally.

## PMEANS Subcommand

PMEANS (predicted means) displays a table of the predicted cell means of the dependent variable, adjusted for the effect of covariates in the cell and unadjusted for covariates. For comparison, PMEANS also displays the observed cell means.

- Output from PMEANS can be computationally expensive.
- PMEANS without any additional specifications displays a table showing, for each cell, the observed mean of the dependent variable, the predicted mean adjusted for the effect of covariates in that cell (*ADJ. MEAN*), the predicted mean unadjusted for covariates (*EST. MEAN*), and the raw and standardized residuals from the estimated means.
- Cells are numbered in output from PMEANS so that the levels vary most rapidly on the factor that is named last in the MANOVA variables specification. A table showing the levels of each factor corresponding to each cell number is displayed at the beginning of the MANOVA output.
- Predicted means are suppressed for any design in which the MUPLUS keyword appears.
- Covariates are not predicted.
- In designs with covariates and multiple error terms, use the ERROR subcommand to designate which error term's regression coefficients are to be used in calculating the standardized residuals.

For univariate analysis, the following keywords are available on the `PMEANS` subcommand:

<b>TABLES</b>	<i>Additional tables showing adjusted predicted means for specified factors or combinations of factors.</i> Enter the names of factors or combinations of factors in parentheses after this keyword. For each factor or combination, <code>MANOVA</code> displays the predicted means (adjusted for covariates) collapsed over all other factors.
<b>PLOT</b>	<i>A plot of the predicted means for each cell.</i>

### Example

```
MANOVA DEP BY A(1,4) B(1,3)
/PMEANS TABLES(A, B, A BY B)
/DESIGN = A, B.
```

- `PMEANS` displays the default table of observed and predicted means for `DEP` and raw and standardized residuals in each of the 12 cells in the model.
- The `TABLES` specification on `PMEANS` displays tables of predicted means for `A` (collapsing over `B`), for `B` (collapsing over `A`), and for all combinations of `A` and `B`.
- Because `A` and `B` are the only factors in the model, the means for `A` by `B` in the `TABLES` specification come from every cell in the model. The means are identical to the adjusted predicted means in the default `PMEANS` table, which always includes all non-empty cells.
- Predicted means for `A` by `B` can be requested in the `TABLES` specification, even though the `A` by `B` effect is not in the design.

## RESIDUALS Subcommand

Use `RESIDUALS` to display and plot casewise values and residuals for your models.

- Use the `ERROR` subcommand to specify an error term other than the default to be used to standardize the residuals.
- If a designated error term does not exist for a given design, no predicted values or residuals are calculated.
- If you specify `RESIDUALS` without any keyword, `CASEWISE` output is displayed.

The following keywords are available:

<b>CASEWISE</b>	<i>A case-by-case listing of the observed, predicted, residual, and standardized residual values for each dependent variable.</i>
<b>PLOT</b>	<i>A plot of observed values, predicted values, and case numbers versus the standardized residuals, plus normal and detrended normal probability plots for the standardized residuals (five plots in all).</i>

## POWER Subcommand

`POWER` requests observed power values based on fixed-effect assumptions for all univariate and multivariate  $F$  tests and  $t$  tests. Both approximate and exact power values can be computed, although exact multivariate power is displayed only when there is one hypothesis degree of

freedom. If `POWER` is specified by itself, with no keywords, `MANOVA` calculates the approximate observed power values of all  $F$  tests at 0.05 significance level.

The following keywords are available on the `POWER` subcommand:

<b>APPROXIMATE</b>	<i>Approximate power values.</i> This setting is the default if <code>POWER</code> is specified without any keyword. Approximate power values for univariate tests are derived from an Edgeworth-type normal approximation to the noncentral beta distribution. Approximate values are normally accurate to three decimal places and are much cheaper to compute than exact values.
<b>EXACT</b>	<i>Exact power values.</i> Exact power values for univariate tests are computed from the noncentral incomplete beta distribution.
<b>F(a)</b>	<i>Alpha level at which the power is to be calculated for F tests.</i> The default is 0.05. To change the default, specify a decimal number between 0 and 1 in parentheses after $F$ . The numbers 0 and 1 themselves are not allowed. $F$ test at 0.05 significance level is the default when <code>POWER</code> is omitted or specified without any keyword.
<b>T(a)</b>	<i>Alpha level at which the power is to be calculated for t tests.</i> The default is 0.05. To change the default, specify a decimal number between 0 and 1 in parentheses after $t$ . The numbers 0 and 1 themselves are not allowed.

- For univariate  $F$  tests and  $t$  tests, `MANOVA` computes a measure of the effect size based on partial  $\eta^2$ :

$$\text{partial } \eta^2 = (ssh)/(ssh + sse)$$

where  $ssh$  is the hypothesis sum of squares and  $sse$  is the error sum of squares. The measure is an overestimate of the actual effect size. However, the measure is consistent and is applicable to all  $F$  tests and  $t$  tests. For a discussion of effect size measures, see (Cohen, 1977) or (Hays, 1981).

## ***CINTERVAL Subcommand***

`CINTERVAL` requests simultaneous confidence intervals for each parameter estimate and regression coefficient. `MANOVA` provides either individual or joint confidence intervals at any desired confidence level. You can compute joint confidence intervals that are using either Scheffé or Bonferroni intervals. Scheffé intervals are based on all possible contrasts, while Bonferroni intervals are based on the number of contrasts that are actually made. For a large number of contrasts, Bonferroni intervals will be larger than Scheffé intervals. Timm (Timm, 1975) provides a good discussion of which intervals are best for certain situations. Both Scheffé and Bonferroni intervals are computed separately for each term in the design. You can request only one type of confidence interval per design.

The following keywords are available on the `CINTERVAL` subcommand. If the subcommand is specified without any keyword, `CINTERVAL` automatically displays individual univariate confidence intervals at the 0.95 level.

<b>INDIVIDUAL(a)</b>	<i>Individual confidence intervals.</i> Specify the desired confidence level in parentheses following the keyword. The desired confidence level can be any decimal number between 0 and 1. When individual intervals are requested, <code>BONFER</code> and <code>SCHEFFE</code> have no effect.
<b>JOINT(a)</b>	<i>Joint confidence intervals.</i> Specify the desired confidence level in parentheses after the keyword. The default is 0.95. The desired confidence level can be any decimal number between 0 and 1.
<b>UNIVARIATE(type)</b>	<i>Univariate confidence interval.</i> Specify either <code>SCHEFFE</code> (for Scheffé intervals) or <code>BONFER</code> (for Bonferroni intervals) in parentheses after the keyword. The default specification is <code>SCHEFFE</code> .

## ***PLOT Subcommand***

MANOVA can display a variety of plots that are useful in checking the assumptions that are needed in the analysis. Plots are produced only once in the MANOVA procedure, regardless of how many `DESIGN` subcommands you enter. Use the following keywords on the `PLOT` subcommand to request plots. If the `PLOT` subcommand is specified by itself, the default is `BOXPLOTS`.

<b>BOXPLOTS</b>	<i>Boxplots.</i> Plots are displayed for each continuous variable (dependent or covariate) that is named on the <code>MANOVA</code> variable list. Boxplots provide a simple graphical means of comparing the cells in terms of mean location and spread. The data must be stored in memory for these plots; if there is not enough memory, boxplots are not produced, and a warning message is issued. This setting is the default if the <code>PLOT</code> subcommand is specified without a keyword.
<b>CELLPLOTS</b>	<i>Cell statistics, including a plot of cell means versus cell variances, a plot of cell means versus cell standard deviations, and a histogram of cell means.</i> Plots are produced for each continuous variable (dependent or covariate) that is named on the <code>MANOVA</code> variable list. The first two plots aid in detecting heteroscedasticity (nonhomogeneous variances) and aid in determining an appropriate data transformation (if a transformation is needed). The third plot gives distributional information for the cell means.
<b>NORMAL</b>	<i>Normal and detrended normal plots.</i> Plots are produced for each continuous variable (dependent or covariate) that is named on the <code>MANOVA</code> variable list. MANOVA ranks the scores and then plots the ranks against the expected normal deviate, or detrended expected normal deviate, for that rank. These plots aid in detecting non-normality and outlying observations. All data must be held in memory to compute ranks. If not enough memory is available, MANOVA displays a warning and skips the plots.

- `ZCORR`, an additional plot that is available on the `PLOT` subcommand, is described in [MANOVA: Multivariate](#).
- You can request other plots on `PMEANS` and `RESIDUALS` (see these respective subcommands).

## MISSING Subcommand

By default, cases with missing values for any of the variables on the MANOVA variable list are excluded from the analysis. The MISSING subcommand allows you to include cases with user-missing values. If MISSING is not specified, the defaults are LISTWISE and EXCLUDE.

- The same missing-value treatment is used to process all designs in a single execution of MANOVA.
- If you enter more than one MISSING subcommand, the last subcommand that was entered will be in effect for the entire procedure, including designs that were specified before the last MISSING subcommand.
- Pairwise deletion of missing data is not available in MANOVA.
- Keywords INCLUDE and EXCLUDE are mutually exclusive; either keyword can be specified with LISTWISE.

<b>LISTWISE</b>	<i>Cases with missing values for any variable that is named on the MANOVA variable list are excluded from the analysis.</i> This process is always true in the MANOVA procedure.
<b>EXCLUDE</b>	<i>Both user-missing and system-missing values are excluded.</i> This setting is the default when MISSING is not specified.
<b>INCLUDE</b>	<i>User-missing values are treated as valid.</i> For factors, you must include the missing-value codes within the range that is specified on the MANOVA variable list. It may be necessary to recode these values so that they will be adjacent to the other factor values. System-missing values cannot be included in the analysis.

## MATRIX Subcommand

MATRIX reads and writes matrix data files. MATRIX writes correlation matrices that can be read by subsequent MANOVA procedures.

- Either IN or OUT is required to specify the matrix file in parentheses. When both IN and OUT are used on the same MANOVA procedure, they can be specified on separate MATRIX subcommands or on the same subcommand.
- The matrix materials include the  $N$ , mean, and standard deviation. Documents from the file that form the matrix are not included in the matrix data file.
- MATRIX=IN cannot be used in place of GET or DATA LIST to begin a new command syntax file. MATRIX is a subcommand on MANOVA, and MANOVA cannot run before an active dataset is defined. To begin a new command file and immediately read a matrix, first use GET to retrieve the matrix file, and then specify IN(\*) on MATRIX.
- Records in the matrix data file that is read by MANOVA can be in any order, with the following exceptions: The order of split-file groups cannot be violated, and all CORR vectors must appear contiguously within each split-file group.
- When MANOVA reads matrix materials, it ignores the record containing the total number of cases. In addition, MANOVA skips unrecognized records. MANOVA does not issue a warning when it skips records.

The following two keywords are available on the `MATRIX` subcommand:

- OUT**      *Write a matrix data file.* Specify either a file or an asterisk, and enclose the specification in parentheses. If you specify a file, the file is stored on disk and can be retrieved at any time. If you specify an asterisk (\*) or leave the parentheses empty, the matrix file replaces the active dataset but is not stored on disk unless you use `SAVE` or `XSAVE`.
- IN**        *Read a matrix data file.* If the matrix file *is not* the current active dataset, specify a file in parentheses. If the matrix file *is* the current active dataset, specify an asterisk (\*) or leave the parentheses empty.

### **Format of the Matrix Data File**

The matrix data file includes two special variables: `ROWTYPE_` and `VARNAME_`.

- Variable `ROWTYPE_` is a short string variable having values `N`, `MEAN`, `CORR` (for Pearson correlation coefficients), and `STDDEV`.
- Variable `VARNAME_` is a short string variable whose values are the names of the variables and covariates that are used to form the correlation matrix. When `ROWTYPE_` is `CORR`, `VARNAME_` gives the variable that is associated with that row of the correlation matrix.
- Between `ROWTYPE_` and `VARNAME_` are the factor variables (if any) that are defined in the `BY` portion of the `MANOVA` variable list. (Factor variables receive the system-missing value on vectors that represent pooled values.)
- Remaining variables are the variables that are used to form the correlation matrix.

### **Split Files and Variable Order**

- When split-file processing is in effect, the first variables in the matrix system file will be the split variables, followed by `ROWTYPE_`, the factor variable(s), `VARNAME_`, and then the variables that are used to form the correlation matrix.
- A full set of matrix materials is written for each subgroup that is defined by the split variable(s).
- A split variable cannot have the same variable name as any other variable that is written to the matrix data file.
- If a split file is in effect when a matrix is written, the same split file must be in effect when that matrix is read into another procedure.

### **Additional Statistics**

In addition to the `CORR` values, `MANOVA` always includes the following with the matrix materials:

- The total weighted number of cases used to compute each correlation coefficient.
- A vector of `N`'s for each cell in the data.
- A vector of `MEAN`'s for each cell in the data.
- A vector of pooled standard deviations, `STDDEV`, which is the square root of the within-cells mean square error for each variable.

**Example**

```
GET FILE IRIS.
MANOVA SEPALLEN SEPALWID PETALLEN PETALWID BY TYPE(1,3)
  /MATRIX=OUT(MANMTX) .
```

- MANOVA reads data from the data file *IRIS* and writes one set of matrix materials to the file *MANMTX*.
- The active dataset is still *IRIS*. Subsequent commands are executed on the file *IRIS*.

**Example**

```
GET FILE IRIS.
MANOVA SEPALLEN SEPALWID PETALLEN PETALWID BY TYPE(1,3)
  /MATRIX=OUT(*) .
LIST.
```

- MANOVA writes the same matrix as in the example above. However, the matrix file replaces the active dataset. The `LIST` command is executed on the matrix file (not on the file *IRIS*).

**Example**

```
GET FILE=PRSNL.
FREQUENCIES VARIABLE=AGE.

MANOVA SEPALLEN SEPALWID PETALLEN PETALWID BY TYPE(1,3)
  /MATRIX=IN(MANMTX) .
```

- This example assumes that you want to perform a frequencies analysis on the file *PRSNL* and then use MANOVA to read a different file. The file that you want to read is an existing matrix data file. The external matrix file *MANMTX* is specified in parentheses after `IN` on the `MATRIX` subcommand.
- *MANMTX* does not replace *PRSNL* as the active dataset.

**Example**

```
GET FILE=MANMTX.
MANOVA SEPALLEN SEPALWID PETALLEN PETALWID BY TYPE(1,3)
  /MATRIX=IN(*) .
```

- This example assumes that you are starting a new session and want to read an existing matrix data file. `GET` retrieves the matrix file *MANMTX*.
- An asterisk is specified in parentheses after `IN` on the `MATRIX` subcommand to read the active dataset. You can also leave the parentheses empty to indicate the default.
- If the `GET` command is omitted, an error message is issued.
- If you specify *MANMTX* in parentheses after `IN`, an error message is issued.

**ANALYSIS Subcommand**

`ANALYSIS` allows you to work with a subset of the continuous variables (dependent variable and covariates) that you named on the `MANOVA` variable list. In univariate analysis of variance, you can use `ANALYSIS` to allow factor-by-covariate interaction terms in your model (see the `DESIGN`

subcommand below). You can also use `ANALYSIS` to switch the roles of the dependent variable and a covariate.

- In general, `ANALYSIS` gives you complete control over which continuous variables are to be dependent variables, which continuous variables are to be covariates, and which continuous variables are to be neither.
- `ANALYSIS` specifications are like the `MANOVA` variables specification, except that factors are not named. Enter the dependent variable and, if there are covariates, the keyword `WITH` and the covariates.
- Only variables that are listed as dependent variables or covariates on the `MANOVA` variable list can be entered on the `ANALYSIS` subcommand.
- In a univariate analysis of variance, the most important use of `ANALYSIS` is to omit covariates from the analysis list, thereby making them available for inclusion on `DESIGN` (see the example below and the `DESIGN` subcommand examples).
- For more information about `ANALYSIS`, refer to *MANOVA: Multivariate*.

### Example

```
MANOVA DEP BY FACTOR(1,3) WITH COV
  /ANALYSIS DEP
  /DESIGN FACTOR, COV, FACTOR BY COV.
```

- `COV`, a continuous variable, is included on the `MANOVA` variable list as a covariate.
- `COV` is not mentioned on `ANALYSIS`, so it will not be included in the model as a dependent variable or covariate. `COV` can, therefore, be explicitly included on the `DESIGN` subcommand.
- `DESIGN` includes the main effects of `FACTOR` and `COV` and the `FACTOR` by `COV` interaction.

## DESIGN Subcommand

`DESIGN` specifies the effects that are included in a specific model. `DESIGN` must be the last subcommand entered for any model.

The cells in a design are defined by all of the possible combinations of levels of the factors in that design. The number of cells equals the product of the number of levels of all the factors. A design is *balanced* if each cell contains the same number of cases. `MANOVA` can analyze both balanced and unbalanced designs.

- Specify a list of terms to be included in the model, separated by spaces or commas.
- The default design, if the `DESIGN` subcommand is omitted or is specified by itself, is a full factorial model containing all main effects and all orders of factor-by-factor interaction.
- If the last subcommand that is specified is not `DESIGN`, a default full factorial design is estimated.
- To include a term for the main effect of a factor, enter the name of the factor on the `DESIGN` subcommand.
- To include a term for an interaction between factors, use the keyword `BY` to join the factors that are involved in the interaction.



- Terms are entered into the model in the order in which you list them on `DESIGN`. If you have specified `SEQUENTIAL` on the `METHOD` subcommand to partition the sums of squares in a hierarchical fashion, this order may affect the significance tests.
- You can specify other types of terms in the model, as described in the following sections.
- Multiple `DESIGN` subcommands are accepted. An analysis of one model is produced for each `DESIGN` subcommand.

### Example

```
MANOVA Y BY A(1,2) B(1,2) C(1,3)
  /DESIGN
  /DESIGN A, B, C
  /DESIGN A, B, C, A BY B, A BY C.
```

- The first `DESIGN` produces the default full factorial design, with all main effects and interactions for factors *A*, *B*, and *C*.
- The second `DESIGN` produces an analysis with main effects only for *A*, *B*, and *C*.
- The third `DESIGN` produces an analysis with main effects and the interactions between *A* and the other two factors. The interaction between *B* and *C* is not in the design, nor is the interaction between all three factors.

## Partitioned Effects: Number in Parentheses

You can specify a number in parentheses following a factor name on the `DESIGN` subcommand to identify individual degrees of freedom or partitions of the degrees of freedom that are associated with an effect.

- If you specify `PARTITION`, the number refers to a partition. Partitions can include more than one degree of freedom (see [PARTITION Subcommand on p. 1077](#)). For example, if the first partition of *SEED* includes two degrees of freedom, the term `SEED(1)` on a `DESIGN` subcommand tests the two degrees of freedom.
- If you do not use `PARTITION`, the number refers to a single degree of freedom that is associated with the effect.
- The number refers to an individual level for a factor if that factor follows the keyword `WITHIN` or `MWITHIN` (see the sections about nested effects and pooled effects below).
- A factor has one less degree of freedom than it has levels or values.

### Example

```
MANOVA YIELD BY SEED(1,4) WITH RAINFALL
  /PARTITION(SEED)=(2,1)
  /DESIGN=SEED(1) SEED(2).
```

- Factor *SEED* is subdivided into two partitions, one partition containing the first two degrees of freedom and the other partition containing the last degree of freedom.
- The two partitions of *SEED* are treated as independent effects.

### ***Nested Effects: WITHIN Keyword***

Use the `WITHIN` keyword (alias `w`) to nest the effects of one factor within the effects of another factor or an interaction term.

#### ***Example***

```
MANOVA YIELD BY SEED(1,4) FERT(1,3) PLOT (1,4)
  /DESIGN = FERT WITHIN SEED BY PLOT.
```

- The three factors in this example are type of seed (*SEED*), type of fertilizer (*FERT*), and location of plots (*PLOT*).
- The `DESIGN` subcommand nests the effects of *FERT* within the interaction term of *SEED* by *PLOT*. The levels of *FERT* are considered distinct for each combination of levels of *SEED* and *PLOT*.

### ***Simple Effects: WITHIN and MWITHIN Keywords***

A factor can be nested within one specific level of another factor by indicating the level in parentheses. This process allows you to estimate simple effects or the effect of one factor within only one level of another factor. Simple effects can be obtained for higher-order interactions as well.

Use `WITHIN` to request simple effects of between-subjects factors.

#### ***Example***

```
MANOVA YIELD BY SEED(2,4) FERT(1,3) PLOT (1,4)
  /DESIGN = FERT WITHIN SEED (1).
```

- This example requests the simple effect of *FERT* within the first level of *SEED*.
- The number (*n*) specified after a `WITHIN` factor refers to the level of that factor. The value is the ordinal position, which is not necessarily the value of that level. In this example, the first level is associated with value 2.
- The number does *not* refer to the number of partitioned effects (see [PARTITION Subcommand on p. 1077](#)).

#### ***Example***

```
MANOVA YIELD BY SEED(2,4) FERT(1,3) PLOT (3,5)
  /DESIGN = FERT WITHIN PLOT(1) WITHIN SEED(2)
```

- This example requests the effect of *FERT* within the second *SEED* level of the first *PLOT* level.
- The second *SEED* level is associated with value 3, and the first *PLOT* level is associated with value 3.

Use `MWITHIN` to request simple effects of within-subjects factors in repeated measures analysis (see [MWITHIN Keyword for Simple Effects on p. 1116](#)).

### ***Pooled Effects: Plus Sign***

To pool different effects for the purpose of significance testing, join the effects with a plus sign (+). A single test is made for the combined effect of the pooled terms.

- The keyword `BY` is evaluated before effects are pooled together.
- Parentheses are not allowed for changing the order of evaluation. For example, it is illegal to specify `(A + B) BY C`. You must specify `/DESIGN=A BY C + B BY C`.

#### ***Example***

```
MANOVA Y BY A(1,3) B(1,4) WITH X
/ANALYSIS=Y
/DESIGN=A, B, A BY B, A BY X + B BY X + A BY B BY X.
```

- This example shows how to test homogeneity of regressions in a two-way analysis of variance.
- The + signs are used to produce a pooled test of all interactions involving the covariate *X*. If this test is significant, the assumption of homogeneity of variance is questionable.

### ***MUPLUS Keyword***

`MUPLUS` combines the constant term ( $\mu$ ) in the model with the term that is specified after it. The normal use of this specification is to obtain parameter estimates that represent weighted means for the levels of some factor. For example, `MUPLUS SEED` represents the constant, or overall, mean plus the effect for each level of *SEED*. The significance of such effects is usually uninteresting, but the parameter estimates represent the weighted means for each level of *SEED*, adjusted for any covariates in the model.

- `MUPLUS` cannot appear more than once on a given `DESIGN` subcommand.
- `MUPLUS` is the only way to get standard errors for the predicted mean for each level of the specified factor.
- Parameter estimates are not displayed by default; you must explicitly request them on the `PRINT` subcommand or via a `CONTRAST` subcommand.
- You can obtain the unweighted mean by specifying the full factorial model, excluding those terms that are contained by an effect, and prefixing the effect whose mean is to be found by `MUPLUS`.

### ***Effects of Continuous Variables***

Usually you name factors but not covariates on the `DESIGN` subcommand. The linear effects of covariates are removed from the dependent variable before the design is tested. However, the design can include variables that are measured at the interval level and originally named as covariates or as additional dependent variables.

- Continuous variables on a `DESIGN` subcommand must be named as dependents or covariates on the `MANOVA` variable list.

- Before you can name a continuous variable on a `DESIGN` subcommand, you must supply an `ANALYSIS` subcommand that does *not* name the variable. This action excludes it from the analysis as a dependent variable or covariate and makes it eligible for inclusion on `DESIGN`.
- You can use the keyword `POOL(varlist)` to pool more than one continuous variable into a single effect (provided that the continuous variables are all excluded on an `ANALYSIS` subcommand). For a single continuous variable, `POOL(VAR)` is equivalent to `VAR`.
- The `TO` convention in the variable list for `POOL` refers to the order of continuous variables (dependent variables and covariates) on the original `MANOVA` variable list, which is not necessarily their order on the active dataset. This use is the only allowable use of the keyword `TO` on a `DESIGN` subcommand.
- You can specify interaction terms between factors and continuous variables. If `FAC` is a factor and `COV` is a covariate that has been omitted from an `ANALYSIS` subcommand, `FAC BY COV` is a valid specification on a `DESIGN` statement.
- You cannot specify an interaction between two continuous variables. Use the `COMPUTE` command to create a variable representing the interaction prior to `MANOVA`.

### Example

\* This example tests whether the regression of the dependent variable `Y` on the two variables `X1` and `X2` is the same across all the categories of the factors `AGE` and `TREATMNT`.

```
MANOVA Y BY AGE(1,5) TREATMNT(1,3) WITH X1, X2
  /ANALYSIS = Y
  /DESIGN = POOL(X1,X2),
            AGE, TREATMNT, AGE BY TREATMNT,
            POOL(X1,X2) BY AGE + POOL(X1,X2) BY TREATMNT
            + POOL(X1,X2) BY AGE BY TREATMNT.
```

- `ANALYSIS` excludes `X1` and `X2` from the standard treatment of covariates so that they can be used in the design.
- `DESIGN` includes five terms. `POOL(X1,X2)`, the overall regression of the dependent variable on `X1` and `X2`, is entered first, followed by the two factors and their interaction.
- The last term is the test for equal regressions. It consists of three factor-by-continuous-variable interactions pooled together. `POOL(X1,X2) BY AGE` is the interaction between `AGE` and the combined effect of the continuous variables `X1` and `X2`. It is combined with similar interactions between `TREATMNT` and the continuous variables and between the `AGE` by `TREATMNT` interaction and the continuous variables.
- If the last term is not statistically significant, there is no evidence that the regression of `Y` on `X1` and `X2` is different across any combination of the categories of `AGE` and `TREATMNT`.

### Error Terms for Individual Effects

The “error” sum of squares against which terms in the design are tested is specified on the `ERROR` subcommand. For any particular term on a `DESIGN` subcommand, you can specify a different error term to be used in the analysis of variance. To do so, name the term followed by the keyword `VS` (or `AGAINST`) and the error term keyword.

- To test a term against only the within-cells sum of squares, specify the term followed by `VS WITHIN` on the `DESIGN` subcommand. For example, `GROUP VS WITHIN` tests the effect of the factor *GROUP* against only the within-cells sum of squares. For most analyses, this term is the default error term.
- To test a term against only the residual sum of squares (the sum of squares for all terms that are not included in your `DESIGN`), specify the term followed by `VS RESIDUAL`.
- To test against the combined within-cells and residual sums of squares, specify the term followed by `VS WITHIN+RESIDUAL`.
- To test against any other sum of squares in the analysis of variance, include a term corresponding to the desired sum of squares in the design and assign it to an integer between 1 and 10. You can then test against the number of the error term. It is often convenient to test against the term before you define it. This process is perfectly acceptable as long as you define the error term on the same `DESIGN` subcommand.

### Example

```
MANOVA DEP BY A, B, C (1,3)
  /DESIGN=A VS 1,
    B WITHIN A = 1 VS 2,
    C WITHIN B WITHIN A = 2 VS WITHIN.
```

- In this example, the factors *A*, *B*, and *C* are completely nested; levels of *C* occur within levels of *B*, which occur within levels of *A*. Each factor is tested against everything within it.
- *A*, the outermost factor, is tested against the *B* within *A* sum of squares, to see if it contributes anything beyond the effects of *B* within each of its levels. The *B* within *A* sum of squares is defined as error term number 1.
- *B* nested within *A*, in turn, is tested against error term number 2, which is defined as the *C* within *B* within *A* sum of squares.
- Finally, *C* nested within *B* nested within *A* is tested against the within-cells sum of squares.

User-defined error terms are specified by simply inserting = *n* after a term, where *n* is an integer from 1 to 10. The equals sign is required. Keywords that are used in building a design term, such as `BY` or `WITHIN`, are evaluated first. For example, error term number 2 in the above example consists of the entire term `C WITHIN B WITHIN A`. An error-term *number*, but not an error-term *definition*, can follow the keyword `VS`.

### CONSTANT Keyword

By default, the constant (grand mean) term is included as the first term in the model.

- If you have specified `NOCONSTANT` on the `METHOD` subcommand, a constant term will not be included in any design unless you request it with the `CONSTANT` keyword on `DESIGN`.
- You can specify an error term for the constant.
- A factor named `CONSTANT` will not be recognized on the `DESIGN` subcommand.

## **References**

Bock, R. D. 1985. *Multivariate statistical methods in behavioral research*. Chicago: Scientific Software, Inc..

Cohen, J. 1977. *Statistical power analysis for the behavioral sciences*. San Diego, California: Academic Press.

Hays, W. L. 1981. *Statistics*, 3rd ed. New York: Holt, Rinehart, and Winston.

Timm, N. H. 1975. *Multivariate statistics: With applications in education and psychology*. Monterey, California: Brooks/Cole.

# MANOVA: Multivariate

MANOVA is available in the Advanced Statistics option.

```
MANOVA dependent varlist [BY factor list (min,max) [factor list...]]
      [WITH covariate list]

[/TRANSFORM [(dependent varlist [/dependent varlist])]=
      [ORTHONORM] [{CONTRAST}] {DEVIATIONS (refcat)      }
      {BASIS      } {DIFFERENCE }
      {HELMERT    }
      {SIMPLE (refcat) }
      {REPEATED   }
      {POLYNOMIAL[({1,2,3...})] }
      {metric     }
      {SPECIAL (matrix)  }

[/RENAME={newname} {newname}...]
      {*          } {*          }

[/{PRINT  }=[HOMOGENEITY [(BOXM)]]
  {NOPRINT} [ERROR [(COV) [COR] [SSCP] [STDDEV]]]
  [SIGNIF [(MULTIV**) [EIGEN] [DIMENR]
          [UNIV**] [HYPOTH] [STEPDOWN] [BRIEF]]]
  [TRANSFORM] ]

[/PCOMPS=[COR] [COV] [ROTATE(rotttype)]
  [NCOMP(n)] [MINEIGEN(eigen-cut)] [ALL]]

[/PLOT=[ZCORR]]

[/DISCRIM [RAW] [STAN] [ESTIM] [COR] [ALL]
  [ROTATE(rotttype)] [ALPHA(0.25**)]]
  {a          }

[/POWER=[T(0.05**)] [F(0.05**)] [{APPROXIMATE}]
  {a          } {a          } {EXACT          }

[/CINTERVAL=[MULTIVARIATE ({ROY          })]
  {PILLAI          }
  {BONFER          }
  {HOTELLING       }
  {WILKS          }

[/ANALYSIS [(UNCONDITIONAL**)]]=[(dependent varlist
  {CONDITIONAL      } [WITH covariate varlist]
  [/dependent varlist...])][WITH varlist]]

[/DESIGN...]*
```

\* The DESIGN subcommand has the same syntax as is described in *MANOVA: Univariate*.

\*\*Default if the subcommand or keyword is omitted.

## Example

```
MANOVA SCORE1 TO SCORE4 BY METHOD(1,3).
```

## Overview

This section discusses the subcommands that are used in multivariate analysis of variance and covariance designs with several interrelated dependent variables. The discussion focuses on subcommands and keywords that do not apply, or apply in different manners, to univariate analyses. It does not contain information on all of the subcommands you will need to specify the design. For subcommands not covered here, see *MANOVA: Univariate*.

### Options

**Dependent Variables and Covariates.** You can specify subsets and reorder the dependent variables and covariates using the `ANALYSIS` subcommand. You can specify linear transformations of the dependent variables and covariates using the `TRANSFORM` subcommand. When transformations are performed, you can rename the variables using the `RENAME` subcommand and request the display of a transposed transformation matrix currently in effect using the `PRINT` subcommand.

**Optional Output.** You can request or suppress output on the `PRINT` and `NOPRINT` subcommands. Additional output appropriate to multivariate analysis includes error term matrices, Box's  $M$  statistic, multivariate and univariate  $F$  tests, and other significance analyses. You can also request predicted cell means for specific dependent variables on the `PMEANS` subcommand, produce a canonical discriminant analysis for each effect in your model with the `DISCRIM` subcommand, specify a principal components analysis of each error sum-of-squares and cross-product matrix in a multivariate analysis on the `PCOMPS` subcommand, display multivariate confidence intervals using the `CINTERVAL` subcommand, and generate a half-normal plot of the within-cells correlations among the dependent variables with the `PLOT` subcommand.

### Basic Specification

- The basic specification is a variable list identifying the dependent variables, with the factors (if any) named after `BY` and the covariates (if any) named after `WITH`.
- By default, `MANOVA` produces multivariate and univariate  $F$  tests.

### Subcommand Order

- The variable list must be specified first.
- Subcommands applicable to a specific design must be specified before that `DESIGN` subcommand. Otherwise, subcommands can be used in any order.

### Syntax Rules

- All syntax rules applicable to univariate analysis also apply to multivariate analysis.
- If you enter one of the multivariate specifications in a univariate analysis, `MANOVA` ignores it.

### Limitations

- Maximum of 20 factors.
- Memory requirements depend primarily on the number of cells in the design. For the default full factorial model, the number of cells equals the product of the number of levels or categories in each factor.



## **MANOVA Variable List**

- Multivariate MANOVA calculates statistical tests that are valid for analyses of dependent variables that are correlated with one another. The dependent variables must be specified first.
- The factor and covariate lists follow the same rules as in univariate analyses.
- If the dependent variables are uncorrelated, the univariate significance tests have greater statistical power.

## **TRANSFORM Subcommand**

TRANSFORM performs linear transformations of some or all of the continuous variables (dependent variables and covariates). Specifications on TRANSFORM include an optional list of variables to be transformed, optional keywords to describe how to generate a transformation matrix from the specified contrasts, and a required keyword specifying the transformation contrasts.

- Transformations apply to all subsequent designs unless replaced by another TRANSFORM subcommand.
- TRANSFORM subcommands are not cumulative. Only the transformation specified most recently is in effect at any time. You can restore the original variables in later designs by specifying SPECIAL with an identity matrix.
- You should not use TRANSFORM when you use the WSFACTORS subcommand to request repeated measures analysis; a transformation is automatically performed in repeated measures analysis (see *MANOVA: Repeated Measures* on p. 1110).
- Transformations are in effect for the duration of the MANOVA procedure only. After the procedure is complete, the original variables remain in the active dataset.
- By default, the transformation matrix is not displayed. Specify the keyword TRANSFORM on the PRINT subcommand to see the matrix generated by the TRANSFORM subcommand.
- If you do not use the RENAME subcommand with TRANSFORM, the variables specified on TRANSFORM are renamed temporarily (for the duration of the procedure) as *T1*, *T2*, and so on. Explicit use of RENAME is recommended.
- Subsequent references to transformed variables should use the new names. The only exception is when you supply a VARIABLES specification on the OMEANS subcommand after using TRANSFORM. In this case, specify the original names. OMEANS displays observed means of original variables. See *OMEANS Subcommand* on p. 1082.

## **Variable Lists**

- By default, MANOVA applies the transformation you request to all continuous variables (dependent variables and covariates).
- You can enter a variable list in parentheses following the TRANSFORM subcommand. If you do, only the listed variables are transformed.

- You can enter multiple variable lists, separated by slashes, within a single set of parentheses. Each list must have the same number of variables, and the lists must not overlap. The transformation is applied separately to the variables on each list.
- In designs with covariates, transform only the dependent variables, or, in some designs, apply the same transformation separately to the dependent variables and the covariates.

### ***CONTRAST, BASIS, and ORTHONORM Keywords***

You can control how the transformation matrix is to be generated from the specified contrasts. If none of these three keywords is specified on TRANSFORM, the default is CONTRAST.

<b>CONTRAST</b>	<i>Generate the transformation matrix directly from the contrast matrix specified (see <a href="#">CONTRAST Subcommand on p. 1075</a>). This is the default.</i>
<b>BASIS</b>	<i>Generate the transformation matrix from the one-way basis matrix corresponding to the specified contrast matrix. BASIS makes a difference only if the transformation contrasts are not orthogonal.</i>
<b>ORTHONORM</b>	<i>Orthonormalize the transformation matrix by rows before use. MANOVA eliminates redundant rows. By default, orthonormalization is not done.</i>

- CONTRAST and BASIS are alternatives and are mutually exclusive.
- ORTHONORM is independent of the CONTRAST/BASIS choice; you can enter it before or after either of those keywords.

### ***Transformation Methods***

To specify a transformation method, use one of the following keywords available on the TRANSFORM subcommand. Note that these are identical to the keywords available for the CONTRAST subcommand (see [CONTRAST Subcommand on p. 1075](#)). However, in univariate designs, they are applied to the different levels of a factor. Here they are applied to the continuous variables in the analysis. This reflects the fact that the different dependent variables in a multivariate MANOVA setup can often be thought of as corresponding to different levels of some factor.

- The transformation keyword (and its specifications, if any) must follow all other specifications on the TRANSFORM subcommand.

<b>DEVIATION</b>	<i>Deviations from the mean of the variables being transformed. The first transformed variable is the mean of all variables in the transformation. Other transformed variables represent deviations of individual variables from the mean. One of the original variables (by default, the last) is omitted as redundant. To omit a variable other than the last, specify the number of the variable to be omitted in parentheses after the DEVIATION keyword. For example, /TRANSFORM (A B C) = DEVIATION(1) omits A and creates variables representing the mean, the deviation of B from the mean, and the deviation of C from the mean. A DEVIATION transformation is not orthogonal.</i>
<b>DIFFERENCE</b>	<i>Difference or reverse Helmert transformation. The first transformed variable is the mean of the original variables. Each of the original variables except the first is then transformed by subtracting the mean of those (original) variables that precede it. A DIFFERENCE transformation is orthogonal.</i>

<b>HELMERT</b>	<i>Helmert transformation.</i> The first transformed variable is the mean of the original variables. Each of the original variables except the last is then transformed by subtracting the mean of those (original) variables that follow it. A HELMERT transformation is orthogonal.
<b>SIMPLE</b>	<i>Each original variable, except the last, is compared to the last of the original variables.</i> To use a variable other than the last as the omitted reference variable, specify its number in parentheses following the keyword SIMPLE. For example, <code>/TRANSFORM(A B C) = SIMPLE(2)</code> specifies the second variable, <i>B</i> , as the reference variable. The three transformed variables represent the mean of <i>A</i> , <i>B</i> , and <i>C</i> , the difference between <i>A</i> and <i>B</i> , and the difference between <i>C</i> and <i>B</i> . A SIMPLE transformation is not orthogonal.
<b>POLYNOMIAL</b>	<i>Orthogonal polynomial transformation.</i> The first transformed variable represents the mean of the original variables. Other transformed variables represent the linear, quadratic, and higher-degree components. By default, values of the original variables are assumed to represent equally spaced points. You can specify unequal spacing by entering a metric consisting of one integer for each variable in parentheses after the keyword POLYNOMIAL. For example, <code>/TRANSFORM(RESP1 RESP2 RESP3) = POLYNOMIAL(1,2,4)</code> might indicate that three response variables correspond to levels of some stimulus that are in the proportion 1:2:4. The default metric is always (1,2,..., <i>k</i> ), where <i>k</i> variables are involved. Only the relative differences between the terms of the metric matter: (1,2,4) is the same metric as (2,3,5) or (20,30,50) because in each instance the difference between the second and third numbers is twice the difference between the first and second.
<b>REPEATED</b>	<i>Comparison of adjacent variables.</i> The first transformed variable is the mean of the original variables. Each additional transformed variable is the difference between one of the original variables and the original variable that followed it. Such transformed variables are often called <i>difference scores</i> . A REPEATED transformation is not orthogonal.
<b>SPECIAL</b>	<i>A user-defined transformation.</i> After the keyword SPECIAL, enter a square matrix in parentheses with as many rows and columns as there are variables to transform. MANOVA multiplies this matrix by the vector of original variables to obtain the transformed variables (see the examples below).

**Example**

```
MANOVA X1 TO X3 BY A(1,4)
  /TRANSFORM(X1 X2 X3) = SPECIAL( 1  1  1,
                                1  0 -1,
                                2 -1 -1)
  /DESIGN.
```

- The given matrix will be post-multiplied by the three continuous variables (considered as a column vector) to yield the transformed variables. The first transformed variable will therefore equal  $X1 + X2 + X3$ , the second will equal  $X1 - X3$ , and the third will equal  $2X1 - X2 - X3$ .
- The variable list is optional in this example since all three interval-level variables are transformed.
- You do not need to enter the matrix one row at a time, as shown above. For example, `/TRANSFORM = SPECIAL(1 1 1 1 0 -1 2 -1 -1)` is equivalent to the TRANSFORM specification in the above example.
- You can specify a repetition factor followed by an asterisk to indicate multiple consecutive elements of a SPECIAL transformation matrix. For example,

```
/TRANSFORM = SPECIAL (4*1 0 -1 2 2*-1)
```

is again equivalent to the TRANSFORM specification above.

### Example

```
MANOVA X1 TO X3, Y1 TO Y3 BY A(1,4)
/TRANSFORM (X1 X2 X3/Y1 Y2 Y3) = SPECIAL( 1 1 1,
                                             1 0 -1,
                                             2 -1 -1)
/DESIGN.
```

- Here the same transformation shown in the previous example is applied to  $X1$ ,  $X2$ ,  $X3$  and to  $Y1$ ,  $Y2$ ,  $Y3$ .

## RENAME Subcommand

Use RENAME to assign new names to transformed variables. Renaming variables after a transformation is strongly recommended. If you transform but do not rename the variables, the names  $T1$ ,  $T2$ , ...,  $Tn$  are used as names for the transformed variables.

- Follow RENAME with a list of new variable names.
- You must enter a new name for each dependent variable and covariate on the MANOVA variable list.
- Enter the new names in the order in which the original variables appeared on the MANOVA variable list.
- To retain the original name for one or more of the interval variables, you can either enter an asterisk or reenter the old name as the new name.
- References to dependent variables and covariates on subcommands following RENAME must use the new names. The original names will not be recognized within the MANOVA procedure. The only exception is the OMEANS subcommand, which displays observed means of the original (untransformed) variables. Use the original names on OMEANS.
- The new names exist only during the MANOVA procedure that created them. They do not remain in the active dataset after the procedure is complete.

### Example

```
MANOVA A, B, C, V4, V5 BY TREATMNT(1,3)
/TRANSFORM(A, B, C) = REPEATED
/RENAME = MEANABC, AMINUSB, BMINUSC, *, *
/DESIGN.
```

- The REPEATED transformation produces three transformed variables, which are then assigned mnemonic names *MEANABC*, *AMINUSB*, and *BMINUSC*.
- $V4$  and  $V5$  retain their original names.

### Example

```
MANOVA WT1, WT2, WT3, WT4 BY TREATMNT(1,3) WITH COV
/TRANSFORM (WT1 TO WT4) = POLYNOMIAL
/RENAME = MEAN, LINEAR, QUAD, CUBIC, *
```

```
/ANALYSIS = MEAN, LINEAR, QUAD WITH COV
/DESIGN.
```

- After the polynomial transformation of the four *WT* variables, `RENAME` assigns appropriate names to the various trends.
- Even though only four variables were transformed, `RENAME` applies to all five continuous variables. An asterisk is required to retain the original name for *COV*.
- The `ANALYSIS` subcommand following `RENAME` refers to the interval variables by their new names.

## ***PRINT and NOPRINT Subcommands***

All of the `PRINT` specifications described in *MANOVA: Univariate* are available in multivariate analyses. The following additional output can be requested. To suppress any optional output, specify the appropriate keyword on `NOPRINT`.

<b>ERROR</b>	<i>Error matrices.</i> Three types of matrices are available.
<b>SIGNIF</b>	<i>Significance tests.</i>
<b>TRANSFORM</b>	<i>Transformation matrix.</i> It is available if you have transformed the dependent variables with the <code>TRANSFORM</code> subcommand.
<b>HOMOGENEITY</b>	<i>Test for homogeneity of variance.</i> <code>BOXM</code> is available for multivariate analyses.

## ***ERROR Keyword***

In multivariate analysis, error terms consist of entire matrices, not single values. You can display any of the following error matrices on a `PRINT` subcommand by requesting them in parentheses following the keyword `ERROR`. If you specify `ERROR` by itself, without further specifications, the default is to display `COV` and `COR`.

<b>SSCP</b>	<i>Error sums-of-squares and cross-products matrix.</i>
<b>COV</b>	<i>Error variance-covariance matrix.</i>
<b>COR</b>	<i>Error correlation matrix with standard deviations on the diagonal.</i> This also displays the determinant of the matrix and Bartlett's test of sphericity, a test of whether the error correlation matrix is significantly different from an identity matrix.

## ***SIGNIF Keyword***

You can request any of the optional output listed below by entering the appropriate specification in parentheses after the keyword `SIGNIF` on the `PRINT` subcommand. Further specifications for `SIGNIF` are described in *MANOVA: Repeated Measures*.

<b>MULTIV</b>	<i>Multivariate F tests for group differences.</i> <code>MULTIV</code> is always printed unless explicitly suppressed with the <code>NOPRINT</code> subcommand.
<b>EIGEN</b>	<i>Eigenvalues of the <math>S_k S_e^{-1}</math> matrix.</i> This matrix is the product of the hypothesis sums-of-squares and cross-products ( <code>SSCP</code> ) matrix and the inverse of the error <code>SSCP</code> matrix. To print <code>EIGEN</code> , request it on the <code>PRINT</code> subcommand.

<b>DIMENR</b>	<i>A dimension-reduction analysis.</i> To print DIMENR, request it on the PRINT subcommand.
<b>UNIV</b>	<i>Univariate F tests.</i> UNIV is always printed except in repeated measures analysis. If the dependent variables are uncorrelated, univariate tests have greater statistical power. To suppress UNIV, use the NOPRINT subcommand.
<b>HYPOTH</b>	<i>The hypothesis SSCP matrix.</i> To print HYPOTH, request it on the PRINT subcommand.
<b>STEPDOWN</b>	<i>Roy-Bargmann stepdown F tests.</i> To print STEPDOWN, request it on the PRINT subcommand.
<b>BRIEF</b>	<i>Abbreviated multivariate output.</i> This is similar to a univariate analysis of variance table but with Wilks' multivariate <i>F</i> approximation (lambda) replacing the univariate <i>F</i> . BRIEF overrides any of the SIGNIF specifications listed above.
<b>SINGLEDF</b>	<i>Significance tests for the single degree of freedom making up each effect for ANOVA tables.</i> Results are displayed separately corresponding to each hypothesis degree of freedom. For more information, see the topic SIGNIF Keyword on p. 1081.

- If neither PRINT nor NOPRINT is specified, MANOVA displays the results corresponding to MULTIV and UNIV for a multivariate analysis not involving repeated measures.
- If you enter any specification except BRIEF or SINGLEDF for SIGNIF on the PRINT subcommand, the requested output is displayed in addition to the default.
- To suppress the default, specify the keyword(s) on the NOPRINT subcommand.

### **TRANSFORM Keyword**

The keyword TRANSFORM specified on PRINT displays the transposed transformation matrix in use for each subsequent design. This matrix is helpful in interpreting a multivariate analysis in which the interval-level variables have been transformed with either TRANSFORM or WSFACORS.

- The matrix displayed by this option is the transpose of the transformation matrix.
- Original variables correspond to the rows of the matrix, and transformed variables correspond to the columns.
- A **transformed variable** is a linear combination of the original variables using the coefficients displayed in the column corresponding to that transformed variable.

### **HOMOGENEITY Keyword**

In addition to the BARTLETT and COCHRAN specifications described in *MANOVA: Univariate*, the following test for homogeneity is available for multivariate analyses:

<b>BOXM</b>	<i>Box's M statistic.</i> BOXM requires at least two dependent variables. If there is only one dependent variable when BOXM is requested, MANOVA prints Bartlett-Box <i>F</i> test statistic and issues a note.
-------------	---

## ***PLOT Subcommand***

In addition to the plots described in *MANOVA: Univariate*, the following is available for multivariate analyses:

**ZCORR**      *A half-normal plot of the within-cells correlations among the dependent variables. MANOVA first transforms the correlations using Fisher's Z transformation. If errors for the dependent variables are uncorrelated, the plotted points should lie close to a straight line.*

## ***PCOMPS Subcommand***

PCOMPS requests a principal components analysis of each error matrix in a multivariate analysis. You can display the principal components of the error correlation matrix, the error variance-covariance matrix, or both. These principal components are corrected for differences due to the factors and covariates in the MANOVA analysis. They tend to be more useful than principal components extracted from the raw correlation or covariance matrix when there are significant group differences between the levels of the factors or when a significant amount of error variance is accounted for by the covariates. You can specify any of the keywords listed below on PCOMPS.

**COR**                      *Principal components analysis of the error correlation matrix.*  
**COV**                      *Principal components analysis of the error variance-covariance matrix.*  
**ROTATE**                *Rotate the principal components solution. By default, no rotation is performed. Specify a rotation type (either VARIMAX, EQUAMAX, or QUARTIMAX) in parentheses after the keyword ROTATE. To cancel a rotation specified for a previous design, enter NOROTATE in the parentheses after ROTATE.*  
**NCOMP(n)**              *The number of principal components to rotate. Specify a number in parentheses. The default is the number of dependent variables.*  
**MINEIGEN(n)**         *The minimum eigenvalue for principal component extraction. Specify a cutoff value in parentheses. Components with eigenvalues below the cutoff will not be retained in the solution. The default is 0; all components (or the number specified on NCOMP) are extracted.*  
**ALL**                      *COR, COV, and ROTATE.*

- You must specify either COR or COV (or both). Otherwise, MANOVA will not produce any principal components.
- Both NCOMP and MINEIGEN limit the number of components that are rotated.
- If the number specified on NCOMP is less than two, two components are rotated provided that at least two components have eigenvalues greater than any value specified on MINEIGEN.
- Principal components analysis is computationally expensive if the number of dependent variables is large.

## ***DISCRIM Subcommand***

DISCRIM produces a canonical discriminant analysis for each effect in a design. (For covariates, DISCRIM produces a canonical correlation analysis.) These analyses aid in the interpretation of multivariate effects. You can request the following statistics by entering the appropriate keywords after the subcommand DISCRIM:

<b>RAW</b>	<i>Raw discriminant function coefficients.</i>
<b>STAN</b>	<i>Standardized discriminant function coefficients.</i>
<b>ESTIM</b>	<i>Effect estimates in discriminant function space.</i>
<b>COR</b>	<i>Correlations between the dependent variables and the canonical variables defined by the discriminant functions.</i>
<b>ROTATE</b>	<i>Rotation of the matrix of correlations between dependent and canonical variables. Specify rotation type VARIMAX, EQUAMAX, or QUARTIMAX in parentheses after this keyword.</i>
<b>ALL</b>	<i>RAW, STAN, ESTIM, COR, and ROTATE.</i>

By default, the significance level required for the extraction of a canonical variable is 0.25. You can change this value by specifying the keyword ALPHA and a value between 0 and 1 in parentheses:

<b>ALPHA</b>	<i>The significance level required before a canonical variable is extracted. The default is 0.25. To change the default, specify a decimal number between 0 and 1 in parentheses after ALPHA.</i>
--------------	---

- The correlations between dependent variables and canonical functions are not rotated unless at least two functions are significant at the level defined by ALPHA.
- If you set ALPHA to 1.0, all discriminant functions are reported (and rotated, if you so request).
- If you set ALPHA to 0, no discriminant functions are reported.

## ***POWER Subcommand***

The following specifications are available for POWER in multivariate analysis. For applications of POWER in univariate analysis, see [MANOVA: Univariate](#).

<b>APPROXIMATE</b>	<i>Approximate power values.</i> This is the default. Approximate power values for multivariate tests are derived from procedures presented by Muller and Peterson (Muller and Peterson, 1984). Approximate values are normally accurate to three decimal places and are much cheaper to compute than exact values.
<b>EXACT</b>	<i>Exact power values.</i> Exact power values for multivariate tests are computed from the noncentral <i>F</i> distribution. Exact multivariate power values will be displayed only if there is one hypothesis degree of freedom, where all the multivariate criteria have identical power.

- For information on the multivariate generalizations of power and effect size, see (Muller et al., 1984), (Green, 1978), and (Huberty, 1972).



## **CINTERVAL Subcommand**

In addition to the specifications described in *MANOVA: Univariate*, the keyword `MULTIVARIATE` is available for multivariate analysis. You can specify a type in parentheses after the `MULTIVARIATE` keyword. The following type keywords are available on `MULTIVARIATE`:

<b>ROY</b>	<i>Roy's largest root.</i> An approximation given by Pillai (Pillai, 1967) is used. This approximation is accurate for upper percentage points (0.95 to 1), but it is not as good for lower percentage points. Thus, for Roy intervals, the user is restricted to the range 0.95 to 1.
<b>PILLAI</b>	<i>Pillai's trace.</i> The intervals are computed by approximating the percentage points with percentage points of the $F$ distribution.
<b>WILKS</b>	<i>Wilks' lambda.</i> The intervals are computed by approximating the percentage points with percentage points of the $F$ distribution.
<b>HOTELLING</b>	<i>Hotelling's trace.</i> The intervals are computed by approximating the percentage points with percentage points of the $F$ distribution.
<b>BONFER</b>	<i>Bonferroni intervals.</i> This approximation is based on Student's $t$ distribution.

- The Wilks', Pillai's, and Hotelling's approximate confidence intervals are thought to match exact intervals across a wide range of alpha levels, especially for large sample sizes (Burns, 1984). Use of these intervals, however, has not been widely investigated.
- To obtain multivariate intervals separately for each parameter, choose individual multivariate intervals. For individual multivariate confidence intervals, the hypothesis degree of freedom is set to 1, in which case Hotelling's, Pillai's, Wilks', and Roy's intervals will be identical and equivalent to those computed from percentage points of Hotelling's  $T^2$  distribution. Individual Bonferroni intervals will differ and, for a small number of dependent variables, will generally be shorter.
- If you specify `MULTIVARIATE` on `CINTERVAL`, you must specify a type keyword. If you specify `CINTERVAL` without any keyword, the default is the same as with univariate analysis—`CINTERVAL` displays individual-univariate confidence intervals at the 0.95 level.

## **ANALYSIS Subcommand**

`ANALYSIS` is discussed in *MANOVA: Univariate* as a means of obtaining factor-by-covariate interaction terms. In multivariate analyses, it is considerably more useful.

- `ANALYSIS` specifies a subset of the continuous variables (dependent variables and covariates) listed on the `MANOVA` variable list and completely redefines which variables are dependent and which are covariates.
- All variables named on an `ANALYSIS` subcommand must have been named on the `MANOVA` variable list. It does not matter whether they were named as dependent variables or as covariates.
- Factors cannot be named on an `ANALYSIS` subcommand.
- After the keyword `ANALYSIS`, specify the names of one or more dependent variables and, optionally, the keyword `WITH` followed by one or more covariates.
- An `ANALYSIS` specification remains in effect for all designs until you enter another `ANALYSIS` subcommand.

- Continuous variables named on the MANOVA variable list but omitted from the ANALYSIS subcommand currently in effect can be specified on the DESIGN subcommand. [For more information, see the topic DESIGN Subcommand on p. 1090.](#)
- You can use an ANALYSIS subcommand to request analyses of several groups of variables provided that the groups do not overlap. Separate the groups of variables with slashes and enclose the entire ANALYSIS specification in parentheses.

### **CONDITIONAL and UNCONDITIONAL Keywords**

When several analysis groups are specified on a single ANALYSIS subcommand, you can control how each list is to be processed by specifying CONDITIONAL or UNCONDITIONAL in the parentheses immediately following the ANALYSIS subcommand. The default is UNCONDITIONAL.

<b>UNCONDITIONAL</b>	<i>Process each analysis group separately, without regard to other lists. This is the default.</i>
<b>CONDITIONAL</b>	<i>Use variables specified in one analysis group as covariates in subsequent analysis groups.</i>

- CONDITIONAL analysis is not carried over from one ANALYSIS subcommand to another.
- You can specify a final covariate list outside the parentheses. These covariates apply to every list within the parentheses, regardless of whether you specify CONDITIONAL or UNCONDITIONAL. The variables on this global covariate list must not be specified in any individual lists.

#### **Example**

```
MANOVA A B C BY FAC(1,4) WITH D, E
  /ANALYSIS = (A, B / C / D WITH E)
  /DESIGN.
```

- The first analysis uses *A* and *B* as dependent variables and uses no covariates.
- The second analysis uses *C* as a dependent variable and uses no covariates.
- The third analysis uses *D* as the dependent variable and uses *E* as a covariate.

#### **Example**

```
MANOVA A, B, C, D, E BY FAC(1,4) WITH F G
  /ANALYSIS = (A, B / C / D WITH E) WITH F G
  /DESIGN.
```

- A final covariate list WITH F G is specified outside the parentheses. The covariates apply to every list within the parentheses.
- The first analysis uses *A* and *B*, with *F* and *G* as covariates.
- The second analysis uses *C*, with *F* and *G* as covariates.
- The third analysis uses *D*, with *E*, *F*, and *G* as covariates.
- Factoring out *F* and *G* is the only way to use them as covariates in all three analyses, since no variable can be named more than once on an ANALYSIS subcommand.

**Example**

```
MANOVA A B C BY FAC(1,3)
  /ANALYSIS(CONDITIONAL) = (A WITH B / C)
  /DESIGN.
```

- In the first analysis,  $A$  is the dependent variable,  $B$  is a covariate, and  $C$  is not used.
- In the second analysis,  $C$  is the dependent variable, and both  $A$  and  $B$  are covariates.

# MANOVA: Repeated Measures

MANOVA is available in the Advanced Statistics option.

```
MANOVA dependent varlist [BY factor list (min,max)[factor list...]  
      [WITH [varying covariate list] [(constant covariate list)]]  
  
      /WSFACTORS = varname (levels) [varname...]  
  
      [/WSDSIGN = [effect effect...]  
  
      [/MEASURE = newname newname...]  
  
      [/RENAME = newname newname...]  
  
      [/{PRINT }=[SIGNIF({AVERF**}) (HF) (GG) (EFSIZE)]]  
      {NOPRINT}          {AVONLY }  
  
      [/DESIGN] *
```

\* The DESIGN subcommand has the same syntax as is described in *MANOVA: Univariate*.

\*\* Default if the subcommand or keyword is omitted.

## Example

```
MANOVA Y1 TO Y4 BY GROUP(1,2)  
      /WSFACTORS=YEAR(4) .
```

## Overview

This section discusses the subcommands that are used in repeated measures designs, in which the dependent variables represent measurements of the same variable (or variables) at different times. This section does not contain information on all subcommands you will need to specify the design. For some subcommands or keywords not covered here, such as DESIGN, see *MANOVA: Univariate*. For information on optional output and the multivariate significance tests available, see *MANOVA: Multivariate*.

- In a simple repeated measures analysis, all dependent variables represent different measurements of the same variable for different values (or levels) of a within-subjects factor. Between-subjects factors and covariates can also be included in the model, just as in analyses not involving repeated measures.
- A **within-subjects factor** is simply a factor that distinguishes measurements made on the same subject or case, rather than distinguishing different subjects or cases.
- MANOVA permits more complex analyses, in which the dependent variables represent levels of two or more within-subjects factors.
- MANOVA also permits analyses in which the dependent variables represent measurements of several variables for the different levels of the within-subjects factors. These are known as **doubly multivariate designs**.

- A repeated measures analysis includes a within-subjects design describing the model to be tested with the within-subjects factors, as well as the usual between-subjects design describing the effects to be tested with between-subjects factors. The default for both types of design is a full factorial model.
- MANOVA always performs an orthonormal transformation of the dependent variables in a repeated measures analysis. By default, MANOVA renames them as *T1*, *T2*, and so forth.

### **Basic Specification**

- The basic specification is a variable list followed by the `WSFACTORS` subcommand.
- By default, MANOVA performs special repeated measures processing. Default output includes `SIGNIF(AVERF)` but not `SIGNIF(UNIV)`. In addition, for any within-subjects effect involving more than one transformed variable, the Mauchly test of sphericity is displayed to test the assumption that the covariance matrix of the transformed variables is constant on the diagonal and zero off the diagonal. The Greenhouse-Geiser epsilon and the Huynh-Feldt epsilon are also displayed for use in correcting the significance tests in the event that the assumption of sphericity is violated.

### **Subcommand Order**

- The list of dependent variables, factors, and covariates must be first.
- `WSFACTORS` must be the first subcommand used after the variable list.

### **Syntax Rules**

- The `WSFACTORS` (within-subjects factors), `WSDESIGN` (within-subjects design), and `MEASURE` subcommands are used only in repeated measures analysis.
- `WSFACTORS` is required for any repeated measures analysis.
- If `WSDESIGN` is not specified, a full factorial within-subjects design consisting of all main effects and interactions among within-subjects factors is used by default.
- The `MEASURE` subcommand is used for doubly multivariate designs, in which the dependent variables represent repeated measurements of more than one variable.
- Do not use the `TRANSFORM` subcommand with the `WSFACTORS` subcommand because `WSFACTORS` automatically causes an orthonormal transformation of the dependent variables.

### **Limitations**

- Maximum of 20 between-subjects factors. There is no limit on the number of measures for doubly multivariate designs.
- Memory requirements depend primarily on the number of cells in the design. For the default full factorial model, this equals the product of the number of levels or categories in each factor.

## **Example**

```
MANOVA Y1 TO Y4 BY GROUP(1,2)
  /WSFACTORS=YEAR(4)
  /CONTRAST(YEAR)=POLYNOMIAL
  /RENAME=CONST, LINEAR, QUAD, CUBIC
  /PRINT=TRANSFORM PARAM(ESTIM)
```

```
/WSDSIGN=YEAR
/DESIGN=GROUP.
```

- **WSFACTORS** immediately follows the **MANOVA** variable list and specifies a repeated measures analysis in which the four dependent variables represent a single variable measured at four levels of the within-subjects factor. The within-subjects factor is called *YEAR* for the duration of the **MANOVA** procedure.
- **CONTRAST** requests polynomial contrasts for the levels of *YEAR*. Because the four variables, *Y1*, *Y2*, *Y3*, and *Y4*, in the active dataset represent the four levels of *YEAR*, the effect is to perform an orthonormal polynomial transformation of these variables.
- **RENAME** assigns names to the dependent variables to reflect the transformation.
- **PRINT** requests that the transformation matrix and the parameter estimates be displayed.
- **WSDSIGN** specifies a within-subjects design that includes only the effect of the *YEAR* within-subjects factor. Because *YEAR* is the only within-subjects factor specified, this is the default design, and **WSDSIGN** could have been omitted.
- **DESIGN** specifies a between-subjects design that includes only the effect of the *GROUP* between-subjects factor. This subcommand could have been omitted.

## MANOVA Variable List

The list of dependent variables, factors, and covariates must be specified first.

- **WSFACTORS** determines how the dependent variables on the **MANOVA** variable list will be interpreted.
- The number of dependent variables on the **MANOVA** variable list must be a multiple of the number of cells in the within-subjects design. If there are six cells in the within-subjects design, each group of six dependent variables represents a single within-subjects variable that has been measured in each of the six cells.
- Normally, the number of dependent variables should equal the number of cells in the within-subjects design multiplied by the number of variables named on the **MEASURE** subcommand (if one is used). If you have more groups of dependent variables than are accounted for by the **MEASURE** subcommand, **MANOVA** will choose variable names to label the output, which may be difficult to interpret.
- Covariates are specified after the keyword **WITH**. You can specify either varying covariates or constant covariates, or both. **Varying covariates**, similar to dependent variables in a repeated measures analysis, represent measurements of the same variable (or variables) at different times while **constant covariates** represent variables whose values remain the same at each within-subjects measurement.
- If you use varying covariates, the number of covariates specified must be an integer multiple of the number of dependent variables.
- If you use constant covariates, you must specify them in parentheses. If you use both constant and varying covariates, constant variates must be specified after all varying covariates.

### Example

```
MANOVA MATH1 TO MATH4 BY METHOD(1,2) WITH PHYS1 TO PHYS4 (SES)
```

```
/WSFACTORS=SEMESTER(4) .
```

- The four dependent variables represent a score measured four times (corresponding to the four levels of *SEMESTER*).
- The four varying covariates *PHYS1* to *PHYS4* represents four measurements of another score.
- *SES* is a constant covariate. Its value does not change over the time covered by the four levels of *SEMESTER*.
- The default contrast (*POLYNOMIAL*) is used.

## ***WSFACTORS Subcommand***

*WSFACTORS* names the within-subjects factors and specifies the number of levels for each.

- For repeated measures designs, *WSFACTORS* must be the first subcommand after the *MANOVA* variable list.
- Only one *WSFACTORS* subcommand is permitted per execution of *MANOVA*.
- Names for the within-subjects factors are specified on the *WSFACTORS* subcommand. Factor names must not duplicate any of the dependent variables, factors, or covariates named on the *MANOVA* variable list.
- If there are more than one within-subjects factors, they must be named in the order corresponding to the order of the dependent variables on the *MANOVA* variable list. *MANOVA* varies the levels of the last-named within-subjects factor most rapidly when assigning dependent variables to within-subjects cells (see the example below).
- Levels of the factors must be represented in the data by the dependent variables named on the *MANOVA* variable list.
- Enter a number in parentheses after each factor to indicate how many levels the factor has. If two or more adjacent factors have the same number of levels, you can enter the number of levels in parentheses after all of them.
- Enter only the number of levels for within-subjects factors, not a range of values.
- The number of cells in the within-subjects design is the product of the number of levels for all within-subjects factors.

### ***Example***

```
MANOVA X1Y1 X1Y2 X2Y1 X2Y2 X3Y1 X3Y2 BY TREATMNT(1,5) GROUP(1,2)
/WSFACTORS=X(3) Y(2)
/DESIGN.
```

- The *MANOVA* variable list names six dependent variables and two between-subjects factors, *TREATMNT* and *GROUP*.
- *WSFACTORS* identifies two within-subjects factors whose levels distinguish the six dependent variables. *X* has three levels and *Y* has two. Thus, there are  $3 \times 2 = 6$  cells in the within-subjects design, corresponding to the six dependent variables.

- Variable *X1Y1* corresponds to levels 1,1 of the two within-subjects factors; variable *X1Y2* corresponds to levels 1,2; *X2Y1* to levels 2,1; and so on up to *X3Y2*, which corresponds to levels 3,2. The first within-subjects factor named, *X*, varies most slowly, and the last within-subjects factor named, *Y*, varies most rapidly on the list of dependent variables.
- Because there is no *WSDSIGN* subcommand, the within-subjects design will include all main effects and interactions: *X*, *Y*, and *X* by *Y*.
- Likewise, the between-subjects design includes all main effects and interactions: *TREATMNT*, *GROUP*, and *TREATMNT* by *GROUP*.
- In addition, a repeated measures analysis always includes interactions between the within-subjects factors and the between-subjects factors. There are three such interactions for each of the three within-subjects effects.

### **CONTRAST for WSFACTORS**

The levels of a within-subjects factor are represented by different dependent variables. Therefore, contrasts between levels of such a factor compare these dependent variables. Specifying the type of contrast amounts to specifying a transformation to be performed on the dependent variables.

- An orthonormal transformation is automatically performed on the dependent variables in a repeated measures analysis.
- To specify the type of orthonormal transformation, use the *CONTRAST* subcommand for the within-subjects factors.
- Regardless of the contrast type you specify, the transformation matrix is orthonormalized before use.
- If you do not specify a contrast type for within-subjects factors, the default contrast type is orthogonal *POLYNOMIAL*. Intrinsically orthogonal contrast types are recommended for within-subjects factors if you wish to examine each degree-of-freedom test. Other orthogonal contrast types are *DIFFERENCE* and *HELMERT*. *MULTIV* and *AVERF* tests are identical, no matter what contrast was specified.
- To perform non-orthogonal contrasts, you must use the *TRANSFORM* subcommand instead of *CONTRAST*. The *TRANSFORM* subcommand is discussed in *MANOVA: Multivariate*.
- When you implicitly request a transformation of the dependent variables with *CONTRAST* for within-subjects factors, the same transformation is applied to any covariates in the analysis. The number of covariates must be an integer multiple of the number of dependent variables.
- You can display the transpose of the transformation matrix generated by your within-subjects contrast using the keyword *TRANSFORM* on the *PRINT* subcommand.

#### **Example**

```
MANOVA SCORE1 SCORE2 SCORE3 BY GROUP(1,4)
  /WSFACTORS=ROUND(3)
  /CONTRAST(ROUND)=DIFFERENCE
  /CONTRAST(GROUP)=DEVIATION
  /PRINT=TRANSFORM PARAM(ESTIM) .
```



- This analysis has one between-subjects factor, *GROUP*, with levels 1, 2, 3, and 4, and one within-subjects factor, *ROUND*, with three levels that are represented by the three dependent variables.
- The first *CONTRAST* subcommand specifies difference contrasts for *ROUND*, the within-subjects factor.
- There is no *WSDSIGN* subcommand, so a default full factorial within-subjects design is assumed. This could also have been specified as *WSDSIGN=ROUND*, or simply *WSDSIGN*.
- The second *CONTRAST* subcommand specifies deviation contrasts for *GROUP*, the between-subjects factor. This subcommand could have been omitted because deviation contrasts are the default.
- *PRINT* requests the display of the transformation matrix generated by the within-subjects contrast and the parameter estimates for the model.
- There is no *DESIGN* subcommand, so a default full factorial between-subjects design is assumed. This could also have been specified as *DESIGN=GROUP*, or simply *DESIGN*.

### ***PARTITION for WSFACORS***

The *PARTITION* subcommand also applies to factors named on *WSFACTORS*. [For more information, see the topic \*PARTITION Subcommand\* on p. 1077.](#)

### ***WSDSIGN Subcommand***

*WSDSIGN* specifies the design for within-subjects factors. Its specifications are like those of the *DESIGN* subcommand, but it uses the within-subjects factors rather than the between-subjects factors.

- The default *WSDSIGN* is a full factorial design, which includes all main effects and all interactions for within-subjects factors. The default is in effect whenever a design is processed without a preceding *WSDSIGN* or when the preceding *WSDSIGN* subcommand has no specifications.
- A *WSDSIGN* specification can include main effects, factor-by-factor interactions, nested terms (term within term), terms using the keyword *MWITHIN*, and pooled effects using the plus sign. The specification is the same as on the *DESIGN* subcommand but involves only within-subjects factors.
- A *WSDSIGN* specification cannot include between-subjects factors or terms based on them, nor does it accept interval-level variables, the keywords *MUPLUS* or *CONSTANT*, or error-term definitions or references.
- The *WSDSIGN* specification applies to all subsequent within-subjects designs until another *WSDSIGN* subcommand is encountered.

#### ***Example***

```
MANOVA JANLO, JANHI, FEBLO, FEBHI, MARLO, MARHI BY SEX(1,2)
  /WSFACTORS MONTH(3) STIMULUS(2)
  /WSDSIGN MONTH, STIMULUS
  /WSDSIGN
  /DESIGN SEX.
```

- There are six dependent variables, corresponding to three months and two different levels of stimulus.
- The dependent variables are named on the MANOVA variable list in such an order that the level of stimulus varies more rapidly than the month. Thus, *STIMULUS* is named last on the WSFACORS subcommand.
- The first WSDSIGN subcommand specifies only the main effects for within-subjects factors. There is no *MONTH* by *STIMULUS* interaction term.
- The second WSDSIGN subcommand has no specifications and, therefore, invokes the default within-subjects design, which includes the main effects and their interaction.

### **MWITHIN Keyword for Simple Effects**

You can use *MWITHIN* on either the *WSDSIGN* or the *DESIGN* subcommand in a model with both between- and within-subjects factors to estimate simple effects for factors nested within factors of the opposite type.

#### **Example**

```
MANOVA WEIGHT1 WEIGHT2 BY TREAT(1,2)
  /WSFACTORS=WEIGHT(2)
  /DESIGN=MWITHIN TREAT(1) MWITHIN TREAT(2)
MANOVA WEIGHT1 WEIGHT2 BY TREAT(1,2)
  /WSFACTORS=WEIGHT(2)
  /WSDSIGN=MWITHIN WEIGHT(1) MWITHIN WEIGHT(2)
  /DESIGN.
```

- The first *DESIGN* tests the simple effects of *WEIGHT* within each level of *TREAT*.
- The second *DESIGN* tests the simple effects of *TREAT* within each level of *WEIGHT*.

### **MEASURE Subcommand**

In a doubly multivariate analysis, the dependent variables represent multiple variables measured under the different levels of the within-subjects factors. Use *MEASURE* to assign names to the variables that you have measured for the different levels of within-subjects factors.

- Specify a list of one or more variable names to be used in labeling the averaged results. If no within-subjects factor has more than two levels, *MEASURE* has no effect.
- The number of dependent variables on the *DESIGN* subcommand should equal the product of the number of cells in the within-subjects design and the number of names on *MEASURE*.
- If you do not enter a *MEASURE* subcommand and there are more dependent variables than cells in the within-subjects design, MANOVA assigns names (normally *MEAS.1*, *MEAS.2*, and so on) to the different measures.
- All of the dependent variables corresponding to each measure should be listed together and ordered so that the within-subjects factor named last on the *WSFACTORS* subcommand varies most rapidly.

**Example**

```
MANOVA TEMP1 TO TEMP6, WEIGHT1 TO WEIGHT6 BY GROUP(1,2)
  /WSFACTORS=DAY(3) AMPM(2)
  /MEASURE=TEMP WEIGHT
  /WSDSIGN=DAY, AMPM, DAY BY AMPM
  /PRINT=SIGNIF(HYPOTH AVERF)
  /DESIGN.
```

- There are 12 dependent variables: six temperatures and six weights, corresponding to morning and afternoon measurements on three days.
- `WSFACTORS` identifies the two factors (*DAY* and *AMPM*) that distinguish the temperature and weight measurements for each subject. These factors define six within-subjects cells.
- `MEASURE` indicates that the first group of six dependent variables correspond to *TEMP* and the second group of six dependent variables correspond to *WEIGHT*.
- These labels, *TEMP* and *WEIGHT*, are used on the output requested by `PRINT`.
- `WSDSIGN` requests a full factorial within-subjects model. Because this is the default, `WSDSIGN` could have been omitted.

**RENAME Subcommand**

Because any repeated measures analysis involves a transformation of the dependent variables, it is always a good idea to rename the dependent variables. Choose appropriate names depending on the type of contrast specified for within-subjects factors. This is easier to do if you are using one of the orthogonal contrasts. The most reliable way to assign new names is to inspect the transformation matrix.

**Example**

```
MANOVA LOW1 LOW2 LOW3 HI1 HI2 HI3
  /WSFACTORS=LEVEL(2) TRIAL(3)
  /CONTRAST(TRIAL)=DIFFERENCE
  /RENAME=CONST LEVELDIF TRIAL21 TRIAL312 INTER1 INTER2
  /PRINT=TRANSFORM
  /DESIGN.
```

- This analysis has two within-subjects factors and no between-subjects factors.
- Difference contrasts are requested for *TRIAL*, which has three levels.
- Because all orthonormal contrasts produce the same *F* test for a factor with two levels, there is no point in specifying a contrast type for *LEVEL*.
- New names are assigned to the transformed variables based on the transformation matrix. These names correspond to the meaning of the transformed variables: the mean or constant, the average difference between levels, the average effect of trial 2 compared to 1, the average effect of trial 3 compared to 1 and 2; and the two interactions between *LEVEL* and *TRIAL*.
- The transformation matrix requested by the `PRINT` subcommand looks like the following table.

Table 131-1  
Transformation matrix

	CONST	LEVELDIF	TRIAL1	TRIAL2	INTER1	INTER2
LOW1	0.408	0.408	-0.500	-0.289	-0.500	-0.289

	CONST	LEVELDIF	TRIAL1	TRIAL2	INTER1	INTER2
LOW2	0.408	0.408	0.500	-0.289	0.500	-0.289
LOW3	0.408	0.408	0.000	0.577	0.000	0.577
HI1	0.408	-0.408	-0.500	-0.289	0.500	0.289
HI2	0.408	-0.408	0.500	-0.289	-0.500	0.289
HI3	0.408	-0.408	0.000	0.577	0.000	-0.577

## PRINT Subcommand

The following additional specifications on PRINT are useful in repeated measures analysis:

<b>SIGNIF(AVERF)</b>	<i>Averaged F tests for use with repeated measures.</i> This is the default display in repeated measures analysis. The averaged <i>F</i> tests in the multivariate setup for repeated measures are equivalent to the univariate (or split-plot or mixed-model) approach to repeated measures.
<b>SIGNIF(AVONLY)</b>	<i>Only the averaged F test for repeated measures.</i> AVONLY produces the same output as AVERF and suppresses all other SIGNIF output.
<b>SIGNIF(HF)</b>	<i>The Huynh-Feldt corrected significance values for averaged univariate F tests.</i>
<b>SIGNIF(GG)</b>	<i>The Greenhouse-Geisser corrected significance values for averaged univariate F tests.</i>
<b>SIGNIF(EFSIZE)</b>	<i>The effect size for the univariate F and t tests.</i>

- The keywords AVERF and AVONLY are mutually exclusive.
- When you request repeated measures analysis with the WSFACTORS subcommand, the default display includes SIGNIF(AVERF) but does not include the usual SIGNIF(UNIV).
- The averaged *F* tests are appropriate in repeated measures because the dependent variables that are averaged actually represent contrasts of the WSFACTOR variables. When the analysis is not doubly multivariate, as discussed above, you can specify PRINT=SIGNIF(UNIV) to obtain significance tests for each degree of freedom, just as in univariate MANOVA.

## References

- Burns, P. R. 1984. Multiple comparison methods in MANOVA. In: *Proceedings of the 7th SPSS Users and Coordinators Conference*, .
- Green, P. E. 1978. *Analyzing multivariate data*. Hinsdale, Ill.: The Dryden Press.
- Huberty, C. J. 1972. Multivariate indices of strength of association. *Multivariate Behavioral Research*, 7, 516–523.
- Muller, K. E., and B. L. Peterson. 1984. Practical methods for computing power in testing the multivariate general linear hypothesis. *Computational Statistics and Data Analysis*, 2, 143–158.
- Pillai, K. C. S. 1967. Upper percentage points of the largest root of a matrix in multivariate analysis. *Biometrika*, 54, 189–194.

# MATCH FILES

```
MATCH FILES FILE={'savfile'|'dataset'} [TABLE={'savfile'|'dataset'}]
              {*}                }          {*}                }

[/RENAME=(old varnames=new varnames)...]

[/IN=varname]

/FILE=... [TABLE= ...]

[/BY varlist]

[/MAP]

[/KEEP={ALL** }] [/DROP=varlist]
              {varlist}

[/FIRST=varname] [/LAST=varname]
```

\*\*Default if the subcommand is omitted.

## Example

```
MATCH FILES FILE='/data/part1.sav'
           /FILE='/data/part2.sav'
           /FILE=*
```

## Overview

`MATCH FILES` combines variables from 2 up to 50 IBM® SPSS® Statistics data files. `MATCH FILES` can make parallel or nonparallel matches between different files or perform table lookups. **Parallel matches** combine files sequentially by case (they are sometimes referred to as **sequential matches**). **Nonparallel matches** combine files according to the values of one or more key variables. In a table lookup, `MATCH FILES` looks up variables in one file and transfers those variables to a case file.

The files specified on `MATCH FILES` can be SPSS Statistics data files or open datasets in the current session. The combined file becomes the new active dataset.

In general, `MATCH FILES` is used to combine files containing the same cases but different variables. To combine files containing the same variables but different cases, use `ADD FILES`. To update existing SPSS Statistics data files, use `UPDATE`.

## Options

**Variable Selection.** You can specify which variables from each input file are included in the new active dataset using the `DROP` and `KEEP` subcommands.

**Variable Names.** You can rename variables in each input file before combining the files using the `RENAME` subcommand. This permits you to combine variables that are the same but whose names differ in different input files or to separate variables that are different but have the same name.

**Variable Flag.** You can create a variable that indicates whether a case came from a particular input file using `IN`. You can use the `FIRST` or `LAST` subcommands to create a variable that flags the first or last case of a group of cases with the same value for the key variable.

**Variable Map.** You can request a map showing all variables in the new active dataset, their order, and the input files from which they came using the `MAP` subcommand.

### ***Basic Specification***

The basic specification is two or more `FILE` subcommands, each of which specifies a file to be matched. In addition, `BY` is required to specify the key variables for nonparallel matches. Both `BY` and `TABLE` are required to match table-lookup files.

- All variables from all input files are included in the new active dataset unless `DROP` or `KEEP` is specified.

### ***Subcommand Order***

- `RENAME` and `IN` must immediately follow the `FILE` or `TABLE` subcommand to which they apply.
- Any `BY`, `FIRST`, `LAST`, `KEEP`, `DROP`, and `MAP` subcommands must follow all of the `TABLE`, `FILE`, `RENAME`, and `IN` subcommands.

### ***Syntax Rules***

- `RENAME` can be repeated after each `FILE` or `TABLE` subcommand and applies only to variables in the file named on the immediately preceding `FILE` or `TABLE`.
- `IN` can be used only for a nonparallel match or for a table lookup. (Thus, `IN` can be used only if `BY` is specified.)
- `BY` can be specified only once. However, multiple variables can be specified on `BY`. When `BY` is used, all files must be sorted in ascending order of the key variables named on `BY`.
- `MAP` can be repeated as often as desired.

### ***Operations***

- `MATCH FILES` reads all files named on `FILE` or `TABLE` and builds a new active dataset.
- If the current active dataset is included and is specified with an asterisk (`FILE=*`), the new merged dataset replaces the active dataset. If that dataset is a named dataset, the merged dataset retains that name. If the current active dataset is not included or is specified by name (for example, `FILE=Dataset1`), a new unnamed, merged dataset is created, and it becomes the active dataset. For information on naming datasets, see [DATASET NAME](#).
- The new active dataset contains complete dictionary information from the input files, including variable names, labels, print and write formats, and missing-value indicators. The new file also contains the documents from each of the input files. See `DROP DOCUMENTS` for information on deleting documents.
- Variables are copied in order from the first file specified, then from the second file specified, and so on.

- If the same variable name is used in more than one input file, data are taken from the file specified first. Dictionary information is taken from the first file containing value labels, missing values, or a variable label for the common variable. If the first file has no such information, `MATCH FILES` checks the second file, and so on, seeking dictionary information.
- All cases from all input files are included in the combined file. Cases that are absent from one of the input files will be assigned system-missing values for variables unique to that file.
- `BY` specifies that cases should be combined according to a common value on one or more key variables. All input files must be sorted in ascending order of the key variables.
- If `BY` is not used, the program performs a parallel (sequential) match, combining the first case from each file, then the second case from each file, and so on, without regard to any identifying values that may be present.
- If the active dataset is named as an input file, any `N` and `SAMPLE` commands that have been specified are applied to that file before files are matched.

### **Limitations**

- Maximum 50 files can be combined on one `MATCH FILES` command.
- Maximum one `BY` subcommand. However, `BY` can specify multiple variables.
- The `TEMPORARY` command cannot be in effect if the active dataset is used as an input file.

### **Example**

```
MATCH FILES FILE='/data/part1.sav'
           /FILE='/data/part2.sav'
           /FILE=*
```

- `MATCH FILES` combines three files (the active dataset and two external SPSS Statistics data files) in a parallel match. Cases are combined according to their order in each file.
- The new active dataset contains as many cases as are contained in the largest of the three input files.

### **Example**

```
GET FILE='/examples/data/mydata.sav'.
SORT CASES BY ID.
DATASET NAME mydata.
GET DATA /TYPE=XLS
           /FILE='/examples/data/excelfile.xls'.
SORT CASES BY ID.
DATASET NAME excelfile.
GET DATA /TYPE=ODBC /CONNECT=
           'DSN=MS Access Database;DBQ=/examples/data/dm_demo.mdb;'+
           'DriverId=25;FIL=MS Access;MaxBufferSize=2048;PageTimeout=5;'
           /SQL='SELECT * FROM main'.
SORT CASES BY ID.
MATCH FILES
           /FILE='mydata'
           /FILE='excelfile'
           /FILE=*
           /BY ID.
```

- A data file in SPSS Statistics format and assigned the dataset name *mydata*. Since it has been assigned a dataset name, it remains available for subsequent use even after other data sources have been opened.
- An Excel file is then read and assigned the dataset name *exceldata*. Like the SPSS Statistics data file, since it has been assigned a dataset name, it remains available after other data sources have been opened.
- Then a table from a database is read. Since it is the most recently opened or activated dataset, it is the active dataset.
- The three datasets are then merged together with `MATCH FILES` command, using the dataset names on the `FILE` subcommands instead of file names.
- An asterisk (\*) is used to specify the active dataset, which is the database table in this example.
- The files are merged together based on the value of the key variable *ID*, specified on the `BY` subcommand.
- Since all the files being merged need to be sorted in the same order of the key variable(s), `SORT CASES` is performed on each dataset.

## FILE Subcommand

`FILE` identifies the files to be combined (except table files). At least one `FILE` subcommand is required on `MATCH FILES`. A separate `FILE` subcommand must be used for each input file.

- An asterisk can be specified on `FILE` to refer to the active dataset.
- Dataset names instead of filenames can be used to refer to currently open datasets.
- The order in which files are specified determines the order of variables in the new active dataset. In addition, if the same variable name occurs in more than one input file, the variable is taken from the file specified first.
- If the files have unequal numbers of cases, cases are generated from the longest file. Cases that do not exist in the shorter files have system-missing values for variables that are unique to those files.

## Text Data Files

You can add variables from one or more text data files by reading the files into IBM® SPSS® Statistics (with `DATA LIST` or `GET DATA`), defining dataset names for each file (`DATASET NAME` command), and then using `MATCH FILES` to add the cases from each file.

### Example

```
DATA LIST FILE="/data/textdata1.txt"
  /id 1-3 var1 5-7 var2 9-12.
SORT CASES by ID.
DATASET NAME file1.
DATA LIST FILE="/data/textdata2.txt"
  /id 1-3 var3 5-9 var4 11-15.
SORT CASES BY ID.
DATASET NAME file2.
DATA LIST FILE="/data/textdata3.txt"
  /id 1-3 var5 5-6 var6 8-10.
```



```
DATASET NAME file3.  
MATCH FILES FILE='file1'  
  /FILE='file2'  
  /FILE='file3'  
  /BY id.  
SAVE OUTFILE='/data/combined_data.sav'.
```

## ***BY Subcommand***

BY specifies one or more identification, or key, variables that determine which cases are to be combined. When BY is specified, cases from one file are matched only with cases from other files that have the same values for the key variables. BY is required unless all input files are to be matched sequentially according to the order of cases.

- BY must follow the FILE and TABLE subcommands and any associated RENAME and IN subcommands.
- BY specifies the names of one or more key variables. The key variables must exist in all input files. The key variables can be numeric or long or short strings.
- All input files must be sorted in ascending order of the key variables. If necessary, use SORT CASES before MATCH FILES.
- Missing values for key variables are handled like any other values.
- Unmatched cases are assigned system-missing values (for numeric variables) or blanks (for string variables) for variables from files that do not contain a match.

## ***Duplicate Cases***

Duplicate cases are those with the same values for the key variables named on the BY subcommand.

- Duplicate cases are permitted in any input files except table files.
- When there is no table file, the first duplicate case in each file is matched with the first matching case (if any) from the other files; the second duplicate case is matched with a second matching duplicate, if any; and so on. (In effect, a parallel match is performed within groups of duplicate cases.) Unmatched cases are assigned system-missing values (for numeric variables) or blanks (for string variables) for variables from files that do not contain a match.
- The program displays a warning if it encounters duplicate keys in one or more of the files being matched.

## ***TABLE Subcommand***

TABLE specifies a table lookup (or keyed table) file. A lookup file contributes variables but not cases to the new active dataset. Variables from the table file are added to all cases from other files that have matching values for the key variables. FILE specifies the files that supply the cases.

- A separate TABLE subcommand must be used to specify each lookup file, and a separate FILE subcommand must be used to specify each case file.
- The BY subcommand is required when TABLE is used.

- All specified files must be sorted in ascending order of the key variables. If necessary, use `SORT CASES` before `MATCH FILES`.
- A lookup file cannot contain duplicate cases (cases for which the key variable[s] named on `BY` have identical values).
- An asterisk on `TABLE` refers to the active dataset.
- Dataset names instead of file names can be used to refer to currently open datasets.
- Cases in a case file that do not have matches in a table file are assigned system-missing values (for numeric variables) or blanks (for string variables) for variables from that table file.
- Cases in a table file that do not match any cases in a case file are ignored.

**Example**

```
MATCH FILES FILE=*
  /TABLE='/data/master.sav'
  /BY EMP_ID.
```

- `MATCH FILES` combines variables from the IBM® SPSS® Statistics data file *master.sav* with the active dataset, matching cases by the variable *EMP\_ID*.
- No new cases are added to the active dataset as a result of the table lookup.
- Cases whose value for *EMP\_ID* is not included in the *master.sav* file are assigned system-missing values for variables taken from the table.

**RENAME Subcommand**

`RENAME` renames variables on the input files *before* they are processed by `MATCH FILES`. `RENAME` must follow the `FILE` or `TABLE` subcommand that contains the variables to be renamed.

- `RENAME` applies only to the immediately preceding `FILE` or `TABLE` subcommand. To rename variables from more than one input file, specify a `RENAME` subcommand after each `FILE` or `TABLE` subcommand.
- Specifications for `RENAME` consist of a left parenthesis, a list of old variable names, an equals sign, a list of new variable names, and a right parenthesis. The two variable lists must name or imply the same number of variables. If only one variable is renamed, the parentheses are optional.
- More than one rename specification can be specified on a single `RENAME` subcommand, each enclosed in parentheses.
- The `TO` keyword can be used to refer to consecutive variables in the file and to generate new variable names.
- `RENAME` takes effect immediately. Any `KEEP` and `DROP` subcommands entered prior to a `RENAME` must use the old names, while `KEEP` and `DROP` subcommands entered after a `RENAME` must use the new names.
- All specifications within a single set of parentheses take effect simultaneously. For example, the specification `RENAME (A,B = B,A)` swaps the names of the two variables.
- Variables cannot be renamed to scratch variables.

**Example**

```
MATCH FILES FILE='/data/update.sav'  
/RENAME=(NEWID = ID)  
/FILE='/data/master.sav'  
/BY ID.
```

- MATCH FILES matches a master IBM® SPSS® Statistics data file (*master.sav*) with an update data file (*update.sav*).
- The variable *NEWID* in the *update.sav* file is renamed *ID* so that it will have the same name as the identification variable in the master file and can be used on the BY subcommand.

## **DROP and KEEP Subcommands**

DROP and KEEP are used to include a subset of variables in the new active dataset. DROP specifies a set of variables to exclude and KEEP specifies a set of variables to retain.

- DROP and KEEP do not affect the input files on disk.
- DROP and KEEP must follow all FILE, TABLE, and RENAME subcommands.
- DROP and KEEP must specify one or more variables. If RENAME is used to rename variables, specify the new names on DROP and KEEP.
- The keyword ALL can be specified on KEEP. ALL must be the last specification on KEEP, and it refers to all variables not previously named on KEEP.
- DROP cannot be used with variables created by the IN, FIRST, or LAST subcommands.
- KEEP can be used to change the order of variables in the resulting file. By default, MATCH FILES first copies the variables in order from the first file, then copies the variables in order from the second file, and so on. With KEEP, variables are kept in the order in which they are listed on the subcommand. If a variable is named more than once on KEEP, only the first mention of the variable is in effect; all subsequent references to that variable name are ignored.

**Example**

```
MATCH FILES FILE='/data/particle.sav'  
/RENAME=(PARTIC=POLLUTE1)  
/FILE='/data/gas.sav'  
/RENAME=(OZONE TO SULFUR=POLLUTE2 TO POLLUTE4)  
/DROP=POLLUTE4.
```

- The renamed variable *POLLUTE4* is dropped from the resulting file. DROP is specified after all of the FILE and RENAME subcommands, and it refers to the dropped variable by its new name.

## **IN Subcommand**

IN creates a new variable in the resulting file that indicates whether a case came from the input file named on the preceding FILE subcommand. IN applies only to the file specified on the immediately preceding FILE subcommand.

- IN can be used only for a nonparallel match or table lookup.
- IN has only one specification—the name of the flag variable.

- The variable created by `IN` has the value 1 for every case that came from the associated input file and the value 0 if the case came from a different input file.
- Variables created by `IN` are automatically attached to the end of the resulting file and cannot be dropped. If `FIRST` or `LAST` is used, the variable created by `IN` precedes the variables created by `FIRST` or `LAST`.

**Example**

```
MATCH FILES FILE='/data/week10.sav'
           /FILE='/data/week11.sav'
           /IN=INWEEK11
           /BY=EMPID.
```

- `IN` creates the variable `INWEEK11`, which has the value 1 for all cases in the resulting file that had values in the input file `week11.sav` and the value 0 for those cases that were not in file `week11.sav`.

**FIRST and LAST Subcommands**

`FIRST` and `LAST` create logical variables that flag the first or last case of a group of cases with the same value for the `BY` variables.

- `FIRST` and `LAST` must follow all `TABLE` and `FILE` subcommands and any associated `RENAME` and `IN` subcommands.
- `FIRST` and `LAST` have only one specification—the name of the flag variable.
- `FIRST` creates a variable with the value 1 for the first case of each group and the value 0 for all other cases.
- `LAST` creates a variable with the value 1 for the last case of each group and the value 0 for all other cases.
- Variables created by `FIRST` and `LAST` are automatically attached to the end of the resulting file and cannot be dropped.
- If one file has several cases with the same values for the key variables, `FIRST` or `LAST` can be used to create a variable that flags the first or last case of the group.

**Example**

```
MATCH FILES TABLE='/data/house.sav'
           /FILE='/data/persons.sav'
           /BY=HOUSEID /FIRST=HEAD.
```

- The variable `HEAD` contains the value 1 for the first person in each household and the value 0 for all other persons. Assuming that the `persons.sav` file is sorted with the head of household as the first case for each household, the variable `HEAD` identifies the case for the head of household.

**Example**

- \* Using match files with only one file.
- \* This example flags the first of several cases with

the same value for a key variable.

```
MATCH FILES FILE='/data/persons.sav'  
  /BY HOUSEID /FIRST=HEAD.  
SELECT IF (HEAD EQ 1).  
CROSSTABS JOBCAT BY SEX.
```

- `MATCH FILES` is used instead of `GET` to read the data file *persons.sav*. The `BY` subcommand identifies the key variable (*HOUSEID*), and `FIRST` creates the variable *HEAD* with the value 1 for the first case in each household and the value 0 for all other cases.
- `SELECT IF` selects only the cases with the value 1 for *HEAD*, and the `CROSSTABS` procedure is run on these cases.

## ***MAP Subcommand***

`MAP` produces a list of the variables that are in the new active dataset and the file or files from which they came. Variables are listed in the order in which they appear in the resulting file. `MAP` has no specifications and must be placed after all `FILE`, `TABLE`, and `RENAME` subcommands.

- Multiple `MAP` subcommands can be used. Each `MAP` shows the current status of the active dataset and reflects only the subcommands that precede the `MAP` subcommand.
- To obtain a map of the resulting file in its final state, specify `MAP` last.
- If a variable is renamed, its original and new names are listed. Variables created by `IN`, `FIRST`, and `LAST` are not included in the map, since they are automatically attached to the end of the file and cannot be dropped.

# MATRIX-END MATRIX

```
MATRIX
matrix statements
END MATRIX
```

*The following matrix language statements can be used in a matrix program:*

BREAK	DO IF	END LOOP	MSAVE	SAVE
CALL	ELSE	GET	PRINT	WRITE
COMPUTE	ELSE IF	LOOP	READ	
DISPLAY	END IF	MGET	RELEASE	

*The following functions can be used in matrix language statements:*

ABS	Absolute values of matrix elements
ALL	Test if all elements are positive
ANY	Test if any element is positive
ARSIN	Arcsines of matrix elements
ARTAN	Arctangents of matrix elements
BLOCK	Create block diagonal matrix
CDFNORM	Cumulative normal distribution function
CHICDF	Cumulative chi-squared distribution function
CHOL	Cholesky decomposition
CMAX	Column maxima
CMIN	Column minima
COS	Cosines of matrix elements
CSSQ	Column sums of squares
CSUM	Column sums
DESIGN	Create design matrix
DET	Determinant
DIAG	Diagonal of matrix
EOF	Check end of file
EVAL	Eigenvalues of symmetric matrix
EXP	Exponentials of matrix elements
FCDF	Cumulative $F$ distribution function
GINV	Generalized inverse
GRADE	Rank elements in matrix, using sequential integers for ties
GSCH	Gram-Schmidt orthonormal basis
IDENT	Create identity matrix
INV	Inverse
KRONECKER	Kronecker product of two matrices
LG10	Logarithms to base 10 of matrix elements
LN	Logarithms to base $e$ of matrix elements

MAGIC	Create magic square
MAKE	Create a matrix with all elements equal
MDIAG	Create a matrix with the given diagonal
MMAX	Maximum element in matrix
MMIN	Minimum element in matrix
MOD	Remainders after division
MSSQ	Matrix sum of squares
MSUM	Matrix sum
NCOL	Number of columns
NROW	Number of rows
RANK	Matrix rank
RESHAPE	Change shape of matrix
RMAX	Row maxima
RMIN	Row minima
RND	Round off matrix elements to nearest integer
RNKORDER	Rank elements in matrix, averaging ties
RSSQ	Row sums of squares
RSUM	Row sums
SIN	Sines of matrix elements
SOLVE	Solve systems of linear equations
SQRT	Square roots of matrix elements
SSCP	Sums of squares and cross-products
SVAL	Singular values
SWEEP	Perform sweep transformation
T	Synonym for TRANSPOS
TCDF	Cumulative normal $t$ distribution function
TRACE	Calculate trace (sum of diagonal elements)
TRANSPOS	Transposition of matrix
TRUNC	Truncation of matrix elements to integer
UNIFORM	Create matrix of uniform random numbers

### **Example**

```

MATRIX.
READ A /FILE=MATRDATA /SIZE={6,6} /FIELD=1 TO 60.
CALL EIGEN(A, EIGENVEC, EIGENVAL) .
LOOP J=1 TO NROW(EIGENVAL) .
+ DO IF (EIGENVAL(J) > 1.0) .
+   PRINT EIGENVAL(J) / TITLE="Eigenvalue:" /SPACE=3.
+   PRINT T(EIGENVEC(:,J)) / TITLE="Eigenvector:" /SPACE=1.
+ END IF.
END LOOP.
END MATRIX.

```

## Overview

The `MATRIX` and `END MATRIX` commands enclose statements that are executed by the matrix processor. Using matrix programs, you can write your own statistical routines in the compact language of matrix algebra. Matrix programs can include mathematical calculations, control structures, display of results, and reading and writing matrices as character files or data files.

As discussed below, a matrix program is for the most part independent of the rest of the session, although it can read and write IBM® SPSS® Statistics data files, including the active dataset.

This section does not attempt to explain the rules of matrix algebra. Many textbooks teach the application of matrix methods to statistics.

The `MATRIX` procedure was originally developed at the Madison Academic Computing Center, University of Wisconsin.

## Terminology

A variable within a matrix program represents a **matrix**, which is simply a set of values arranged in a rectangular array of rows and columns.

- An  $n \times m$  (read “n by m”) matrix is one that has  $n$  rows and  $m$  columns. The integers  $n$  and  $m$  are the dimensions of the matrix. An  $n \times m$  matrix contains  $n \times m$  elements, or data values.
- An  $n \times 1$  matrix is sometimes called a **column vector**, and a  $1 \times n$  matrix is sometimes called a **row vector**. A vector is a special case of a matrix.
- A  $1 \times 1$  matrix, containing a single data value, is often called a **scalar**. A scalar is also a special case of a matrix.
- An **index** to a matrix or vector is an integer that identifies a specific row or column. Indexes normally appear in printed works as subscripts, as in  $A_{31}$ , but are specified in the matrix language within parentheses, as in  $A(3,1)$ . The row index for a matrix precedes the column index.
- The **main diagonal** of a matrix consists of the elements whose row index equals their column index. It begins at the top left corner of the matrix; in a square matrix, it runs to the bottom right corner.
- The **transpose** of a matrix is the matrix with rows and columns interchanged. The transpose of an  $n \times m$  matrix is an  $m \times n$  matrix.
- A **symmetric matrix** is a square matrix that is unchanged if you flip it about the main diagonal. That is, the element in row  $i$ , column  $j$  equals the element in row  $j$ , column  $i$ . A symmetric matrix equals its transpose.
- Matrices are always rectangular, although it is possible to read or write symmetric matrices in triangular form. Vectors and scalars are considered degenerate rectangles.
- It is an error to try to create a matrix whose rows have different numbers of elements.

A matrix program does not process individual cases unless you so specify, using the control structures of the matrix language. Unlike ordinary IBM® SPSS® Statistics variables, matrix variables do not have distinct values for different cases. A matrix is a single entity.



Vectors in matrix processing should not be confused with the vectors temporarily created by the VECTOR command. The latter are shorthand for a list of SPSS Statistics variables and, like all ordinary SPSS Statistics variables, are unavailable during matrix processing.

## Matrix Variables

A matrix variable is created by a matrix statement that assigns a value to a variable name.

- A matrix variable name follows the same rules as those applicable to an ordinary IBM® SPSS® Statistics variable name, except matrix variable names cannot exceed 8 bytes.
- The names of matrix functions and procedures cannot be used as variable names within a matrix program. (In particular, the letter *T* cannot be used as a variable name because *T* is an alias for the TRANSPOS function.)
- The COMPUTE, READ, GET, MGET, and CALL statements create matrices. An index variable named on a LOOP statement creates a scalar with a value assigned to it.
- A variable name can be redefined within a matrix program without regard to the dimensions of the matrix it represents. The same name can represent scalars, vectors, and full matrices at different points in the matrix program.
- MATRIX-END MATRIX does not include any special processing for missing data. When reading a data matrix from a data file, you must therefore specify whether missing data are to be accepted as valid or excluded from the matrix.

## String Variables in Matrix Programs

Matrix variables can contain short string data. Support for string variables is limited, however.

- MATRIX will attempt to carry out calculations with string variables if you so request. The results will not be meaningful.
- You must specify a format (such as A8) when you display a matrix that contains string data.

## Syntax of Matrix Language

A matrix program consists of statements. Matrix statements must appear in a matrix program, between the MATRIX and END MATRIX commands. They are analogous to IBM® SPSS® Statistics commands and follow the rules of the command language regarding the abbreviation of keywords; the equivalence of upper and lower case; the use of spaces, commas, and equals signs; and the splitting of statements across multiple lines. However, commas are required to separate arguments to matrix functions and procedures and to separate variable names on the RELEASE statement.

Matrix statements are composed of the following elements:

- Keywords, such as the names of matrix statements
- Variable names
- Explicitly written matrices, which are enclosed within braces ({})
- Arithmetic and logical operators

- Matrix functions
- The command terminator, which serves as a statement terminator within a matrix program

### ***Comments in Matrix Programs***

Within a matrix program, you can enter comments in any of the following forms: on lines beginning with the `COMMENT` command, on lines beginning with an asterisk, or between the characters `/*` and `*/` on a command line.

### ***Matrix Notation***

To write a matrix explicitly:

- Enclose the matrix within braces (`{}`).
- Separate the elements of each row by commas.
- Separate the rows by semicolons.
- String elements must be enclosed in either quotes, as is generally true in the command language.

#### ***Example***

```
{1,2,3;4,5,6}
```

- The example represents the following matrix:

$$\begin{bmatrix} 1 & 2 & 3 \\ 4 & 5 & 6 \end{bmatrix}$$

#### ***Example***

```
{1,2,3}
```

- This example represents a row vector:

$$[1 \ 2 \ 3]$$

#### ***Example***

```
{11;12;13}
```

- This example represents a column vector:

$$\begin{bmatrix} 11 \\ 12 \\ 13 \end{bmatrix}$$

#### ***Example***

```
{3}
```

- This example represents a scalar. The braces are optional. You can specify the same scalar as 3.

## Matrix Notation Shorthand

You can simplify the construction of matrices using notation shorthand.

**Consecutive Integers.** Use a colon to indicate a range of consecutive integers. For example, the vector  $\{1, 2, 3, 4, 5, 6\}$  can be written as  $\{1:6\}$ .

**Incremented Ranges of Integers.** Use a second colon followed by an integer to indicate the increment. The matrix  $\{1, 3, 5, 7; 2, 5, 8, 11\}$  can be written as  $\{1:7:2; 2:11:3\}$ , where  $1:7:2$  indicates the integers from 1 to 7 incrementing by 2, and  $2:11:3$  indicates the integers from 2 to 11 incrementing by 3.

- You must use integers when specifying a range in either of these ways. Numbers with fractional parts are truncated to integers.
- If an arithmetic expression is used, it should be enclosed in parentheses.

## Extraction of an Element, a Vector, or a Submatrix

You can use indexes in parentheses to extract an element from a vector or matrix, a vector from a matrix, or a submatrix from a matrix. In the following discussion, an **integer index** refers to an integer expression used as an index, which can be a scalar matrix with an integer value or an integer element extracted from a vector or matrix. Similarly, a **vector index** refers to a vector expression used as an index, which can be a vector matrix or a vector extracted from a matrix.

For example, if  $S$  is a scalar matrix,  $S = [2]$ ,  $R$  is a row vector,  $R = [1 \ 3 \ 5]$ ,  $C$  is a column

vector,  $C = \begin{bmatrix} 2 \\ 3 \\ 4 \end{bmatrix}$ , and  $A$  is a  $5 \times 5$  matrix,  $A = \begin{bmatrix} 11 & 12 & 13 & 14 & 15 \\ 21 & 22 & 23 & 24 & 25 \\ 31 & 32 & 33 & 34 & 35 \\ 41 & 42 & 43 & 44 & 45 \\ 51 & 52 & 53 & 54 & 55 \end{bmatrix}$ , then:

$$R(S) = R(2) = \{3\}$$

$$C(S) = C(2) = \{3\}$$

- An integer index extracts an element from a vector matrix.
- The distinction between a row and a column vector does not matter when an integer index is used to extract an element from it.

$$A(2,3) = A(S,3) = \{23\}$$

- Two integer indexes separated by a comma extract an element from a rectangular matrix.

$$A(R,2) = A(1:5:2,2) = \{12; 32; 52\}$$

$$A(2,R) = A(2,1:5:2) = \{21, 23, 25\}$$

$$A(C,2) = A(2:4,2) = \{22; 32; 42\}$$

$$A(2,C) = A(2,2:4) = \{22, 23, 24\}$$

- An integer and a vector index separated by a comma extract a vector from a matrix.
- The distinction between a row and a column vector does not matter when used as indexes in this way.

$$A(2,:) = A(S,:) = \{21, 22, 23, 24, 25\}$$

$$A(:,2) = A(:,S) = \{12; 22; 32; 42; 52\}$$

- A colon by itself used as an index extracts an entire row or column vector from a matrix.

$$A(R,C) = A(R,2:4) = A(1:5:2,C) = A(1:5:2,2:4) = \{12,13,14;32,33,34;52,53,54\}$$

$$A(C,R) = A(C,1:5:2) = A(2:4,R) = A(2:4,1:5:2) = \{21,23,25;31,33,35;41,43,45\}$$

- Two vector indexes separated by a comma extract a submatrix from a matrix.
- The distinction between a row and a column vector does not matter when used as indexes in this way.

### Construction of a Matrix from Other Matrices

You can use vector or rectangular matrices to construct a new matrix, separating row expressions by semicolons and components of row expressions by commas. If a column vector  $V_c$  has  $n$  elements and matrix  $M$  has the dimensions  $n \times m$ , then  $\{M; V_c\}$  is an  $n \times (m + 1)$  matrix. Similarly, if the row vector  $V_r$  has  $m$  elements and  $M$  is the same, then  $\{M; V_r\}$  is an  $(n + 1) \times m$  matrix. In fact, you can paste together any number of matrices and vectors this way.

- All of the components of each column expression must have the same number of actual rows, and all of the row expressions must have the same number of actual columns.
- The distinction between row vectors and column vectors must be observed carefully when constructing matrices in this way, so that the components will fit together properly.
- Several of the matrix functions are also useful in constructing matrices; see in particular the `MAKE`, `UNIFORM`, and `IDENT` functions in [Matrix Functions on p. 1141](#).

#### Example

```
COMPUTE M={CORNER, COL3; ROW3}.
```

- This example constructs the matrix  $M$  from the matrix  $CORNER$ , the column vector  $COL3$ , and the row vector  $ROW3$ .
- $COL3$  supplies new row components and is separated from  $CORNER$  by a comma.
- $ROW3$  supplies column elements and is separated from previous expressions by a semicolon.
- $COL3$  must have the same number of rows as  $CORNER$ .
- $ROW3$  must have the same number of columns as the matrix resulting from the previous expressions.
- For example, if  $CORNER = \begin{bmatrix} 11 & 12 \\ 21 & 22 \end{bmatrix}$ ,  $COL3 = \begin{bmatrix} 13 \\ 23 \end{bmatrix}$ , and  $ROW3 = [31 \ 32 \ 33]$ , then:

$$M = \begin{bmatrix} 11 & 12 & 13 \\ 21 & 22 & 23 \\ 31 & 32 & 33 \end{bmatrix}$$

### Matrix Operations

You can perform matrix calculations according to the rules of matrix algebra and compare matrices using relational or logical operators.

## Conformable Matrices

Many operations with matrices make sense only if the matrices involved have “suitable” dimensions. Most often, this means that they should be the same size, with the same number of rows and the same number of columns. Matrices that are the right size for an operation are said to be **conformable matrices**. If you attempt to do something in a matrix program with a matrix that is not conformable for that operation—a matrix that has the wrong dimensions—you will receive an error message, and the operation will not be performed. An important exception, where one of the matrices is a scalar, is discussed below.

Requirements for carrying out matrix operations include:

- Matrix addition and subtraction require that the two matrices be the same size.
- The relational and logical operations described below require that the two matrices be the same size.
- Matrix multiplication requires that the number of columns of the first matrix equal the number of rows of the second matrix.
- Raising a matrix to a power can be done only if the matrix is square. This includes the important operation of *inverting* a matrix, where the power is  $-1$ .
- Conformability requirements for matrix functions are noted in [Matrix Functions on p. 1141](#) and in [COMPUTE Statement on p. 1140](#).

## Scalar Expansion

When one of the matrices involved in an operation is a scalar, the scalar is treated as a matrix of the correct size in order to carry out the operation. This internal scalar expansion is performed for the following operations:

- Addition and subtraction.
- Elementwise multiplication, division, and exponentiation. Note that multiplying a matrix elementwise by an expanded scalar is equivalent to ordinary scalar multiplication—each element of the matrix is multiplied by the scalar.
- All relational and logical operators.

## Arithmetic Operators

You can add, subtract, multiply, or exponentiate matrices according to the rules of matrix algebra, or you can perform elementwise arithmetic, in which you multiply, divide, or exponentiate each element of a matrix separately. The arithmetic operators are listed below.

Unary -	<i>Sign reversal.</i> A minus sign placed in front of a matrix reverses the sign of each element. (The unary + is also accepted but has no effect.)
+	<i>Matrix addition.</i> Corresponding elements of the two matrices are added. The matrices must have the same dimensions, or one must be a scalar.
-	<i>Matrix subtraction.</i> Corresponding elements of the two matrices are subtracted. The matrices must have the same dimensions, or one must be a scalar.

*	<i>Multiplication.</i> There are two cases. First, <i>scalar multiplication</i> : if either of the matrices is a scalar, each element of the other matrix is multiplied by that scalar. Second, <i>matrix multiplication</i> : if $A$ is an $m \times n$ matrix and $B$ is an $n \times p$ matrix, $A*B$ is an $m \times p$ matrix in which the element in row $i$ , column $k$ , is equal to $\sum_{j=1}^n A(i, j) \times B(j, k)$
/	<i>Division.</i> The division operator performs elementwise division (described below). True matrix division, the inverse operation of matrix multiplication, is accomplished by taking the <code>INV</code> function (square matrices) or the <code>GINV</code> function (rectangular matrices) of the denominator and multiplying.
**	<i>Matrix exponentiation.</i> A matrix can be raised only to an integer power. The matrix, which must be square, is multiplied by itself as many times as the absolute value of the exponent. If the exponent is negative, the result is then inverted.
&*	<i>Elementwise multiplication.</i> Each element of the matrix is multiplied by the corresponding element of the second matrix. The matrices must have the same dimensions, or one must be a scalar.
&/	<i>Elementwise division.</i> Each element of the matrix is divided by the corresponding element of the second matrix. The matrices must have the same dimensions, or one must be a scalar.
&**	<i>Elementwise exponentiation.</i> Each element of the first matrix is raised to the power of the corresponding element of the second matrix. The matrices must have the same dimensions, or one must be a scalar.
:	<i>Sequential integers.</i> This operator creates a vector of consecutive integers from the value preceding the operator to the value following it. You can specify an optional increment following a second colon. See <a href="#">Matrix Notation Shorthand on p. 1133</a> for the principal use of this operator.

- Use these operators only with numeric matrices. The results are undefined when they are used with string matrices.

## Relational Operators

The relational operators are used to compare two matrices, element by element. The result is a matrix of the same size as the (expanded) operands and containing either 1 or 0. The value of each element, 1 or 0, is determined by whether the comparison between the corresponding element of the first matrix with the corresponding element of the second matrix is true or false—1 for true and 0 for false. The matrices being compared must be of the same dimensions unless one of them is a scalar. The relational operators are listed in the following table.

**Table 133-1**

*Relational operators in matrix programs*

>	GT	Greater than
<	LT	Less than
<> or ~ (=)	NE	Not equal to
<=	LE	Less than or equal to
>=	GE	Greater than or equal to
=	EQ	Equal to

- The symbolic and alphabetic forms of these operators are equivalent.

- The symbols representing NE ( $\sim$  or  $\neg$ ) are system dependent. In general, the tilde ( $\sim$ ) is valid for ASCII systems, while the logical-not sign ( $\neg$ ), or whatever symbol is over the number 6 on the keyboard, is valid for IBM EBCDIC systems.
- Use these operators only with numeric matrices. The results are undefined when they are used with string matrices.

## Logical Operators

Logical operators combine two matrices, normally containing values of 1 (true) or 0 (false). When used with other numerical matrices, they treat all positive values as true and all negative and 0 values as false. The logical operators are:

<b>NOT</b>	<i>Reverses the truth of the matrix that follows it.</i> Positive elements yield 0, and negative or 0 elements yield 1.
<b>AND</b>	<i>Both must be true.</i> The matrix $A$ AND $B$ is 1 where the corresponding elements of $A$ and $B$ are both positive and 0 elsewhere.
<b>OR</b>	<i>Either must be true.</i> The matrix $A$ OR $B$ is 1 where the corresponding element of either $A$ or $B$ is positive and 0 where both elements are negative or 0.
<b>XOR</b>	<i>Either must be true but not both.</i> The matrix $A$ XOR $B$ is 1 where one but not both of the corresponding elements of $A$ and $B$ is positive and 0 where both are positive or neither is positive.

## Precedence of Operators

Parentheses can be used to control the order in which complex expressions are evaluated. When the order of evaluation is not specified by parentheses, operations are carried out in the order listed below. The operations higher on the list take precedence over the operations lower on the list.

+ - (unary)

:

\*\* &\*\*

\* &\* &/

+ - (addition and subtraction)

> >= < <= <>=

NOT

AND

OR XOR

Operations of equal precedence are performed left to right of the expressions.

### Examples

```
COMPUTE A = {1, 2, 3; 4, 5, 6}.
COMPUTE B = A + 4.
COMPUTE C = A &** 2.
```

```

COMPUTE D = 2 &** A.
COMPUTE E = A < 5.
COMPUTE F = (C &/ 2) < B.

```

- The results of these COMPUTE statements are:

$$A = \begin{bmatrix} 1 & 2 & 3 \\ 4 & 5 & 6 \end{bmatrix}$$

$$B = \begin{bmatrix} 5 & 6 & 7 \\ 8 & 9 & 10 \end{bmatrix}$$

$$C = \begin{bmatrix} 1 & 4 & 9 \\ 16 & 25 & 36 \end{bmatrix}$$

$$D = \begin{bmatrix} 2 & 4 & 8 \\ 16 & 32 & 64 \end{bmatrix}$$

$$E = \begin{bmatrix} 1 & 1 & 1 \\ 1 & 0 & 0 \end{bmatrix}$$

$$F = \begin{bmatrix} 1 & 1 & 1 \\ 0 & 0 & 0 \end{bmatrix}$$

## MATRIX and Other Commands

A matrix program is a single procedure within a session.

- No active dataset is needed to run a matrix program. If one exists, it is ignored during matrix processing unless you specifically reference it (with an asterisk) on the GET, SAVE, MGET, or MSAVE statements.
- Variables defined in the active dataset are unavailable during matrix processing, except with the GET or MGET statements.
- Matrix variables are unavailable after the END MATRIX command unless you use SAVE or MSAVE to write them to the active dataset.
- You cannot run a matrix program from a syntax window if split-file processing is in effect. If you save the matrix program into a syntax file, however, you can use the INCLUDE command to run the program even if split-file processing is in effect.

## Matrix Statements

The following table lists all of the statements that are accepted within a matrix program. Most of them have the same name as an analogous IBM® SPSS® Statistics command and perform an exactly analogous function. Use only these statements between the MATRIX and END MATRIX commands. Any command not recognized as a valid matrix statement will be rejected by the matrix processor.

Table 133-2  
Valid matrix statements

BREAK	ELSE IF	MSAVE
CALL	END IF	PRINT
COMPUTE	END LOOP	READ
DISPLAY	GET	RELEASE
DO IF	LOOP	SAVE*
ELSE	MGET	WRITE

\*Maximum of 100 SAVE commands in a matrix program.



## ***Exchanging Data with IBM SPSS Statistics Data Files***

Matrix programs can read and write IBM® SPSS® Statistics data files.

- The `GET` and `SAVE` statements read and write ordinary (case-oriented) SPSS Statistics data files, treating each case as a row of a matrix and each ordinary variable as a column.

A matrix program cannot contain more than 100 `SAVE` commands.

- The `MGET` and `MSAVE` statements read and write matrix-format data files, respecting the structure defined by SPSS Statistics when it creates the file. These statements are discussed below.
- Case weighting in a data file is ignored when the file is read into a matrix program.

## ***Using an Active Dataset***

You can use the `GET` statement to read a case-oriented active dataset into a matrix variable. The result is a rectangular data matrix in which cases have become rows and variables have become columns. Special circumstances can affect the processing of this data matrix.

**Split-File Processing.** After a `SPLIT FILE` command, a matrix program executed with the `INCLUDE` command will read one split-file group with each execution of a `GET` statement. This enables you to process the subgroups separately within the matrix program.

**Case Selection.** When a subset of cases is selected for processing, as the result of a `SELECT IF`, `SAMPLE`, or `N OF CASES` command, only the selected cases will be read by the `GET` statement in a matrix program.

**Temporary Transformations.** The entire matrix program is treated as a single procedure. Temporary transformations—those preceded by the `TEMPORARY` command—entered immediately before a matrix program are in effect throughout that program (even if you `GET` the active dataset repeatedly) and are no longer in effect at the end of the matrix program.

**Case Weighting.** Case weighting in an active dataset is ignored when the file is read into a matrix program.

## ***MATRIX and END MATRIX Commands***

The `MATRIX` command, when encountered in a session, invokes the matrix processor, which reads matrix statements until the `END MATRIX` or `FINISH` command is encountered.

- `MATRIX` is a procedure and cannot be entered inside a transformation structure such as `DO IF` or `LOOP`.
- The `MATRIX` procedure does not require an active dataset.
- Comments are removed before subsequent lines are passed to the matrix processor.
- Macros are expanded before subsequent lines are passed to the matrix processor.

The `END MATRIX` command terminates matrix processing and returns control to the command processor.

- The contents of matrix variables are lost after an `END MATRIX` command.
- The active dataset, if present, becomes available again after an `END MATRIX` command.

## ***COMPUTE Statement***

The `COMPUTE` statement carries out most of the calculations in the matrix program. It closely resembles the `COMPUTE` command in the IBM® SPSS® Statistics transformation language.

- The basic specification is the target variable, an equals sign, and the assignment expression. Values of the target variable are calculated according to the specification on the assignment expression.
- The target variable must be named first, and the equals sign is required. Only one target variable is allowed per `COMPUTE` statement.
- Expressions that extract portions of a matrix, such as  $M(1,:)$  or  $M(1:3,4)$ , are allowed to assign values. (For more information, see the topic [Matrix Notation Shorthand on p. 1133](#).) The target variable must be specified as a variable.
- Matrix functions must specify at least one argument enclosed in parentheses. If an expression has two or more arguments, each argument must be separated by a comma. For a complete discussion of the functions and their arguments, see [Matrix Functions on p. 1141](#).

## ***String Values on COMPUTE Statements***

Matrix variables, unlike those in the transformation language, are not checked for data type (numeric or string) when you use them in a `COMPUTE` statement.

- Numerical calculations with matrices containing string values will produce meaningless results.
- One or more elements of a matrix can be set equal to string constants by enclosing the string constants in quotes on a `COMPUTE` statement.
- String values can be copied from one matrix to another with the `COMPUTE` statement.
- There is no way to display a matrix that contains both numeric and string values, if you compute one for some reason.

### ***Example***

```
COMPUTE LABELS={"Observe", "Predict", "Error"}.  
PRINT LABELS /FORMAT=A7.
```

- `LABELS` is a row vector containing three string values.

## ***Arithmetic Operations and Comparisons***

The expression on a `COMPUTE` statement can be formed from matrix constants and variables, combined with the arithmetic, relational, and logical operators discussed above. Matrix constructions and matrix functions are also allowed.

**Examples**

```

COMPUTE PI = 3.14159265.
COMPUTE RSQ = R * R.
COMPUTE FLAGS = EIGENVAL >= 1.
COMPUTE ESTIM = {OBS, PRED, ERR}.

```

- The first statement computes a scalar. Note that the braces are optional on a scalar constant.
- The second statement computes the square of the matrix *R*. *R* can be any square matrix, including a scalar.
- The third statement computes a vector named *FLAGS*, which has the same dimension as the existing vector *EIGENVAL*. Each element of *FLAGS* equals 1 if the corresponding element of *EIGENVAL* is greater than or equal to 1, and 0 if the corresponding element is less than 1.
- The fourth statement constructs a matrix *ESTIM* by concatenating the three vectors or matrices *OBS*, *PRED*, and *ERR*. The component matrices must have the same number of rows.

**Matrix Functions**

The following functions are available in the matrix program. Except where noted, each takes one or more numeric matrices as arguments and returns a matrix value as its result. The arguments must be enclosed in parentheses, and multiple arguments must be separated by commas.

On the following list, matrix arguments are represented by names beginning with *M*. Unless otherwise noted, these arguments can be vectors or scalars. Arguments that must be vectors are represented by names beginning with *V*, and arguments that must be scalars are represented by names beginning with *S*.

<b>ABS(M)</b>	<i>Absolute value.</i> Takes a single argument. Returns a matrix having the same dimensions as the argument, containing the absolute values of its elements.
<b>ALL(M)</b>	<i>Test for all elements nonzero.</i> Takes a single argument. Returns a scalar: 1 if all elements of the argument are nonzero and 0 if any element is zero.
<b>ANY(M)</b>	<i>Test for any element nonzero.</i> Takes a single argument. Returns a scalar: 1 if any element of the argument is nonzero and 0 if all elements are zero.
<b>ARSIN(M)</b>	<i>Inverse sine.</i> Takes a single argument, whose elements must be between $-1$ and $1$ . Returns a matrix having the same dimensions as the argument, containing the inverse sines (arcsines) of its elements. The results are in radians and are in the range from $-\pi/2$ to $\pi/2$ .
<b>ARTAN(M)</b>	<i>Inverse tangent.</i> Takes a single argument. Returns a matrix having the same dimensions as the argument, containing the inverse tangents (arctangents) of its elements, in radians. To convert radians to degrees, multiply by $180/\pi$ , which you can compute as $45/\text{ARTAN}(1)$ . For example, the statement <code>COMPUTE DEGREES=ARTAN(M)*45/ARTAN(1)</code> returns a matrix containing inverse tangents in degrees.

**BLOCK(M1,M2,...)**

*Create a block diagonal matrix.* Takes any number of arguments. Returns a matrix with as many rows as the sum of the rows in all the arguments, and as many columns as the sum of the columns in all the arguments, with the argument matrices down the diagonal and zeros elsewhere. For example, if:

$$A = \begin{bmatrix} 1 & 1 & 1 \\ 1 & 1 & 1 \end{bmatrix}, B = \begin{bmatrix} 2 & 2 \\ 2 & 2 \end{bmatrix}, C = \begin{bmatrix} 3 & 3 & 3 \\ 3 & 3 & 3 \\ 3 & 3 & 3 \end{bmatrix}, \text{ and } D = [4 \quad 4 \quad 4]$$

then:

$$BLOCK(A, B, C, D) = \begin{bmatrix} 1 & 1 & 1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 1 & 1 & 1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 2 & 2 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 2 & 2 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 3 & 3 & 3 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 3 & 3 & 3 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 3 & 3 & 3 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 3 & 3 & 3 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 4 & 4 & 4 \end{bmatrix}$$

**CDFNORM(M)**

*Standard normal cumulative distribution function of elements.* Takes a single argument. Returns a matrix having the same dimensions as the argument, containing the values of the cumulative normal distribution function for each of its elements. If an element of the argument is  $x$ , the corresponding element of the result is a number between 0 and 1, giving the proportion of a normal distribution that is less than  $x$ . For example,  $CDFNORM(\{-1.96, 0, 1.96\})$  results in, approximately,  $\{.025, .5, .975\}$ .

**CHICDF(M,S)**

*Chi-square cumulative distribution function of elements.* Takes two arguments, a matrix of chi-square values and a scalar giving the degrees of freedom (which must be positive). Returns a matrix having the same dimensions as the first argument, containing the values of the cumulative chi-square distribution function for each of its elements. If an element of the first argument is  $x$  and the second argument is  $S$ , the corresponding element of the result is a number between 0 and 1, giving the proportion of a chi-square distribution with  $S$  degrees of freedom that is less than  $x$ . If  $x$  is not positive, the result is 0.

**CHOL(M)**

*Cholesky decomposition.* Takes a single argument, which must be a symmetric positive-definite matrix (a square matrix, symmetric about the main diagonal, with positive eigenvalues). Returns a matrix having the same dimensions as the argument. If  $M$  is a symmetric positive-definite matrix and  $B=CHOL(M)$ , then  $T(B)*B=M$ , where  $T$  is the transpose function defined below.

**CMAX(M)**

*Column maxima.* Takes a single argument. Returns a row vector with the same number of columns as the argument. Each column of the result contains the maximum value of the corresponding column of the argument.

**CMIN(M)**

*Column minima.* Takes a single argument. Returns a row vector with the same number of columns as the argument. Each column of the result contains the minimum value of the corresponding column of the argument.

- COS(M)** *Cosines.* Takes a single argument. Returns a matrix having the same dimensions as the argument, containing the cosines of the elements of the argument. Elements of the argument matrix are assumed to be measured in radians. To convert degrees to radians, multiply by  $\pi/180$ , which you can compute as `ARTAN(1)/45`. For example, the statement `COMPUTE COSINES=COS(DEGREES*ARTAN(1)/45)` returns cosines from a matrix containing elements measured in degrees.
- CSSQ(M)** *Column sums of squares.* Takes a single argument. Returns a row vector with the same number of columns as the argument. Each column of the result contains the sum of the squared values of the elements in the corresponding column of the argument.
- CSUM(M)** *Column sums.* Takes a single argument. Returns a row vector with the same number of columns as the argument. Each column of the result contains the sum of the elements in the corresponding column of the argument.
- DESIGN(M)** *Main-effects design matrix from the columns of a matrix.* Takes a single argument. Returns a matrix having the same number of rows as the argument, and as many columns as the sum of the numbers of unique values in each column of the argument. Constant columns in the argument are skipped with a warning message. The result contains 1 in the row(s) where the value in question occurs in the argument and 0 otherwise. For example, if:
- $$A = \begin{bmatrix} 1 & 2 & 8 \\ 1 & 3 & 8 \\ 2 & 6 & 5 \\ 3 & 3 & 8 \\ 3 & 6 & 5 \end{bmatrix}, \text{ then:}$$
- $$DESIGN(A) = \begin{bmatrix} 1 & 0 & 0 & 1 & 0 & 0 & 1 & 0 \\ 1 & 0 & 0 & 0 & 1 & 0 & 1 & 0 \\ 0 & 1 & 0 & 0 & 0 & 1 & 0 & 1 \\ 0 & 0 & 1 & 0 & 1 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 & 0 & 1 & 0 & 1 \end{bmatrix}$$
- The first three columns of the result correspond to the three distinct values 1, 2, and 3 in the first column of  $A$ ; the fourth through sixth columns of the result correspond to the three distinct values 2, 3, and 6 in the second column of  $A$ ; and the last two columns of the result correspond to the two distinct values 8 and 5 in the third column of  $A$ .
- DET(M)** *Determinant.* Takes a single argument, which must be a square matrix. Returns a scalar, which is the determinant of the argument.
- DIAG(M)** *Diagonal of a matrix.* Takes a single argument. Returns a column vector with as many rows as the minimum of the number of rows and the number of columns in the argument. The  $i$ th element of the result is the value in row  $i$ , column  $i$  of the argument.
- EOF(file)** *End of file indicator.* Normally used after a `READ` statement. Takes a single argument, which must be either a filename in quotes, or a file handle defined on a `FILE HANDLE` command that precedes the matrix program. Returns a scalar equal to 1 if the last attempt to read that file encountered the last record in the file, and equal to 0 if the last attempt did not encounter the last record in the file. Calling the `EOF` function causes a `REREAD` specification on the `READ` statement to be ignored on the next attempt to read the file.

<b>eval(M)</b>	<i>Eigenvalues of a symmetric matrix.</i> Takes a single argument, which must be a symmetric matrix. Returns a column vector with the same number of rows as the argument, containing the eigenvalues of the argument in decreasing numerical order.
<b>EXP(M)</b>	<i>Exponentials of matrix elements.</i> Takes a single argument. Returns a matrix having the same dimensions as the argument, in which each element equals $e$ raised to the power of the corresponding element in the argument matrix.
<b>FCDF(M,S1,S2)</b>	<i>Cumulative F distribution function of elements.</i> Takes three arguments, a matrix of $F$ values and two scalars giving the degrees of freedom (which must be positive). Returns a matrix having the same dimensions as the first argument $M$ , containing the values of the cumulative $F$ distribution function for each of its elements. If an element of the first argument is $x$ and the second and third arguments are $S1$ and $S2$ , the corresponding element of the result is a number between 0 and 1, giving the proportion of an $F$ distribution with $S1$ and $S2$ degrees of freedom that is less than $x$ . If $x$ is not positive, the result is 0.
<b>GINV(M)</b>	<i>Moore-Penrose generalized inverse of a matrix.</i> Takes a single argument. Returns a matrix with the same dimensions as the transpose of the argument. If $A$ is the generalized inverse of a matrix $M$ , then $M*A*M=M$ and $A*M*A=A$ . Both $A*M$ and $M*A$ are symmetric.
<b>GRADE(M)</b>	<i>Ranks elements in a matrix.</i> Takes a single argument. Uses sequential integers for ties.
<b>GSCH(M)</b>	<i>Gram-Schmidt orthonormal basis for the space spanned by the column vectors of a matrix.</i> Takes a single argument, in which there must be as many linearly independent columns as there are rows. (That is, the rank of the argument must equal the number of rows.) Returns a square matrix with as many rows as the argument. The columns of the result form a basis for the space spanned by the columns of the argument.
<b>IDENT(S1 [,S2])</b>	<i>Create an identity matrix.</i> Takes either one or two arguments, which must be scalars. Returns a matrix with as many rows as the first argument and as many columns as the second argument, if any. If the second argument is omitted, the result is a square matrix. Elements on the main diagonal of the result equal 1, and all other elements equal 0.
<b>INV(M)</b>	<i>Inverse of a matrix.</i> Takes a single argument, which must be square and nonsingular (that is, its determinant must not be 0). Returns a square matrix having the same dimensions as the argument. If $A$ is the inverse of $M$ , then $M*A=A*M=I$ , where $I$ is the identity matrix.
<b>KRONEKER(M1,M2)</b>	<i>Kronecker product of two matrices.</i> Takes two arguments. Returns a matrix whose row dimension is the product of the row dimensions of the arguments and whose column dimension is the product of the column dimensions of the arguments. The Kronecker product of two matrices $A$ and $B$ takes the form of an array of scalar products: $A(1,1)*BA(1,2)* B \dots A(1,N)*B$ $A(2,1)*BA(2,2)* B \dots A(2,N)* B$ <p>...</p> $A(M,1)*BA(M,2)*B \dots A(M, N)*B$
<b>LG10(M)</b>	<i>Base 10 logarithms of the elements.</i> Takes a single argument, all of whose elements must be positive. Returns a matrix having the same dimensions as the argument, in which each element is the logarithm to base 10 of the corresponding element of the argument.

<b>LN(M)</b>	<i>Natural logarithms of the elements.</i> Takes a single argument, all of whose elements must be positive. Returns a matrix having the same dimensions as the argument, in which each element is the logarithm to base $e$ of the corresponding element of the argument.
<b>MAGIC(S)</b>	<i>Magic square.</i> Takes a single scalar, which must be 3 or larger, as an argument. Returns a square matrix with $S$ rows and $S$ columns containing the integers from 1 through $S^2$ . All of the row sums and all of the column sums are equal in the result matrix. (The result matrix is only one of several possible magic squares.)
<b>MAKE(S1,S2,S3)</b>	<i>Create a matrix, all of whose elements equal a specified value.</i> Takes three scalars as arguments. Returns an $S1 \times S2$ matrix, all of whose elements equal $S3$ .
<b>MDIAG(V)</b>	<i>Create a square matrix with a specified main diagonal.</i> Takes a single vector as an argument. Returns a square matrix with as many rows and columns as the dimension of the vector. The elements of the vector appear on the main diagonal of the matrix, and the other matrix elements are all 0.
<b>MMAX(M)</b>	<i>Maximum element in a matrix.</i> Takes a single argument. Returns a scalar equal to the numerically largest element in the argument $M$ .
<b>MMIN(M)</b>	<i>Minimum element in a matrix.</i> Takes a single argument. Returns a scalar equal to the numerically smallest element in the argument $M$ .
<b>MOD(M,S)</b>	<i>Remainders after division by a scalar.</i> Takes two arguments, a matrix and a scalar (which must not be 0). Returns a matrix having the same dimensions as $M$ , each of whose elements is the remainder after the corresponding element of $M$ is divided by $S$ . The sign of each element of the result is the same as the sign of the corresponding element of the matrix argument $M$ .
<b>MSSQ(M)</b>	<i>Matrix sum of squares.</i> Takes a single argument. Returns a scalar that equals the sum of the squared values of all of the elements in the argument.
<b>MSUM(M)</b>	<i>Matrix sum.</i> Takes a single argument. Returns a scalar that equals the sum of all of the elements in the argument.
<b>NCOL(M)</b>	<i>Number of columns in a matrix.</i> Takes a single argument. Returns a scalar that equals the number of columns in the argument.
<b>NROW(M)</b>	<i>Number of rows in a matrix.</i> Takes a single argument. Returns a scalar that equals the number of rows in the argument.
<b>RANK(M)</b>	<i>Rank of a matrix.</i> Takes a single argument. Returns a scalar that equals the number of linearly independent rows or columns in the argument.
<b>RESHAPE(M,S1,S2)</b>	<i>Matrix of different dimensions.</i> Takes three arguments, a matrix and two scalars, whose product must equal the number of elements in the matrix. Returns a matrix whose dimensions are given by the scalar arguments. For example, if $M$ is any matrix with exactly 50 elements, then <code>RESHAPE(M, 5, 10)</code> is a matrix with 5 rows and 10 columns. Elements are assigned to the reshaped matrix in order by row.
<b>RMAX(M)</b>	<i>Row maxima.</i> Takes a single argument. Returns a column vector with the same number of rows as the argument. Each row of the result contains the maximum value of the corresponding row of the argument.
<b>RMIN(M)</b>	<i>Row minima.</i> Takes a single argument. Returns a column vector with the same number of rows as the argument. Each row of the result contains the minimum value of the corresponding row of the argument.
<b>RND(M)</b>	<i>Elements rounded to the nearest integers.</i> Takes a single argument. Returns a matrix having the same dimensions as the argument. Each element of the result equals the corresponding element of the argument rounded to an integer.

- RNKORDER(M)** *Ranking of matrix elements in ascending order.* Takes a single argument. Returns a matrix having the same dimensions as the argument  $M$ . The smallest element of the argument corresponds to a result element of 1, and the largest element of the argument to a result element equal to the number of elements, except that ties (equal elements in  $M$ ) are resolved by assigning a rank equal to the arithmetic mean of the applicable ranks. For example, if:
- $$M = \begin{bmatrix} -1 & -21.7 & 8 \\ 0 & 3.91 & -21.7 \\ 8 & 9 & 10 \end{bmatrix}, \text{ then:}$$
- $$RNKORDER(M) = \begin{bmatrix} 3 & 1.5 & 6.5 \\ 4 & 5 & 1.5 \\ 6.5 & 8 & 9 \end{bmatrix}$$
- RSSQ(M)** *Row sums of squares.* Takes a single argument. Returns a column vector having the same number of rows as the argument. Each row of the result contains the sum of the squared values of the elements in the corresponding row of the argument.
- RSUM(M)** *Row sums.* Takes a single argument. Returns a column vector having the same number of rows as the argument. Each row of the result contains the sum of the elements in the corresponding row of the argument.
- SIN(M)** *Sines.* Takes a single argument. Returns a matrix having the same dimensions as the argument, containing the sines of the elements of the argument. Elements of the argument matrix are assumed to be measured in radians. To convert degrees to radians, multiply by  $\pi/180$ , which you can compute as  $ARTAN(1)/45$ . For example, the statement `COMPUTE SINES=SIN(DEGREES*ARTAN(1)/45)` computes sines from a matrix containing elements measured in degrees.
- SOLVE(M1,M2)** *Solution of systems of linear equations.* Takes two arguments, the first of which must be square and nonsingular (its determinant must be nonzero), and the second of which must have the same number of rows as the first. Returns a matrix with the same dimensions as the second argument. If  $M1*X=M2$ , then  $X=SOLVE(M1, M2)$ . In effect, this function sets its result  $X$  equal to  $INV(M1)*M2$ .
- SQRT(M)** *Square roots of elements.* Takes a single argument whose elements must not be negative. Returns a matrix having the same dimensions as the arguments, whose elements are the positive square roots of the corresponding elements of the argument.
- SSCP(M)** *Sums of squares and cross-products.* Takes a single argument. Returns a square matrix having as many rows (and columns) as the argument has columns.  $SSCP(M)$  equals  $T(M)*M$ , where  $T$  is the transpose function defined below.
- SVAL(M)** *Singular values of a matrix.* Takes a single argument. Returns a column vector containing as many rows as the minimum of the numbers of rows and columns in the argument, containing the singular values of the argument in decreasing numerical order. The singular values of a matrix  $M$  are the square roots of the eigenvalues of  $T(M)*M$ , where  $T$  is the transpose function discussed below.
- SWEEP(M,S)** *Sweep transformation of a matrix.* Takes two arguments, a matrix and a scalar, which must be less than or equal to both the number of rows and the number of columns of the matrix. In other words, the pivot element of the matrix, which is  $M(S,S)$ , must exist. Returns a matrix of the same dimensions as  $M$ . Suppose that  $S=\{k\}$  and  $A=SWEEP(M,S)$ . If  $M(k,k)$  is not 0, then
- $$A(k,k) = 1/M(k,k)$$



$$A(i,k) = -M(i,k)/M(k,k), \text{ for } i \text{ not equal to } k$$

$$A(k,j) = M(k,j)/M(k,k), \text{ for } j \text{ not equal to } k$$

$$A(i,j) = (M(i,j)*M(k,k) - M(i,k)*M(k,j))/M(k,k), \text{ for } i,j \text{ not equal to } k$$

and if  $M(k,k)$  equals 0, then

$$A(i,k) = A(k,i) = 0, \text{ for all } i$$

$$A(i,j) = M(i,j), \text{ for } i,j \text{ not equal to } k$$

<b>TCDF(M,S)</b>	<i>Cumulative t distribution function of elements.</i> Takes two arguments, a matrix of $t$ values and a scalar giving the degrees of freedom (which must be positive). Returns a matrix having the same dimensions as $M$ , containing the values of the cumulative $t$ distribution function for each of its elements. If an element of the first argument is $x$ and the second argument is $S$ , then the corresponding element of the result is a number between 0 and 1, giving the proportion of a $t$ distribution with $S$ degrees of freedom that is less than $x$ .
<b>TRACE(M)</b>	<i>Sum of the main diagonal elements.</i> Takes a single argument. Returns a scalar, which equals the sum of the elements on the main diagonal of the argument.
<b>TRANSPOS(M)</b>	<i>Transpose of the matrix.</i> Takes a single argument. Returns the transpose of the argument. TRANSPOS can be shortened to T.
<b>TRUNC(M)</b>	<i>Truncation of elements to integers.</i> Takes a single argument. Returns a matrix having the same dimensions as the argument, whose elements equal the corresponding elements of the argument truncated to integers.
<b>UNIFORM(S1,S2)</b>	<i>Uniformly distributed pseudo-random numbers between 0 and 1.</i> Takes two scalars as arguments. Returns a matrix with the number of rows specified by the first argument and the number of columns specified by the second argument, containing pseudo-random numbers uniformly distributed between 0 and 1.

## CALL Statement

Closely related to the matrix functions are the matrix procedures, which are invoked with the CALL statement. Procedures, similarly to functions, accept arguments enclosed in parentheses and separated by commas. They return their result in one or more of the arguments as noted in the individual descriptions below. They are implemented as procedures rather than as functions so that they can return more than one value or (in the case of SETDIAG) modify a matrix without making a copy of it.

<b>EIGEN(M,var1,var2)</b>	<i>Eigenvectors and eigenvalues of a symmetric matrix.</i> Takes three arguments: a symmetric matrix and two valid variable names to which the results are assigned. If $M$ is a symmetric matrix, the statement CALL EIGEN(M, A, B) will assign to $A$ a matrix having the same dimensions as $M$ , containing the eigenvectors of $M$ as its columns, and will assign to $B$ a column vector having as many rows as $M$ , containing the eigenvalues of $M$ in descending numerical order. The eigenvectors in $A$ are ordered to correspond with the eigenvalues in $B$ ; thus, the first column corresponds to the largest eigenvalue, the second to the second largest, and so on.
<b>SETDIAG(M,V)</b>	<i>Set the main diagonal of a matrix.</i> Takes two arguments, a matrix and a vector. Elements on the main diagonal of $M$ are set equal to the corresponding elements of $V$ . If $V$ is a scalar, all the diagonal elements are set equal to that scalar. Otherwise, if $V$ has fewer elements than the main diagonal of $M$ , remaining elements on the main diagonal are

**SVD(M,var1,var2,var3)**

unchanged. If  $V$  has more elements than are needed, the extra elements are not used. See also the `MDIAG` matrix function.

*Singular value decomposition of a matrix.* Takes four arguments: a matrix and three valid variable names to which the results are assigned. If  $M$  is a matrix, the statement `CALL SVD(M,U,Q,V)` will assign to  $Q$  a diagonal matrix of the same dimensions as  $M$ , and to  $U$  and  $V$  unitary matrices (matrices whose inverses equal their transposes) of appropriate dimensions, such that  $M=U*Q*T(V)$ , where  $T$  is the transpose function defined above. The singular values of  $M$  are in the main diagonal of  $Q$ .

**PRINT Statement**

The `PRINT` statement displays matrices or matrix expressions. Its syntax is as follows:

```
PRINT [matrix expression]
      [/FORMAT="format descriptor"]
      [/TITLE="title"]
      [/SPACE={NEWPAGE}]
           {n          }
      [{/RLABELS=list of quoted names}]
      {/RNAMEs=vector of names      }

      [{/CLABELS=list of quoted names}]
      {/CNAMES=vector of names      }
```

**Matrix Expression**

**Matrix expression** is a single matrix variable name or an expression that evaluates to a matrix. `PRINT` displays the specified matrix.

- The matrix specification must precede any other specifications on the `PRINT` statement. If no matrix is specified, no data will be displayed, but the `TITLE` and `SPACE` specifications will be honored.
- You can specify a matrix name, a matrix raised to a power, or a matrix function (with its arguments in parentheses) by itself, but you must enclose other matrix expressions in parentheses. For example, `PRINT A`, `PRINT INV(A)`, and `PRINT B**DET(T(C)*D)` are all legal, but `PRINT A+B` is not. You must specify `PRINT (A+B)`.
- Constant expressions are allowed.
- A matrix program can consist entirely of `PRINT` statements, without defining any matrix variables.

**FORMAT Keyword**

`FORMAT` specifies a single format descriptor for display of the matrix data.

- All matrix elements are displayed with the same format.
- You can use any printable numeric format (for numeric matrices) or string format (for string matrices) as defined in `FORMATS`.
- The matrix processor will choose a suitable numeric format if you omit the `FORMAT` specification, but a string format such as `A8` is essential when displaying a matrix containing string data.

- String values exceeding the width of a string format are truncated.
- See [Scaling Factor in Displays on p. 1150](#) for default formatting of matrices containing large or small values.

### **TITLE Keyword**

TITLE specifies a title for the matrix displayed. The title must be enclosed in quotes. If it exceeds the maximum display width, it is truncated. The slash preceding TITLE is required, even if it is the only specification on the PRINT statement. If you omit the TITLE specification, the matrix name or expression from the PRINT statement is used as a default title.

### **SPACE Keyword**

SPACE controls output spacing before printing the title and the matrix. You can specify either a positive number or the keyword NEWPAGE. The slash preceding SPACE is required, even if it is the only specification on the PRINT statement.

NEWPAGE	<i>Start a new page before printing the title.</i>
n	<i>Skip n lines before displaying the title.</i>

### **RLABELS Keyword**

RLABELS allows you to supply row labels for the matrix.

- The labels must be separated by commas.
- Enclose individual labels in quotes if they contain embedded commas or if you want to preserve lowercase letters. Otherwise, quotes are optional.
- If too many names are supplied, the extras are ignored. If not enough names are supplied, the last rows remain unlabeled.

### **RNAMES Keyword**

RNAMES allows you to supply the name of a vector or a vector expression containing row labels for the matrix.

- Either a row vector or a column vector can be used, but the vector must contain string data.
- If too many names are supplied, the extras are ignored. If not enough names are supplied, the last rows remain unlabeled.

### **CLABELS Keyword**

CLABELS allows you to supply column labels for the matrix.

- The labels must be separated by commas.

- Enclose individual labels in quotes if they contain embedded commas or if you want to preserve lowercase letters. Otherwise, quotes are optional.
- If too many names are supplied, the extras are ignored. If not enough names are supplied, the last columns remain unlabeled.

## **CNAMES Keyword**

CNAMES allows you to supply the name of a vector or a vector expression containing column labels for the matrix.

- Either a row vector or a column vector can be used, but the vector must contain string data.
- If too many names are supplied, the extras are ignored. If not enough names are supplied, the last columns remain unlabeled.

## **Scaling Factor in Displays**

When a matrix contains very large or very small numbers, it may be necessary to use scientific notation to display the data. If you do not specify a display format, the matrix processor chooses a power-of-10 multiplier that will allow the largest value to be displayed, and it displays this multiplier on a heading line before the data. The multiplier is not displayed for each element in the matrix. The displayed values, multiplied by the power of 10 that is indicated in the heading, equal the actual values (possibly rounded).

- Values that are very small, relative to the multiplier, are displayed as 0.
- If you explicitly specify a scientific-notation format (`Ew.d`), each matrix element is displayed using that format. This permits you to display very large and very small numbers in the same matrix without losing precision.

### **Example**

```
COMPUTE M = {.0000000001357, 2.468, 3690000000}.
PRINT M /TITLE "Default format".
PRINT M /FORMAT "E13" /TITLE "Explicit exponential format".
```

- The first PRINT subcommand uses the default format with  $10^9$  as the multiplier for each element of the matrix. This results in the following output:

Figure 133-1

```
Default format
10 ** 9 X
.000000000 .000000002 3.690000000
```

Note that the first element is displayed as 0 and the second is rounded to one significant digit.

- An explicitly specified exponential format on the second PRINT subcommand allows each element to be displayed with full precision, as the following output shows:

Figure 133-2

```
Explicit exponential format
1.E-010 2.E+000 4.E+009
```

## Matrix Control Structures

The matrix language includes two structures that allow you to alter the flow of control within a matrix program.

- The `DO IF` statement tests a logical expression to determine whether one or more subsequent matrix statements should be executed.
- The `LOOP` statement defines the beginning of a block of matrix statements that should be executed repeatedly until a termination criterion is satisfied or a `BREAK` statement is executed.

These statements closely resemble the `DO IF` and `LOOP` commands in the IBM® SPSS® Statistics transformation language. In particular, these structures can be nested within one another as deeply as the available memory allows.

### DO IF Structures

A `DO IF` structure in a matrix program affects the flow of control exactly as the analogous commands affect a transformation program, except that missing-value considerations do not arise in a matrix program. The syntax of the `DO IF` structure is as follows:

```
DO IF [(logical expression)]
    matrix statements
[ELSE IF [(logical expression)]]
    matrix statements
[ELSE IF...]
.
.
.
[ELSE]
    matrix statements
END IF.
```

- The `DO IF` statement marks the beginning of the structure, and the `END IF` statement marks its end.
- The `ELSE IF` statement is optional and can be repeated as many times as desired within the structure.
- The `ELSE` statement is optional. It can be used only once and must follow any `ELSE IF` statements.
- The `END IF` statement must follow any `ELSE IF` and `ELSE` statements.
- The `DO IF` and `ELSE IF` statements must contain a logical expression, normally one involving the relational operators `EQ`, `GT`, and so on. However, the matrix language allows any expression that evaluates to a scalar to be used as the logical expression. Scalars greater than 0 are considered true, and scalars less than or equal to 0 are considered false.

A `DO IF` structure affects the flow of control within a matrix program as follows:

- If the logical expression on the `DO IF` statement is true, the statements immediately following the `DO IF` are executed up to the next `ELSE IF` or `ELSE` in the structure. Control then passes to the first statement following the `END IF` for that structure.
- If the expression on the `DO IF` statement is false, control passes to the first `ELSE IF`, where the logical expression is evaluated. If this expression is true, statements following the `ELSE IF` are executed up to the next `ELSE IF` or `ELSE` statement, and control passes to the first statement following the `END IF` for that structure.
- If the expressions on the `DO IF` and the first `ELSE IF` statements are both false, control passes to the next `ELSE IF`, where that logical expression is evaluated. If none of the expressions is true on any of the `ELSE IF` statements, statements following the `ELSE` statement are executed up to the `END IF` statement, and control falls out of the structure.
- If none of the expressions on the `DO IF` statement or the `ELSE IF` statements is true and there is no `ELSE` statement, control passes to the first statement following the `END IF` for that structure.

## **LOOP Structures**

A `LOOP` structure in a matrix program affects the flow of control exactly as the analogous commands affect transformation program, except that missing-value considerations do not arise in a matrix program. Its syntax is as follows:

```
LOOP [varname=n TO m [BY k]] [IF [(logical expression)]
matrix statements
[BREAK]
matrix statements
END LOOP [IF [(logical expression)]]
```

The matrix statements specified between `LOOP` and `END LOOP` are executed repeatedly until one of the following conditions is met:

- A logical expression on the `IF` clause of the `LOOP` statement is evaluated as false.
- An index variable used on the `LOOP` statement passes beyond its terminal value.
- A logical expression on the `IF` clause of the `END LOOP` statement is evaluated as true.
- A `BREAK` statement is executed within the loop structure (but outside of any nested loop structures).

Note: Unlike the `LOOP` command (outside the matrix language), the index value of a matrix `LOOP` structure does not override the maximum number of loops controlled by `SET MXLOOPS`. You must explicitly set the `MXLOOPS` value to a value high enough to accommodate the index value. [For more information, see the topic `MXLOOPS` Subcommand on p. 1871.](#)

### ***Index Clause on the LOOP Statement***

An index clause on a `LOOP` statement creates an index variable whose name is specified immediately after the keyword `LOOP`. The variable is assigned an initial value of  $n$ . Each time through the loop, the variable is tested against the terminal value  $m$  and incremented by the increment value  $k$  if  $k$  is specified or by 1 if  $k$  is not specified. When the index variable is greater than  $m$  for positive increments or less than  $m$  for negative increments, control passes to the statement after the `END LOOP` statement.

- Both the index clause and the `IF` clause are optional. If both are present, the index clause must appear first.
- The index variable must be scalar with a valid matrix variable name.
- The initial value,  $n$ , the terminal value,  $m$ , and the increment,  $k$  (if present), must be scalars or matrix expressions evaluating to scalars. Non-integer values are truncated to integers before use.
- If the keyword `BY` and the increment  $k$  are absent, an increment of 1 is used.

### ***IF Clause on the LOOP Statement***

The logical expression is evaluated before each iteration of the loop structure. If it is false, the loop terminates and control passes to the statement after `END LOOP`.

- The `IF` clause is optional. If both the index clause and the `IF` clause are present, the index clause must appear first.
- As in the `DO IF` structure, the logical expression of the `IF` clause is evaluated as scalar, with positive values being treated as true and 0 or negative values, as false.

### ***IF Clause on the END LOOP Statement***

When an `IF` clause is present on an `END LOOP` statement, the logical expression is evaluated after each iteration of the loop structure. If it is true, the loop terminates and control passes to the statement following the `END LOOP` statement.

- The `IF` clause is optional.
- As in the `LOOP` statement, the logical expression of the `IF` clause is evaluated as scalar, with positive values being treated as true and 0 or negative values, as false.

### ***BREAK Statement***

The `BREAK` statement within a loop structure transfers control immediately to the statement following the (next) `END LOOP` statement. It is normally placed within a `DO IF` structure inside the `LOOP` structure to exit the loop when the specified conditions are met.

#### ***Example***

```
LOOP LOCATION = 1, NROW(VEC).  
+ DO IF (VEC(LOCATION) = TARGET).  
+   BREAK.
```

```
+ END IF.
END LOOP.
```

- This loop searches for the (first) location of a specific value, *TARGET*, in a vector, *VEC*.
- The `DO IF` statement checks whether the vector element indexed by *LOCATION* equals the target.
- If so, the `BREAK` statement transfers control out of the loop, leaving *LOCATION* as the index of *TARGET* in *VEC*.

## ***READ Statement: Reading Character Data***

The `READ` statement reads data into a matrix or submatrix from a character-format file—that is, a file containing ordinary numbers or words in readable form. The syntax for the `READ` statement is:

```
READ variable reference
  [/FILE = file reference]
  [/FIELD = c1 TO c2 [BY w]]
  [/SIZE = size expression]
  [/MODE = {RECTANGULAR}]
           {SYMMETRIC }
  [/REREAD]
  [/FORMAT = format descriptor]
```

- The file can contain values in freefield or fixed-column format. The data can appear in any of the field formats supported by `DATA LIST`.
- More than one matrix can be read from a single input record by rereading the record.
- If the end of the file is encountered during a `READ` operation (that is, fewer values are available than the number of elements required by the specified matrix size), a warning message is displayed and the contents of the unread elements of the matrix are unpredictable.

## ***Variable Specification***

The variable reference on the `READ` statement is a matrix variable name, with or without indexes.

For a name without indexes:

- `READ` creates the specified matrix variable.
- The matrix need not exist when `READ` is executed.
- If the matrix already exists, it is replaced by the matrix read from the file.
- You must specify the size of the matrix using the `SIZE` specification.

For an indexed name:

- `READ` creates a submatrix from an existing matrix.
- The matrix variable named must already exist.



- You can define any submatrix with indexes; for example,  $M(:, I)$ . To define an entire existing matrix, specify  $M(:, :)$ .
- The `SIZE` specification can be omitted. If specified, its value must match the size of the specified submatrix.

### **FILE Specification**

`FILE` designates the character file containing the data. It can be an actual filename in quotes, or a file handle defined on a `FILE HANDLE` command that precedes the matrix program.

- The filename or handle must specify an existing file containing character data, not a data file in IBM® SPSS® Statistics format or a specially formatted file of another kind, such as a spreadsheet file.
- The `FILE` specification is required on the first `READ` statement in a matrix program (first in order of appearance, not necessarily in order of execution). If you omit the `FILE` specification from a later `READ` statement, the statement uses the most recently named file (in order of appearance) on a `READ` statement in the same matrix program.

### **FIELD Specification**

`FIELD` specifies the column positions of a fixed-format record where the data for matrix elements are located.

- The `FIELD` specification is required.
- `startcol` is the number of the leftmost column of the input area.
- `endcol` is the number of the rightmost column of the input area.
- Both `startcol` and `endcol` are required and both must be constants. For example, `FIELD = 9 TO 72` specifies that values to be read appear between columns 9 and 72 (inclusive) of each input record.
- The `BY` clause, if present, indicates that each value appears within a fixed set of columns on the input record; that is, one value is separated from the next by its column position rather than by a space or comma. `Width` is the width of the area designated for each value. For example, `FIELD = 1 TO 80 BY 10` indicates that there are eight possible values per record and that one will appear between columns 1 and 10 (inclusive), another between columns 11 and 20, and so on, up to columns 71 and 80. The `BY` value must evenly divide the length of the field. That is, `endcol-startcol+1` must be a multiple of the width.
- You can use the `FORMAT` specification to supply the same information as the `BY` clause of the `FIELD` specification. If you omit the `BY` clause and do not specify a format on the `FORMAT` specification, `READ` assumes that values are separated by blanks or commas within the designated field.

### **SIZE Specification**

The `SIZE` specification is a matrix expression that, when evaluated, specifies the size of the matrix to be read.

- The expression should evaluate to a two-element row or column vector. The first element designates the number of rows in the matrix to be read; the second element gives the number of columns.
- Values of the `SIZE` specification are truncated to integers if necessary.
- The size expression may be a constant, such as `{5;5}`, or a matrix variable name, such as `MSIZE`, or any valid expression, such as `INFO(1, :)`.
- If you use a scalar as the size expression, a column vector containing that number of rows is read. Thus, `SIZE=1` reads a scalar, and `SIZE=3` reads a  $3 \times 1$  column vector.

You must include a `SIZE` specification whenever you name an entire matrix (rather than a submatrix) on the `READ` statement. If you specify a submatrix, the `SIZE` specification is optional but, if included, must agree with the size of the specified submatrix.

### **MODE Specification**

`MODE` specifies the format of the matrix to be read in. It can be either rectangular or symmetric. If the `MODE` specification is omitted, the default is `RECTANGULAR`.

<b>RECTANGULAR</b>	<i>Matrix is completely represented in file.</i> Each row begins on a new record, and all entries in that row are present on that and (possibly) succeeding records. This is the default if the <code>MODE</code> specification is omitted.
<b>SYMMETRIC</b>	<i>Elements of the matrix below the main diagonal are the same as those above it.</i> Only matrix elements on and below the main diagonal are read; elements above the diagonal are set equal to the corresponding symmetric elements below the diagonal. Each row is read beginning on a new record, although it may span more than one record. Only a single value is read from the first record, two values are read from the second, and so on.

- If `SYMMETRIC` is specified, the matrix processor first checks that the number of rows and the number of columns are the same. If the numbers, specified either on `SIZE` or on the variable reference, are not the same, an error message is displayed and the command is not executed.

### **REREAD Specification**

The `REREAD` specification indicates that the current `READ` statement should begin with the last record read by a previous `READ` statement.

- `REREAD` has no further specifications.
- `REREAD` cannot be used on the first `READ` statement to read from a file.
- If you omit `REREAD`, the `READ` statement begins with the first record following the last one read by the previous `READ` statement.
- The `REREAD` specification is ignored on the first `READ` statement following a call to the `EOF` function for the same file.

## **FORMAT Specification**

FORMAT specifies how the matrix processor should interpret the input data. The format descriptor can be any valid IBM® SPSS® Statistics data format, such as F6, E12.2, or A6, or it can be a type code; for example, F, E, or A.

- If you omit the FORMAT specification, the default is F.
- You can specify the width of fixed-size data fields with either a FORMAT specification or a BY clause on a FIELD specification. You can include it in both places only if you specify the same value.
- If you do not include either a FORMAT or a BY clause on FIELD, READ expects values separated by blanks or commas.
- An additional way of specifying the width is to supply a repetition factor without a width (for example, 10F, 5COMMA, or 3E). The field width is then calculated by dividing the width of the whole input area on the FIELD specification by the repetition factor. A format with a digit for the repetition factor must be enclosed in quotes.
- Only one format can be specified. A specification such as `FORMAT='5F2.0 3F3.0 F2.0'` is invalid.

## **WRITE Statement: Writing Character Data**

WRITE writes the value of a matrix expression to an external file. The syntax of the WRITE statement is:

```
WRITE matrix expression
  [/OUTFILE = file reference]
  /FIELD = startcol TO endcol [BY width]
  [/MODE = {RECTANGULAR}]
           {TRIANGULAR }
  [/HOLD]
  [/FORMAT = format descriptor]
```

## **Matrix Expression Specification**

Specify any matrix expression that evaluates to the value(s) to be written.

- The matrix specification must precede any other specifications on the WRITE statement.
- You can specify a matrix name, a matrix raised to a power, or a matrix function (with its arguments in parentheses) by itself, but you must enclose other matrix expressions in parentheses. For example, `WRITE A`, `WRITE INV(A)`, or `WRITE B**DET(T(C)*D)` is legal, but `WRITE A+B` is not. You must specify `WRITE (A+B)`.
- Constant expressions are allowed.

## **OUTFILE Specification**

OUTFILE designates the character file to which the matrix expression is to be written. The file reference can be an actual filename in quotes or a file handle defined on a FILE HANDLE command that precedes the matrix program. The filename or file handle must be a valid file specification.

- The `OUTFILE` specification is required on the first `WRITE` statement in a matrix program (first in order of appearance, not necessarily in order of execution).
- If you omit the `OUTFILE` specification from a later `WRITE` statement, the statement uses the most recently named file (in order of appearance) on a `WRITE` statement in the same matrix program.

## ***FIELD Specification***

`FIELD` specifies the column positions of a fixed-format record to which the data should be written.

- The `FIELD` specification is required.
- The start column, `c1`, is the number of the leftmost column of the output area.
- The end column, `c2`, is the number of the rightmost column of the output area.
- Both `c1` and `c2` are required, and both must be constants. For example, `FIELD = 9 TO 72` specifies that values should be written between columns 9 and 72 (inclusive) of each output record.
- The `BY` clause, if present, indicates how many characters should be allocated to the output value of a single matrix element. The value `w` is the width of the area designated for each value. For example, `FIELD = 1 TO 80 BY 10` indicates that up to eight values should be written per record, and that one should go between columns 1 and 10 (inclusive), another between columns 11 and 20, and so on up to columns 71 and 80. The value on the `BY` clause must evenly divide the length of the field. That is,  $c2 - c1 + 1$  must be a multiple of `w`.
- You can use the `FORMAT` specification (see below) to supply the same information as the `BY` clause. If you omit the `BY` clause from the `FIELD` specification and do not specify a format on the `FORMAT` specification, `WRITE` uses freefield format, separating matrix elements by single blank spaces.

## ***MODE Specification***

`MODE` specifies the format of the matrix to be written. If `MODE` is not specified, the default is `RECTANGULAR`.

<b>RECTANGULAR</b>	<i>Write the entire matrix.</i> Each row starts a new record, and all of the values in that row are present in that and (possibly) subsequent records. This is the default if the <code>MODE</code> specification is omitted.
<b>TRIANGULAR</b>	<i>Write only the lower triangular entries and the main diagonal.</i> Each row begins a new record and may span more than one record. This mode may save file space.

- A matrix written with `MODE = TRIANGULAR` must be square, but it need not be symmetric. If it is not, values in the upper triangle are not written.
- A matrix written with `MODE = TRIANGULAR` may be read with `MODE = SYMMETRIC`.

## **HOLD Specification**

HOLD causes the last line written by the current WRITE statement to be held so that the next WRITE to that file will write on the same line. Use HOLD to write more than one matrix on a line.

## **FORMAT Specification**

FORMAT indicates how the internal (binary) values of matrix elements should be converted to character format for output.

- The format descriptor is any valid IBM® SPSS® Statistics data format, such as F6, E12.2, or A6, or it can be a format type code, such as F, E, or A. It specifies how the written data are encoded and, if a width is specified, how wide the fields containing the data are. (See FORMATS for valid formats.)
- If you omit the FORMAT specification, the default is F.
- The data field widths may be specified either here or after BY on the FIELD specification. You may specify the width in both places only if you give the same value.
- An additional way of specifying the width is to supply a repetition factor without a width (for example, 10F or 5COMMA). The field width is then calculated by dividing the width of the whole output area on the FIELD specification by the repetition factor. A format with a digit for the repetition factor must be enclosed in quotes.
- If the field width is not specified in any of these ways, then the freefield format is used—matrix values are written separated by one blank, and each value occupies as many positions as necessary to avoid the loss of precision. Each row of the matrix is written starting with a new output record.
- Only one format descriptor can be specified. Do *not* try to specify more than one format; for example, '5F2.0 3F3.0 F2.0' is invalid as a FORMAT specification on WRITE.

## **GET Statement: Reading IBM SPSS Statistics Data Files**

GET reads matrices from an external IBM® SPSS® Statistics data file or from the active dataset. The syntax of GET is as follows:

```
GET variable reference
  [/FILE={file reference}]
  {*
  }
  [/VARIABLES = variable list]
  [/NAMES = names vector]
  [/MISSING = {ACCEPT}]
  {OMIT }
  {value }
  [/SYSMIS = {OMIT }]
  {value}
```

## **Variable Specification**

The variable reference on the GET statement is a matrix variable name with or without indexes.

For a name without indexes:

- GET creates the specified matrix variable.

- The size of the matrix is determined by the amount of data read from the IBM® SPSS® Statistics data file or the active dataset.
- If the matrix already exists, it is replaced by the matrix read from the file.

For an indexed name:

- GET creates a submatrix from an existing matrix.
- The matrix variable named must already exist.
- You can define any submatrix with indexes; for example,  $M(:, I)$ . To define an entire existing matrix, specify  $M(:, :)$ .
- The indexes, along with the size of the existing matrix, specify completely the size of the submatrix, which must agree with the dimensions of the data read from the SPSS Statistics data file.
- The specified submatrix is replaced by the matrix elements read from the SPSS Statistics data file.

### **FILE Specification**

FILE designates the IBM® SPSS® Statistics data file to be read. Use an asterisk, or simply omit the FILE specification, to designate the current active dataset.

- The file reference can be either a filename enclosed in quotes, a previously declared dataset in the current session, or a file handle defined on a FILE HANDLE command that precedes the matrix program.
- If you omit the FILE specification, the active dataset is used.
- In a matrix program executed with the INCLUDE command, if a SPLIT FILE command is in effect, a GET statement that references the active dataset will read a single split-file group of cases. (A matrix program cannot be executed from a syntax window if a SPLIT FILE command is in effect.)

### **VARIABLES Specification**

VARIABLES specifies a list of variables to be read from the IBM® SPSS® Statistics data file.

- The keyword TO can be used to reference consecutive variables on the data file.
- The variable list can consist of the keyword ALL to get all the variables in the data file. ALL is the default if the VARIABLES specification is omitted.
- All variables read from the data file should be numeric. If a string variable is specified, a warning message is issued and the string variable is skipped.

#### **Example**

```
GET M /VARIABLES = AGE, RESIDE, INCOME TO HEALTH.
```

- The variables *AGE*, *RESIDE*, and *INCOME TO HEALTH* from the active dataset will form the columns of the matrix *M*.

## ***NAMES Specification***

NAMES specifies a vector to store the variable names from the data file.

- If you omit the NAMES specification, the variable names are not available to the MATRIX procedure.

## ***MISSING Specification***

MISSING specifies how missing values declared for the IBM® SPSS® Statistics data file should be handled.

- The MISSING specification is required if the data file contains missing values for any variable being read.
- If you omit the MISSING specification and a missing value is encountered for a variable being read, an error message is displayed and the GET statement is not executed.

The following keywords are available on the MISSING specification. There is no default.

<b>ACCEPT</b>	<i>Accept user-missing values for entry.</i> If the system-missing value exists for a variable to be read, you must specify SYSMIS to indicate how the system-missing value should be handled.
<b>OMIT</b>	<i>Skip an entire observation when a variable with a missing value is encountered.</i>
<b>value</b>	<i>Recode all missing values encountered (including the system-missing value) to the specified value for entry.</i> The replacement value can be any numeric constant.

## ***SYSMIS Specification***

SYSMIS specifies how system-missing values should be handled when you have specified ACCEPT on MISSING.

- The SYSMIS specification is ignored unless ACCEPT is specified on MISSING.
- If you specify ACCEPT on MISSING but omit the SYSMIS specification, and a system-missing value is encountered for a variable being read, an error message is displayed and the GET statement is not executed.

The following keywords are available on the SYSMIS specification. There is no default.

<b>OMIT</b>	<i>Skip an entire observation when a variable with a system-missing value is encountered.</i>
<b>value</b>	<i>Recode all system-missing values encountered to the specified value for entry.</i> The replacement value can be any numeric constant.

### ***Example***

```
GET SCORES
/VARIABLES = TEST1, TEST2, TEST3
/NAMES = VARNAMES
/MISSING = ACCEPT
/SYSMIS = -1.0.
```

- A matrix named *SCORES* is read from the active dataset.

- The variables *TEST1*, *TEST2*, and *TEST3* form the columns of the matrix, while the cases in the active dataset form the rows.
- A vector named *VARNAMES*, whose three elements contain the variable names *TEST1*, *TEST2*, and *TEST3*, is created.
- User-missing values defined in the active dataset are accepted into the matrix *SCORES*.
- System-missing values in the active dataset are converted to the value  $-1$  in the matrix *SCORES*.

## ***SAVE Statement: Writing IBM SPSS Statistics Data Files***

SAVE writes matrices to an external IBM® SPSS® Statistics data file or to the current active dataset. The rows of the matrix expression become cases, and the columns become variables. The syntax of the SAVE statement is as follows:

```
SAVE matrix expression
  [/OUTFILE = {file reference}]
  { *
  [/VARIABLES = variable list]
  [/NAMES = names vector]
  [/STRINGS = variable list]
```

### ***Matrix Expression Specification***

The matrix expression following the keyword SAVE is any matrix language expression that evaluates to the value(s) to be written to a data file.

- The matrix specification must precede any other specifications on the SAVE statement.
- You can specify a matrix name, a matrix raised to a power, or a matrix function (with its arguments in parentheses) by itself, but you must enclose other matrix expressions in parentheses. For example, SAVE A, SAVE INV(A), or SAVE B\*\*DET(T(C)\*D) is legal, but SAVE A+B is not. You must specify SAVE (A+B).
- Constant expressions are allowed.

### ***OUTFILE Specification***

OUTFILE designates the file to which the matrix expression is to be written. It can be an actual filename in quotes, a previously declared dataset in the current session, or a file handle defined on a FILE HANDLE command that precedes the matrix program.

- To save a matrix expression as the active dataset, specify an asterisk (\*). If there is no active dataset, one will be created; if there is one, it is replaced by the saved matrices.
- The OUTFILE specification is required on the first SAVE statement in a matrix program (first in order of appearance, not necessarily in order of execution). If you omit the OUTFILE specification from a later SAVE statement, the statement uses the most recently named file (in order of appearance) on a SAVE statement in the same matrix program.



- If more than one `SAVE` statement writes to the active dataset in a single matrix program, the dictionary of the new active dataset is written on the basis of the information given by the first such `SAVE`. All of the subsequently saved matrices are appended to the new active dataset as additional cases. If the number of columns differs, an error occurs.
- When you execute a matrix program with the `INCLUDE` command, the `SAVE` statement creates a new IBM® SPSS® Statistics data file at the end of the matrix program's execution, so any attempt to `GET` the data file obtains the original data file, if any.
- When you execute a matrix program from a syntax window, `SAVE` creates a new SPSS Statistics data file immediately, but the file remains open, so you cannot `GET` it until after the `END MATRIX` statement.

## **VARIABLES Specification**

You can provide variable names for the IBM® SPSS® Statistics data file with the `VARIABLES` specification. The variable list is a list of valid variable names separated by commas.

- You can use the `TO` convention, as shown in the example below.
- You can also use the `NAMES` specification, discussed below, to provide variable names.

### **Example**

```
SAVE {A,B,X,Y} /OUTFILE=*  
  /VARIABLES = A,B,X1 TO X50,Y1,Y2.
```

- The matrix expression on the `SAVE` statement constructs a matrix from two column vectors  $A$  and  $B$  and two matrices  $X$  and  $Y$ . All four matrix variables must have the same number of rows so that this matrix construction will be valid.
- The `VARIABLES` specification provides descriptive names so that the variable names in the new active dataset will resemble the names used in the matrix program.

## **NAMES Specification**

As an alternative to the explicit list on the `VARIABLES` specification, you can specify a name list with a vector containing string values. The elements of this vector are used as names for the variables.

- The `NAMES` specification on `SAVE` is designed to complement the `NAMES` specification on the `GET` statement. Names extracted from IBM® SPSS® Statistics data files can be used in a new data file by specifying the same vector name on both `NAMES` specifications.
- If you specify both `VARIABLES` and `NAMES`, a warning message is displayed and the `VARIABLES` specification is used.
- If you omit both the `VARIABLES` and `NAMES` specifications, or if you do not specify names for all columns of the matrix, the `MATRIX` procedure creates default names. The names have the form  $COLn$ , where  $n$  is the column number.

### ***STRINGS Specification***

The `STRINGS` specification provides the names of variables that contain short string data rather than numeric data.

- By default, all variables are assumed to be numeric.
- The variable list specification following `STRINGS` consists of a list of variable names separated by commas. The names must be among those used by `SAVE`.

### ***MGET Statement: Reading Matrix Data Files***

`MGET` reads a matrix-format data file. `MGET` puts the data it reads into separate matrix variables. It also names these new variables automatically. The syntax of `MGET` is as follows:

```
MGET [ [/] FILE = file reference]
      [/TYPE = {COV   } ]
              {CORR  }
              {MEAN  }
              {STDDEV}
              {N     }
              {COUNT }
```

- Since `MGET` assigns names to the matrices it reads, do not specify matrix names on the `MGET` statement.

### ***FILE Specification***

`FILE` designates a matrix-format data file. See [MATRIX DATA on p. 1171](#) for a discussion of matrix-format data files. To designate the active dataset (if it is a matrix-format data file), use an asterisk, or simply omit the `FILE` specification.

- The file reference can be either a filename enclosed in quotes or a file handle defined on a `FILE HANDLE` command that precedes the matrix program.
- The same matrix-format data file can be read more than once.
- If you omit the `FILE` specification, the current active dataset is used.
- `MGET` ignores the `SPLIT FILE` command when reading the active dataset. It does honor the split-file groups that were in effect when the matrix-format data file was created.
- The maximum number of split-file groups that can be read is 99.
- The maximum number of cells that can be read is 99.

### ***TYPE Specification***

`TYPE` specifies the rowtype(s) to read from the matrix-format data file.

- By default, records of all rowtypes are read.
- If the matrix-format data file does not contain rows of the requested type, an error occurs.

Valid keywords on the TYPE specification are:

<b>COV</b>	<i>A matrix of covariances.</i>
<b>CORR</b>	<i>A matrix of correlation coefficients.</i>
<b>MEAN</b>	<i>A vector of means.</i>
<b>STDDEV</b>	<i>A vector of standard deviations.</i>
<b>N</b>	<i>A vector of numbers of cases.</i>
<b>COUNT</b>	<i>A vector of counts.</i>

### **Names of Matrix Variables from MGET**

- The MGET statement automatically creates matrix variable names for the matrices it reads.
- All new variables created by MGET are reported to the user.
- If a matrix variable already exists with the same name that MGET chose for a new variable, the new variable is not created and a warning is issued. The RELEASE statement can be used to get rid of a variable. A COMPUTE statement followed by RELEASE can be used to change the name of an existing matrix variable.

MGET constructs variable names in the following manner:

- The first two characters of the name identify the row type. If there are no cells and no split file groups, these two characters constitute the name:

<i>CV</i>	A covariance matrix (rowtype COV)
<i>CR</i>	A correlation matrix (rowtype CORR)
<i>MN</i>	A vector of means (rowtype MEAN)
<i>SD</i>	A vector of standard deviations (rowtype STDDEV)
<i>NC</i>	A vector of numbers of cases (rowtype N)
<i>CN</i>	A vector of counts (rowtype COUNT)

- Characters 3–5 of the variable name identify the cell number or the split-group number. Cell identifiers consist of the letter *F* and a two-digit cell number. Split-group identifiers consist of the letter *S* and a two-digit split-group number; for example, *MNF12* or *SDS22*.
- If there are both cells and split groups, characters 3–5 identify the cell and characters 6–8 identify the split group. The same convention for cell or split-file numbers is used; for example, *CRF12S21*.
- After the name is constructed as described above, any leading zeros are removed from the cell number and the split-group number; for example, *CNF2S99* or *CVF2S1*.

### **MSAVE Statement: Writing Matrix Data Files**

The MSAVE statement writes matrix expressions to a matrix-format data file that can be used as matrix input to other procedures. (See [MATRIX DATA](#) on p. 1171 for a discussion of matrix-format data files.) The syntax of MSAVE is as follows:

```
MSAVE matrix expression
  /TYPE = {COV  }
         {CORR }
```

```

        {MEAN  }
        {STDDEV}
        {N     }
        {COUNT}
[/OUTFILE = {file reference}]
        {*     }
[/VARIABLES = variable list]
[/SNAMES = variable list]
[/SPLIT = split vector]
[/FNAMES = variable list]
[/FACTOR = factor vector]

```

- Only one matrix-format data file can be saved in a single matrix program.
- Each MSAVE statement writes records of a single rowtype. Therefore, several MSAVE statements will normally be required to write a complete matrix-format data file.
- Most specifications are retained from one MSAVE statement to the next so that it is not necessary to repeat the same specifications on a series of MSAVE statements. The exception is the FACTOR specification, as noted below.

### Example

```

MSAVE M /TYPE=MEAN /OUTFILE=CORRMAT /VARIABLES=V1 TO V8.
MSAVE S /TYPE STDDEV.
MSAVE MAKE(1,8,24) /TYPE N.
MSAVE C /TYPE CORR.

```

- The series of MSAVE statements save the matrix variables *M*, *S*, and *C*, which contain, respectively, vectors of means and standard deviations and a matrix of correlation coefficients. The matrix-format data file thus created is suitable for use in a procedure such as FACTOR.
- The first MSAVE statement saves *M* as a vector of means. This statement specifies OUTFILE, a previously defined file handle, and VARIABLES, a list of variable names to be used in the IBM® SPSS® Statistics data file.
- The second MSAVE statement saves *S* as a vector of standard deviations. Note that the OUTFILE and VARIABLES specifications do not have to be repeated.
- The third MSAVE statement saves a vector of case counts. The matrix function MAKE constructs an eight-element vector with values equal to the case count (24 in this example).
- The last MSAVE statement saves *C*, an  $8 \times 8$  matrix, as the correlation matrix.

### Matrix Expression Specification

- The matrix expression must be specified first on the MSAVE statement.
- The matrix expression specification can be any matrix language expression that evaluates to the value(s) to be written to the matrix-format file.
- You can specify a matrix name, a matrix raised to a power, or a matrix function (with its arguments in parentheses) by itself, but you must enclose other matrix expressions in parentheses. For example, MSAVE A, SAVE INV(A), or MSAVE B\*\*DET(T(C)\*D) is legal, but MSAVE N \* WT is not. You must specify MSAVE (N \* WT).
- Constant expressions are allowed.

## **TYPE Specification**

TYPE specifies the rowtype to write to the matrix-format data file. Only a single rowtype can be written by any one MSAVE statement. Valid keywords on the TYPE specification are:

<b>COV</b>	<i>A matrix of covariances.</i>
<b>CORR</b>	<i>A matrix of correlation coefficients.</i>
<b>MEAN</b>	<i>A vector of means.</i>
<b>STDDEV</b>	<i>A vector of standard deviations.</i>
<b>N</b>	<i>A vector of numbers of cases.</i>
<b>COUNT</b>	<i>A vector of counts.</i>

## **OUTFILE Specification**

OUTFILE designates the matrix-format data file to which the matrices are to be written. It can be an asterisk, an actual filename in quotes, or a file handle defined on a FILE HANDLE command that precedes the matrix program. The filename or handle must be a valid file specification.

- The OUTFILE specification is required on the first MSAVE statement in a matrix program.
- To save a matrix expression as the active dataset (replacing any active dataset created before the matrix program), specify an asterisk (\*).
- Since only one matrix-format data file can be written in a single matrix program, any OUTFILE specification on the second and later MSAVE statements in one matrix program must be the same as that on the first MSAVE statement.

## **VARIABLES Specification**

You can provide variable names for the matrix-format data file with the VARIABLES specification. The variable list is a list of valid variable names separated by commas. You can use the TO convention.

- The VARIABLES specification names only the data variables in the matrix. Split-file variables and grouping or factor variables are named on the SNames and FNames specifications.
- The names in the VARIABLES specification become the values of the special variable VARNAME\_ in the matrix-format data file for rowtypes of CORR and COV.
- You cannot specify the reserved names ROWTYPE\_ and VARNAME\_ on the VARIABLES specification.
- If you omit the VARIABLES specification, the default names COL1, COL2, ..., etc., are used.

## **FACTOR Specification**

To write a matrix-format data file with factor or group codes, you must use the FACTOR specification to provide a row matrix containing the values of each of the factors or group variables for the matrix expression being written by the current MSAVE statement.

- The factor vector must have the same number of columns as there are factors in the matrix data file being written. You can use a scalar when the groups are defined by a single variable. For example, `FACTOR=1` indicates that the matrix data being written are for the value 1 of the factor variable.
- The values of the factor vector are written to the matrix-format data file as values of the factors in the file.
- To create a complete matrix-format data file with factors, you must execute an `MSAVE` statement for every combination of values of the factors or grouping variables (in other words, for every group). If split-file variables are also present, you must execute an `MSAVE` statement for every combination of factor codes within every combination of values of the split-file variables.

### Example

```
MSAVE M11 /TYPE=MEAN /OUTFILE=CORRMAT /VARIABLES=V1 TO V8
      /FNAMES=SEX, GROUP /FACTOR={1,1}.
MSAVE S11 /TYPE STDDEV.
MSAVE MAKE(1,8,N(1,1)) /TYPE N.
MSAVE C11 /TYPE CORR.

MSAVE M12 /TYPE=MEAN /FACTOR={1,2}.
MSAVE S12 /TYPE STDDEV.
MSAVE MAKE(1,8,N(1,2)) /TYPE N.
MSAVE C12 /TYPE CORR.

MSAVE M21 /TYPE=MEAN /FACTOR={2,1}.
MSAVE S21 /TYPE STDDEV.
MSAVE MAKE(1,8,N(2,1)) /TYPE N.
MSAVE C21 /TYPE CORR.

MSAVE M22 /TYPE=MEAN /FACTOR={2,2}.
MSAVE S22 /TYPE STDDEV.
MSAVE MAKE(1,8,N(2,2)) /TYPE N.
MSAVE C22 /TYPE CORR.
```

- The first four `MSAVE` statements provide data for a group defined by the variables *SEX* and *GROUP*, with both factors having the value 1.
- The second, third, and fourth groups of four `MSAVE` statements provide the corresponding data for the other groups, in which *SEX* and *GROUP*, respectively, equal 1 and 2, 2 and 1, and 2 and 2.
- Within each group of `MSAVE` statements, a suitable number-of-cases vector is created with the matrix function `MAKE`.

### ***FNAMES Specification***

To write a matrix-format data file with factor or group codes, you can use the `FNAMES` specification to provide variable names for the grouping or factor variables.

- The variable list following the keyword `FNAMES` is a list of valid variable names, separated by commas.
- If you omit the `FNAMES` specification, the default names *FAC1*, *FAC2*, ..., etc., are used.

### ***SPLIT Specification***

To write a matrix-format data file with split-file groups, you must use the `SPLIT` specification to provide a row matrix containing the values of each of the split-file variables for the matrix expression being written by the current `MSAVE` statement.

- The split vector must have the same number of columns as there are split-file variables in the matrix data file being written. You can use a scalar when there is only one split-file variable. For example, `SPLIT=3` indicates that the matrix data being written are for the value 3 of the split-file variable.
- The values of the split vector are written to the matrix-format data file as values of the split-file variable(s).
- To create a complete matrix-format data file with split-file variables, you must execute `MSAVE` statements for every combination of values of the split-file variables. (If factor variables are present, you must execute `MSAVE` statements for every combination of factor codes within every combination of values of the split-file variables.)

### ***SNAMES Specification***

To write a matrix-format data file with split-file groups, you can use the `SNAMES` specification to provide variable names for the split-file variables.

- The variable list following the keyword `SNAMES` is a list of valid variable names separated by commas.
- If you omit the `SNAMES` specification, the default names *SPL1*, *SPL2*, ..., etc., are used.

### ***DISPLAY Statement***

`DISPLAY` provides information on the matrix variables currently defined in a matrix program and on usage of internal memory by the matrix processor. Two keywords are available on `DISPLAY`:

- |                   |  |
|-------------------|--|
| <b>DICTIONARY</b> | <i>Display variable name and row and column dimensions for each matrix variable currently defined.</i>   |
| <b>STATUS</b>     | <i>Display the status and size of internal tables. This display is intended as a debugging aid when writing large matrix programs that approach the memory limitations of your system.</i> |

If you enter the `DISPLAY` statement with no specifications, both `DICTIONARY` and `STATUS` information is displayed.

## ***RELEASE Statement***

Use the `RELEASE` statement to release the work areas in memory assigned to matrix variables that are no longer needed.

- Specify a list of currently defined matrix variables. Variable names on the list must be separated by commas.
- `RELEASE` discards the contents of the named matrix variables. Releasing a large matrix when it is no longer needed makes memory available for additional matrix variables.
- All matrix variables are released when the `END MATRIX` statement is encountered.

## ***Macros Using the Matrix Language***

Macro expansion (see [DEFINE-!ENDDEFINE on p. 595](#)) occurs before command lines are passed to the matrix processor. Therefore, previously defined macro names can be used within a matrix program. If the macro name expands to one or more valid matrix statements, the matrix processor will execute those statements. Similarly, you can define an entire matrix program, including the `MATRIX` and `END MATRIX` commands, as a macro, but you cannot define a macro within a matrix program, since `DEFINE` and `END DEFINE` are not valid matrix statements.



# MATRIX DATA

```
MATRIX DATA VARIABLES=varlist  [/FILE={INLINE**}]
                               {file  }

[/FORMAT={ LIST** }  [ LOWER** ]  [ DIAGONAL** ] ]
          {FREE  }    {UPPER  }    {NODIAGONAL}
          {FULL  }

[/SPLIT=varlist]  [/FACTORS=varlist]

[/CELLS=number of cells]  [/N=sample size]

[/CONTENTS= CORR**  [COV]  [MAT]  [MSE]  [DFE]  [MEAN]  [PROX]
            [ STDDEV ]  [N_SCALAR]  [ N_VECTOR ]  [N_MATRIX]  [COUNT] ]
            {SD  }          {N  }          }
```

**\*\***Default if the subcommand is omitted.

## Example

```
MATRIX DATA VARIABLES=ROWTYPE_ SAVINGS POP15 POP75 INCOME GROWTH.
BEGIN DATA
MEAN 9.6710 35.0896 2.2930 1106.7784 3.7576
STDDEV 4.4804 9.1517 1.2907 990.8511 2.8699
N 50 50 50 50 50
CORR 1
CORR -.4555 1
CORR .3165 -.9085 1
CORR .2203 -.7562 .7870 1
CORR .3048 -.0478 .0253 -.1295 1
END DATA.
```

## Overview

MATRIX DATA reads raw matrix materials and converts them to a matrix data file that can be read by procedures that handle matrix materials. The data can include vector statistics, such as means and standard deviations, as well as matrices.

MATRIX DATA is similar to a DATA LIST command: it defines variable names and their order in a raw data file. However, MATRIX DATA can read only data that conform to the general format of matrix data files.

### Matrix Files

Like the matrix data files created by procedures, the file that MATRIX DATA creates contains the following variables in the indicated order. If the variables are in a different order in the raw data file, MATRIX DATA rearranges them in the active dataset.

- *Split-file variables.* These optional variables define split files. There can be up to eight split variables, and they must have numeric values. Split-file variables will appear in the order in which they are specified on the SPLIT subcommand.

- *ROWTYPE\_*. This is a string variable with A8 format. Its values define the data type for each record. For example, it might identify a row of values as means, standard deviations, or correlation coefficients. Every matrix data file has a *ROWTYPE\_* variable.
- *Factor variables*. There can be any number of factors. They occur only if the data include within-cells information, such as the within-cells means. Factors have the system-missing value on records that define pooled information. Factor variables appear in the order in which they are specified on the FACTORS subcommand.
- *VARNAME\_*. This is a string variable with A8 format. MATRIX DATA automatically generates *VARNAME\_* and its values based on the variables named on VARIABLES. You never enter values for *VARNAME\_*. Values for *VARNAME\_* are blank for records that define vector information. Every matrix in the program has a *VARNAME\_* variable.
- *Continuous variables*. These are the variables that were used to generate the correlation coefficients or other aggregated data. There can be any number of them. Continuous variables appear in the order in which they are specified on VARIABLES.

### **Options**

**Data Files.** You can define both inline data and data in an external file.

**Data Format.** By default, data are assumed to be entered in freefield format with each vector or row beginning on a new record (the keyword LIST on the FORMAT subcommand). If each vector or row does not begin on a new record, use the keyword FREE. You can also use FORMAT to indicate whether matrices are entered in upper or lower triangular or full square or rectangular format and whether or not they include diagonal values.

**Variable Types.** You can specify split-file and factor variables using the SPLIT and FACTORS subcommands. You can identify record types by specifying ROWTYPE\_ on the VARIABLES subcommand if ROWTYPE\_ values are included in the data or by implying ROWTYPE\_ values on CONTENTS.

### **Basic Specification**

The basic specification is VARIABLES and a list of variables. Additional specifications are required as follows:

- FILE is required to specify the data file if the data are not inline.
- If data are in any format other than lower triangular with diagonal values included, FORMAT is required.
- If the data contain values in addition to matrix coefficients, such as the mean and standard deviation, either the variable ROWTYPE\_ must be specified on VARIABLES and ROWTYPE\_ values must be included in the data or CONTENTS must be used to describe the data.
- If the data include split-file variables, SPLIT is required. If there are factors, FACTORS is required.

Specifications on most MATRIX DATA subcommands depend on whether ROWTYPE\_ is included in the data and specified on VARIABLES or whether it is implied using CONTENTS.

Table 134-1  
Subcommand requirements in relation to ROWTYPE\_

Subcommand	Implicit ROWTYPE_ using CONTENTS	Explicit ROWTYPE_ on VARIABLES
FILE	Defaults to INLINE	Defaults to INLINE
VARIABLES	Required	Required
FORMAT	Defaults to LOWER DIAG	Defaults to LOWER DIAG
SPLIT	Required if split files*	Required if split files
FACTORS	Required if factors	Required if factors
CELLS	Required if factors	Inapplicable
CONTENTS	Defaults to CORR	Optional
N	Optional	Optional

\* If the data do not contain values for the split-file variables, this subcommand can specify a single variable, which is not specified on the VARIABLES subcommand.

### Subcommand Order

- SPLIT and FACTORS, when used, must follow VARIABLES.
- The remaining subcommands can be specified in any order.

### Syntax Rules

- No commands can be specified between MATRIX DATA and BEGIN DATA, not even a VARIABLE LABELS or FORMAT command. Data transformations cannot be used until after MATRIX DATA is executed.

## Examples

### Reading a Correlation Matrix

```
MATRIX DATA
  VARIABLES=ROWTYPE_ SAVINGS POP15 POP75 INCOME GROWTH.
BEGIN DATA
MEAN 9.6710 35.0896 2.2930 1106.7784 3.7576
STDDEV 4.4804 9.1517 1.2907 990.8511 2.8699
N 50 50 50 50 50
CORR 1
CORR -.4555 1
CORR .3165 -.9085 1
CORR .2203 -.7562 .7870 1
CORR .3048 -.0478 .0253 -.1295 1
END DATA.
```

- The variable ROWTYPE\_ is specified on VARIABLES. ROWTYPE\_ values are included in the data.
- No other specifications are required.

### MATRIX DATA with DISCRIMINANT

```
MATRIX DATA VARIABLES=WORLD ROWTYPE_ FOOD APPL SERVICE RENT
/FACTORS=WORLD.
```

## MATRIX DATA

```

BEGIN DATA
1 N      25 25 25 25
1 MEAN   76.64 77.32 81.52 101.40
2 N      7 7 7 7
2 MEAN   76.1428571 85.2857143 60.8571429 249.571429
3 N      13 13 13 13
3 MEAN   55.5384615 76 63.4615385 86.3076923
. SD     16.4634139 22.5509310 16.8086768 77.1085326
. CORR   1
. CORR   .1425366 1
. CORR   .5644693 .2762615 1
. CORR   .2133413 -.0499003 .0417468 1
END DATA.

DISCRIMINANT GROUPS=WORLD(1,3)
/VARIABLES=FOOD APPL SERVICE RENT /METHOD=WILKS /MATRIX=IN(*).

```

- MATRIX DATA is used to generate a active dataset that DISCRIMINANT can read. DISCRIMINANT reads the mean, count (unweighted  $N$ ), and  $N$  (weighted  $N$ ) for each cell in the data, as well as the pooled values for the standard deviation and correlation coefficients. If count equals  $N$ , only  $N$  needs to be supplied.
- ROWTYPE\_ is specified on VARIABLES to identify record types in the data. Though CONTENTS and CELLS can be used to identify record types and distinguish between within-cells data and pooled values, it is usually easier to specify ROWTYPE\_ on VARIABLES and enter the ROWTYPE\_ values in the data.
- Because factors are present in the data, the continuous variables (*FOOD*, *APPL*, *SERVICE*, and *RENT*) must be specified last on VARIABLES and must be last in the data.
- The FACTORS subcommand identifies *WORLD* as the factor variable.
- BEGIN DATA immediately follows MATRIX DATA.
- $N$  and MEAN values for each cell are entered in the data.
- ROWTYPE\_ values for the pooled records are SD and COR. MATRIX DATA assigns the values STDDEV and CORR to the corresponding vectors in the matrix. Records with pooled information have the system-missing value (.) for the factors.
- The DISCRIMINANT procedure reads the data matrix. An asterisk (\*) is specified as the input file on the MATRIX subcommand because the data are in the active dataset.

**MATRIX DATA with REGRESSION**

```

MATRIX DATA VARIABLES=SAVINGS POP15 POP75 INCOME GROWTH
/CONTENTS=MEAN SD N CORR /FORMAT=UPPER NODIAGONAL.

BEGIN DATA
9.6710 35.0896 2.2930 1106.7784 3.7576
4.4804 9.1517 1.2908 990.8511 2.8699
50 50 50 50 50
-.4555 .3165 .2203 .3048
-.9085 -.7562 -.0478
.7870 .0253
-.1295
END DATA.

REGRESSION MATRIX=IN(*) /VARIABLES=SAVINGS TO GROWTH
/DEP=SAVINGS /ENTER.

```

- `MATRIX DATA` is used to generate a matrix that `REGRESSION` can read. `REGRESSION` reads and writes matrices that always contain the mean, standard deviation,  $N$ , and Pearson correlation coefficients. Data in this example do not have `ROWTYPE_` values, and the correlation values are from the upper triangle of the matrix without the diagonal values.
- `ROWTYPE_` is not specified on `VARIABLES` because its values are not included in the data.
- Because there are no `ROWTYPE_` values, `CONTENTS` is required to define the record types and the order of the records in the file.
- By default, `MATRIX DATA` reads values from the lower triangle of the matrix, including the diagonal values. `FORMAT` is required in this example to indicate that the data are in the upper triangle and do not include diagonal values.
- `BEGIN DATA` immediately follows the `MATRIX DATA` command.
- The `REGRESSION` procedure reads the data matrix. An asterisk (\*) is specified as the input file on the `MATRIX` subcommand because the data are in the active dataset. Since there is a single vector of  $N$ 's in the data, missing values are handled listwise (the default for `REGRESSION`).

### **MATRIX DATA with ONEWAY**

```
MATRIX DATA VARIABLES=EDUC ROWTYPE_ WELL /FACTORS=EDUC.
BEGIN DATA
1 N 65
2 N 95
3 N 181
4 N 82
5 N 40
6 N 37
1 MEAN 2.6462
2 MEAN 2.7737
3 MEAN 4.1796
4 MEAN 4.5610
5 MEAN 4.6625
6 MEAN 5.2297
. MSE 6.2699
. DFE 494
END DATA.

ONEWAY WELL BY EDUC (1,6) /MATRIX=IN(*)
```

- One of the two types of matrices that the `ONEWAY` procedure reads includes a vector of frequencies for each factor level, a vector of means for each factor level, a record containing the pooled variance (within-group mean square error), and the degrees of freedom for the mean square error. `MATRIX DATA` is used to generate an active dataset containing this type of matrix data for the `ONEWAY` procedure.
- `ROWTYPE_` is explicit on `VARIABLES` and identifies record types.
- Because factors are present in the data, the continuous variables (`WELL`) must be specified last on `VARIABLES` and must be last in the data.
- The `FACTORS` subcommand identifies `EDUC` as the factor variable.
- `MSE` is entered in the data as the `ROWTYPE_` value for the vector of square pooled standard deviations.
- `DFE` is entered in the data as the `ROWTYPE_` value for the vector of degrees of freedom.
- Records with pooled information have the system-missing value (.) for the factors.

## Operations

- `MATRIX DATA` defines and writes data in one step.
- `MATRIX DATA` clears the active dataset and defines a new active dataset.
- If `CONTENTS` is not specified and `ROWTYPE_` is not specified on `VARIABLES`, `MATRIX DATA` assumes that the data contain only `CORR` values and issues warning messages to alert you to its assumptions.
- With the default format, data values, including diagonal values, must be in the lower triangle of the matrix. If `MATRIX DATA` encounters values in the upper triangle, it ignores those values and issues a series of warnings.
- With the default format, if any matrix rows span records in the data file, `MATRIX DATA` cannot form the matrix properly.
- `MATRIX DATA` does not allow format specifications for matrix materials. The procedure assigns the formats shown in the following table. To change data formats, execute `MATRIX DATA` and then assign new formats with the `FORMATS`, `PRINT FORMATS`, or `WRITE FORMATS` command.

Table 134-2  
Print and write formats for matrix variables

Variable type	Format
<code>ROWTYPE_</code> , <code>VARNAME_</code>	A8
Split-file variables	F4.0
Factors	F4.0
Continuous variables	F10.4

## Format of the Raw Matrix Data File

- If `LIST` is in effect on the `FORMAT` subcommand, the data are entered in freefield format, with blanks and commas used as separators and each scalar, vector, or row of the matrix beginning on a new record. Unlike `LIST` format with `DATA LIST`, a vector or row of the matrix can be contained on multiple records. The continuation records do not have a value for `ROWTYPE_`.
- `ROWTYPE_` values can be enclosed in quotes.
- The order of variables in the raw data file must match the order in which they are specified on `VARIABLES`. However, this order does not have to correspond to the order of variables in the resulting matrix data file.
- The way records are entered for pooled vectors or matrices when factors are present depends upon whether `ROWTYPE_` is specified on the `VARIABLES` subcommand. [For more information, see the topic FACTORS Subcommand on p. 1182.](#)
- `MATRIX DATA` recognizes plus and minus signs as field separators when they are not preceded by the letter *D* or *E*. This allows `MATRIX DATA` to read scientific notation as well as correlation matrices written by FORTRAN in F10.8 format. A plus sign preceded by a *D* or *E* is read as part of the number in scientific notation.

## VARIABLES Subcommand

`VARIABLES` specifies the names of the variables in the raw data and the order in which they occur.

- `VARIABLES` is required.
- There is no limit to the number of variables that can be specified.
- If `ROWTYPE_` is specified on `VARIABLES`, the continuous variables must be the last variables specified on the subcommand and must be last in the data.
- If split-file variables are present, they must also be specified on `SPLIT`.
- If factor variables are present, they must also be specified on `FACTORS`.

When either of the following is true, the only variables that must be specified on `VARIABLES` are the continuous variables:

1. The data contain only correlation coefficients. There can be no additional information, such as the mean and standard deviation, and no factor information or split-file variables. `MATRIX DATA` assigns the record type `CORR` to all records.
2. `CONTENTS` is used to define all record types. The data can then contain information such as the mean and standard deviation, but no factor, split-file, or `ROWTYPE_` variables. `MATRIX DATA` assigns the record types defined on the `CONTENTS` subcommand.

### Variable `VARNAME_`

`VARNAME_` cannot be specified on the `VARIABLES` subcommand or anywhere on `MATRIX DATA`, and its values cannot be included in the data. The `MATRIX DATA` command generates the variable `VARNAME_` automatically.

### Variable `ROWTYPE_`

- `ROWTYPE_` is a string variable with `A8` format. Its values define the data types. All matrix data files contain a `ROWTYPE_` variable.
- If `ROWTYPE_` is specified on `VARIABLES` and its values are entered in the data, `MATRIX DATA` is primarily used to define the names and order of the variables in the raw data file.
- `ROWTYPE_` must precede the continuous variables.
- Valid values for `ROWTYPE_` are `CORR`, `COV`, `MAT`, `MSE`, `DFE`, `MEAN`, `STDDEV` (or `SD`), `N_VECTOR` (or `N`), `N_SCALAR`, `N_MATRIX`, `COUNT`, or `PROX`. For definitions of these values. [For more information, see the topic CONTENTS Subcommand on p. 1184.](#) Three-character abbreviations for these values are permitted. These values can also be enclosed in quotation marks or apostrophes.
- If `ROWTYPE_` is not specified on `VARIABLES`, `CONTENTS` must be used to define the order in which the records occur within the file. `MATRIX DATA` follows these specifications strictly and generates a `ROWTYPE_` variable according to the `CONTENTS` specifications. A data-entry error, especially skipping a record, can cause the procedure to assign the wrong values to the wrong records.

### Example

```
* ROWTYPE_ is specified on VARIABLES.
MATRIX DATA
```

## MATRIX DATA

```

      VARIABLES=ROWTYPE_ SAVINGS POP15 POP75 INCOME GROWTH.
BEGIN DATA
MEAN 9.6710 35.0896 2.2930 1106.7784 3.7576
STDDEV 4.4804 9.1517 1.2907 990.8511 2.8699
N 50 50 50 50 50
CORR 1
CORR -.4555 1
CORR .3165 -.9085 1
CORR .2203 -.7562 .7870 1
CORR .3048 -.0478 .0253 -.1295 1
END DATA.

```

- *ROWTYPE\_* is specified on *VARIABLES*. *ROWTYPE\_* values in the data identify each record type.
- Note that *VARNAME\_* is not specified on *VARIABLES*, and its values are not entered in the data.

**Example**

```

* ROWTYPE_ is specified on VARIABLES.

MATRIX DATA
      VARIABLES=ROWTYPE_ SAVINGS POP15 POP75 INCOME GROWTH.
BEGIN DATA
'MEAN ' 9.6710 35.0896 2.2930 1106.7784 3.7576
'SD ' 4.4804 9.1517 1.2907 990.8511 2.8699
'N ' 50 50 50 50 50
"CORR " 1
"CORR " -.4555 1
"CORR " .3165 -.9085 1
"CORR " .2203 -.7562 .7870 1
"CORR " .3048 -.0478 .0253 -.1295 1
END DATA.

```

- *ROWTYPE\_* values for the mean, standard deviation, *N*, and Pearson correlation coefficients are abbreviated and enclosed in quotes.

**Example**

```

* ROWTYPE_ is not specified on VARIABLES.

MATRIX DATA VARIABLES=SAVINGS POP15 POP75 INCOME GROWTH
  /CONTENTS=MEAN SD N CORR.
BEGIN DATA
9.6710 35.0896 2.2930 1106.7784 3.7576
4.4804 9.1517 1.2907 990.8511 2.8699
50 50 50 50 50
1
-.4555 1
.3165 -.9085 1
.2203 -.7562 .7870 1
.3048 -.0478 .0253 -.1295 1
END DATA.

```

- *ROWTYPE\_* is not specified on *VARIABLES*, and its values are not included in the data.
- *CONTENTS* is required to define the record types and the order of the records in the file.



## **FILE Subcommand**

FILE specifies the matrix file containing the data. The default specification is `INLINE`, which indicates that the data are included within the command sequence between the `BEGIN DATA` and `END DATA` commands.

- If the data are in an external file, FILE must specify the file.
- If the FILE subcommand is omitted, the data must be inline.

### **Example**

```
MATRIX DATA FILE=RAWMTX /VARIABLES=varlist.
```

- FILE indicates that the data are in the file *RAWMTX*.

## **FORMAT Subcommand**

FORMAT indicates how the matrix data are formatted. It applies only to matrix values in the data, not to vector values, such as the mean and standard deviation.

- FORMAT can specify up to three keywords: one to specify the data-entry format, one to specify matrix shape, and one to specify whether the data include diagonal values.
- The minimum specification is a single keyword.
- Default settings remain in effect unless explicitly overridden.

### **Data-Entry Format**

FORMAT has two keywords that specify the data-entry format:

<b>LIST</b>	<i>Each scalar, vector, and matrix row must begin on a new record. A vector or row of the matrix may be continued on multiple records. This is the default.</i>
<b>FREE</b>	<i>Matrix rows do not need to begin on a new record. Any item can begin in the middle of a record.</i>

### **Matrix Shape**

FORMAT has three keywords that specify the matrix shape. With either triangular shape, no values—not even missing indicators—are entered for the implied values in the matrix.

<b>LOWER</b>	<i>Read data values from the lower triangle. This is the default.</i>
<b>UPPER</b>	<i>Read data values from the upper triangle.</i>
<b>FULL</b>	<i>Read the full square matrix of data values. FULL cannot be specified with NODIAGONAL.</i>

## Diagonal Values

FORMAT has two keywords that refer to the diagonal values:

<b>DIAGONAL</b>	<i>Data include the diagonal values.</i> This is the default.
<b>NODIAGONAL</b>	<i>Data do not include diagonal values.</i> The diagonal value is set to the system-missing value for all matrices except the correlation matrices. For correlation matrices, the diagonal value is set to 1. NODIAGONAL cannot be specified with FULL.

The following table shows how data might be entered for each combination of FORMAT settings that govern matrix shape and diagonal values. With UPPER NODIAGONAL and LOWER NODIAGONAL, you do not enter the matrix row that has blank values for the continuous variables. If you enter that row, MATRIX DATA cannot properly form the matrix.

Table 134-3  
Various FORMAT settings

FULL	UPPER DIAGONAL	UPPER NODIAGONAL	LOWER DIAGONAL	LOWER NODIAGONAL
MEAN 5 4 3	MEAN 5 4 3	MEAN 5 4 3	MEAN 5 4 3	MEAN 5 4 3
SD 3 2 1	SD 3 2 1	SD 3 2 1	SD 3 2 1	SD 3 2 1
N 9 9 9	N 9 9 9	N 9 9 9	N 9 9 9	N 9 9 9
CORR 1 .6 .7	CORR 1 .6 .7	CORR .6 .7	CORR 1	CORR .6
CORR .6 1 .8	CORR 1 .8	CORR .8	CORR .6 1	CORR .7 .8
CORR .7 .8 1	CORR 1		CORR .7 .8 1	

### Example

```
MATRIX DATA VARIABLES=ROWTYPE_ V1 TO V3
  /FORMAT=UPPER NODIAGONAL.
BEGIN DATA
MEAN  5  4  3
SD    3  2  1
N     9  9  9
CORR  .6 .7
CORR  .8
END DATA.
LIST.
```

- FORMAT specifies the upper-triangle format with no diagonal values. The default LIST is in effect for the data-entry format.

### Example

```
MATRIX DATA VARIABLES=ROWTYPE_ V1 TO V3
  /FORMAT=UPPER NODIAGONAL.
BEGIN DATA
MEAN 5 4 3
SD 3 2 1
N 9 9 9
CORR .6 .7
CORR .8
END DATA.
```

LIST.

- This example is identical to the previous example. It shows that data do not have to be aligned in columns. Data throughout this section are aligned in columns to emphasize the matrix format.

## ***SPLIT Subcommand***

**SPLIT** specifies the variables whose values define the split files. **SPLIT** must follow the **VARIABLES** subcommand.

- **SPLIT** can specify a subset of up to eight of the variables named on **VARIABLES**. All split variables must be numeric. The keyword **TO** can be used to imply variables in the order in which they are named on **VARIABLES**.
- A separate matrix must be included in the data for each value of each split variable. **MATRIX DATA** generates a complete set of matrix materials for each.
- If the data contain neither **ROWTYPE\_** nor split-file variables, a single split-file variable can be specified on **SPLIT**. This variable is *not* specified on the **VARIABLES** subcommand. **MATRIX DATA** generates a complete set of matrix materials for each set of matrix materials in the data and assigns values 1, 2, 3, etc., to the split variable until the end of the data is encountered.

### ***Example***

```
MATRIX DATA VARIABLES=S1 ROWTYPE_ V1 TO V3 /SPLIT=S1.
BEGIN DATA
0 MEAN 5 4 3
0 SD 1 2 3
0 N 9 9 9
0 CORR 1
0 CORR .6 1
0 CORR .7 .8 1
1 MEAN 9 8 7
1 SD 5 6 7
1 N 9 9 9
1 CORR 1
1 CORR .4 1
1 CORR .3 .2 1
END DATA.
LIST.
```

- The split variable *S1* has two values: 0 and 1. Two separate matrices are entered in the data, one for each value *S1*.
- *S1* must be specified on both **VARIABLES** and **SPLIT**.

### ***Example***

```
MATRIX DATA VARIABLES=V1 TO V3 /CONTENTS=MEAN SD N CORR
/SPLIT=SPL.
BEGIN DATA
5 4 3
1 2 3
9 9 9
1
.6 1
.7 .8 1
```

## MATRIX DATA

```

9 8 7
5 6 7
9 9 9
1
.4 1
.3 .2 1
END DATA.
LIST.

```

- The split variable *SPL* is not specified on *VARIABLES*, and values for *SPL* are not included in the data.
- Two sets of matrix materials are included in the data. *MATRIX DATA* therefore assigns values 1 and 2 to variable *SPL* and generates two matrices in the matrix data file.

## FACTORS Subcommand

*FACTORS* specifies the variables whose values define the cells represented by the within-cells data. *FACTORS* must follow the *VARIABLES* subcommand.

- *FACTORS* specifies a subset of the variables named on the *VARIABLES* subcommand. The keyword *TO* can be used to imply variables in the order in which they are named on *VARIABLES*.
- If *ROWTYPE\_* is explicit on *VARIABLES* and its values are included in the data, records that represent pooled information have the system-missing value (indicated by a period) for the factors, since the values of *ROWTYPE\_* are ambiguous.
- If *ROWTYPE\_* is not specified on *VARIABLES* and its values are not in the data, enter data values for the factors only for records that represent within-cells information. Enter nothing for the factors for records that represent pooled information. *CELLS* must be specified to indicate the number of within-cells records, and *CONTENTS* must be specified to indicate which record types have within-cells data.

### Example

```

* Rowtype is explicit.

MATRIX DATA VARIABLES=ROWTYPE_ F1 F2  VAR1 TO VAR3
  /FACTORS=F1 F2.
BEGIN DATA
MEAN 1 1 1 2 3
SD 1 1 5 4 3
N 1 1 9 9 9
MEAN 1 2 4 5 6
SD 1 2 6 5 4
N 1 2 9 9 9
MEAN 2 1 7 8 9
SD 2 1 7 6 5
N 2 1 9 9 9
MEAN 2 2 9 8 7
SD 2 2 8 7 6
N 2 2 9 9 9
CORR . . .1
CORR . . .6 1
CORR . . .7 .8 1
END DATA.

```

- *ROWTYPE\_* is specified on *VARIABLES*.

- Factor variables must be specified on both `VARIABLES` and `FACTORS`.
- Periods in the data represent missing values for the `CORR` factor values.

### Example

```
* Rowtype is implicit.
MATRIX DATA VARIABLES=F1 F2 VAR1 TO VAR3
  /FACTORS=F1 F2 /CONTENTS=(MEAN SD N) CORR /CELLS=4.
BEGIN DATA
1 1 1 2 3
1 1 5 4 3
1 1 9 9 9
1 2 4 5 6
1 2 6 5 4
1 2 9 9 9
2 1 7 8 9
2 1 7 6 5
2 1 9 9 9
2 2 9 8 7
2 2 8 7 6
2 2 9 9 9
      1
      .6 1
      .7 .8 1
END DATA.
```

- `ROWTYPE_` is not specified on `VARIABLES`.
- Nothing is entered for the `CORR` factor values because the records contain pooled information.
- `CELLS` is required because there are factors in the data and `ROWTYPE_` is implicit.
- `CONTENTS` is required to define the record types and to differentiate between the within-cells and pooled types.

## CELLS Subcommand

`CELLS` specifies the number of within-cells records in the data. The only valid specification for `CELLS` is a single integer, which indicates the number of sets of within-cells information that `MATRIX DATA` must read.

- `CELLS` is required when there are factors in the data and `ROWTYPE_` is implicit.
- If `CELLS` is used when `ROWTYPE_` is specified on `VARIABLES`, `MATRIX DATA` issues a warning and ignores the `CELLS` subcommand.

### Example

```
MATRIX DATA VARIABLES=F1 VAR1 TO VAR3 /FACTORS=F1 /CELLS=2
  /CONTENTS=(MEAN SD N) CORR.
BEGIN DATA
1 5 4 3
1 3 2 1
1 9 9 9
2 8 7 6
2 6 7 8
2 9 9 9
      1
      .6 1
```

```
.7 .8 1
END DATA.
```

- The specification for `CELLS` is 2 because the factor variable `FI` has two values (1 and 2) and there are therefore two sets of within-cells information.
- If there were two factor variables, `FI` and `F2`, and each had two values, 1 and 2, `CELLS` would equal 4 to account for all four possible factor combinations (assuming all that 4 combinations are present in the data).

## CONTENTS Subcommand

`CONTENTS` defines the record types when `ROWTYPE_` is not included in the data. The minimum specification is a single keyword indicating a type of record. The default is `CORR`.

- `CONTENTS` is required to define record types and record order whenever `ROWTYPE_` is not specified on `VARIABLES` and its values are not in the data. The only exception to this rule is the rare situation in which all data values represent pooled correlation records and there are no factors. In that case, `MATRIX DATA` reads the data values and assigns the default `ROWTYPE_` of `CORR` to all records.
- The order in which keywords are specified on `CONTENTS` must correspond to the order in which records appear in the data. If the keywords on `CONTENTS` are in the wrong order, `MATRIX DATA` will incorrectly assign values.

<b>CORR</b>	<i>Matrix of correlation coefficients.</i> This is the default. If <code>ROWTYPE_</code> is not specified on the <code>VARIABLES</code> subcommand and you omit the <code>CONTENTS</code> subcommand, <code>MATRIX DATA</code> assigns the <code>ROWTYPE_</code> value <code>CORR</code> to all matrix rows.
<b>COV</b>	<i>Matrix of covariance coefficients.</i>
<b>MAT</b>	<i>Generic square matrix.</i>
<b>MSE</b>	<i>Vector of mean squared errors.</i>
<b>DFE</b>	<i>Vector of degrees of freedom.</i>
<b>MEAN</b>	<i>Vector of means.</i>
<b>STDDEV</b>	<i>Vector of standard deviations.</i> <code>SD</code> is a synonym for <code>STDDEV</code> . <code>MATRIX DATA</code> assigns the <code>ROWTYPE_</code> value <code>STDDEV</code> to the record if either <code>STDDEV</code> or <code>SD</code> is specified.
<b>N_VECTOR</b>	<i>Vector of counts.</i> <code>N</code> is a synonym for <code>N_VECTOR</code> . <code>MATRIX DATA</code> assigns the <code>ROWTYPE_</code> value <code>N</code> to the record.
<b>N_SCALAR</b>	<i>Count.</i> Scalars are a shorthand mechanism for representing vectors in which all elements have the same value, such as when a vector of <code>N</code> 's is calculated using listwise deletion of missing values. Enter <code>N_SCALAR</code> as the <code>ROWTYPE_</code> value in the data and then the <code>N_SCALAR</code> value for the first continuous variable only. <code>MATRIX DATA</code> assigns the <code>ROWTYPE_</code> value <code>N</code> to the record and copies the specified <code>N_SCALAR</code> value across all of the continuous variables.
<b>N_MATRIX</b>	<i>Square matrix of counts.</i> Enter <code>N_MATRIX</code> as the <code>ROWTYPE_</code> value for each row of counts in the data. <code>MATRIX DATA</code> assigns the <code>ROWTYPE_</code> value <code>N</code> to each of those rows.

<b>COUNT</b>	Count vector accepted by procedure <i>DISCRIMINANT</i> . This contains unweighted <i>N</i> 's.
<b>PROX</b>	Matrix produced by <i>PROXIMITIES</i> . Any proximity matrix can be used with <i>PROXIMITIES</i> or <i>CLUSTER</i> . A value label of <i>SIMILARITY</i> or <i>DISSIMILARITY</i> should be specified for <i>PROX</i> by using the <i>VALUE LABELS</i> command after <i>END DATA</i> .

**Example**

```
MATRIX DATA VARIABLES=V1 TO V3 /CONTENTS=MEAN SD N_SCALAR CORR.
BEGIN DATA
  5  4  3
  3  2  1
  9
  1
  .6  1
  .7 .8  1
END DATA.
LIST.
```

- *ROWTYPE\_* is not specified on *VARIABLES*, and *ROWTYPE\_* values are not in the data. *CONTENTS* is therefore required to identify record types.
- *CONTENTS* indicates that the matrix records are in the following order: mean, standard deviation, *N*, and correlation coefficients.
- The *N\_SCALAR* value is entered for the first continuous variable only.

**Example**

```
MATRIX DATA VARIABLES=V1 TO V3 /CONTENTS=PROX.
BEGIN DATA

data records

END DATA.
VALUE LABELS ROWTYPE_ 'PROX' 'DISSIMILARITY'.
```

- *CONTENTS* specifies *PROX* to read a raw matrix and create a matrix data file in the same format as one produced by procedure *PROXIMITIES*. *PROX* is assigned the value label *DISSIMILARITY*.

**Within-Cells Record Definition**

When the data include factors and *ROWTYPE\_* is not specified, *CONTENTS* distinguishes between within-cells and pooled records by enclosing the keywords for within-cells records in parentheses.

- If the records associated with the within-cells keywords appear together for each set of factor values, enclose the keywords together within a single set of parentheses.
- If the records associated with each within-cells keyword are grouped together across factor values, enclose the keyword within its own parentheses.

**Example**

```
MATRIX DATA VARIABLES=F1 VAR1 TO VAR3 /FACTORS=F1 /CELLS=2
/CONTENTS=(MEAN SD N) CORR.
```

- MEAN, SD, and N contain within-cells information and are therefore specified within parentheses. CORR is outside the parentheses because it identifies pooled records.
- CELLS is required because there is a factor specified and *ROWTYPE\_* is implicit.

**Example**

```
MATRIX DATA VARIABLES=F1 VAR1 TO VAR3 /FACTORS=F1 /CELLS=2
/CONTENTS=(MEAN SD N) CORR.
BEGIN DATA
1 5 4 3
1 3 2 1
1 9 9 9
2 4 5 6
2 6 5 4
2 9 9 9
1
.6 1
.7 .8 1
END DATA.
```

- The parentheses around the CONTENTS keywords indicate that the mean, standard deviation, and *N* for value 1 of factor *F1* are together, followed by the mean, standard deviation, and *N* for value 2 of factor *F1*.

**Example**

```
MATRIX DATA VARIABLES=F1 VAR1 TO VAR3 /FACTORS=F1 /CELLS=2
/CONTENTS=(MEAN) (SD) (N) CORR.
BEGIN DATA
1 5 4 3
2 4 5 6
1 3 2 1
2 6 5 4
1 9 9 9
2 9 9 9
1
.6 1
.7 .8 1
END DATA.
```

- The parentheses around each CONTENTS keyword indicate that the data include the means for all cells, followed by the standard deviations for all cells, followed by the *N* values for all cells.

**Example**

```
MATRIX DATA VARIABLES=F1 VAR1 TO VAR3 /FACTORS=F1 /CELLS=2
/CONTENTS=(MEAN SD) (N) CORR.
BEGIN DATA
1 5 4 3
1 3 2 1
2 4 5 6
2 6 5 4
1 9 9 9
2 9 9 9
1
.6 1
.7 .8 1
```



END DATA.

- The parentheses around the CONTENTS keywords indicate that the data include the mean and standard deviation for value 1 of *F1*, followed by the mean and standard deviation for value 2 of *F1*, followed by the *N* values for all cells.

### Optional Specification When ROWTYPE\_ Is Explicit

When *ROWTYPE\_* is explicitly named on VARIABLES, MATRIX DATA uses *ROWTYPE\_* values to determine record types.

- When *ROWTYPE\_* is explicitly named on VARIABLES, CONTENTS can be used for informational purposes. However, *ROWTYPE\_* values in the data determine record types.
- If MATRIX DATA reads values for *ROWTYPE\_* that are not specified on CONTENTS, it issues a warning.
- Missing values for factors are entered as periods, even though CONTENTS is specified. [For more information, see the topic FACTORS Subcommand on p. 1182.](#)

#### Example

```
MATRIX DATA VARIABLES=ROWTYPE_ F1 F2 VAR1 TO VAR3
  /FACTORS=F1 F2 /CONTENTS=(MEAN SD N) CORR.
BEGIN DATA
MEAN 1 1 1 2 3
SD 1 1 5 4 3
N 1 1 9 9 9
MEAN 1 2 4 5 6
SD 1 2 6 5 4
N 1 2 9 9 9
CORR . . 1
CORR . . .6 1
CORR . . .7 .8 1
END DATA.
```

- *ROWTYPE\_* is specified on VARIABLES. MATRIX DATA therefore uses *ROWTYPE\_* values in the data to identify record types.
- Because *ROWTYPE\_* is specified on VARIABLES, CONTENTS is optional. However, CONTENTS is specified for informational purposes. This is most useful when data are in an external file and the *ROWTYPE\_* values cannot be seen in the data.
- Missing values for factors are entered as periods, even though CONTENTS is specified.

## N Subcommand

*N* specifies the population *N* when the data do not include it. The only valid specification is an integer, which indicates the population *N*.

- MATRIX DATA generates one record with a *ROWTYPE\_* of *N* for each split file, and it uses the specified *N* value for each continuous variable.

#### Example

```
MATRIX DATA VARIABLES=V1 TO V3 /CONTENTS=MEAN SD CORR
```

---

*MATRIX DATA*

```
      /N=99.  
BEGIN DATA  
  5  4  3  
  3  4  5  
  1  
  .6  1  
  .7  .8  1  
END DATA.
```

- *MATRIX DATA* uses 99 as the *N* value for all continuous variables.

# MCONVERT

```
MCONVERT [ [/MATRIX=] [IN({* })] [OUT({* })]]  
          {file}          {file}  
          [ {/REPLACE} ]  
          [ {/APPEND } ]
```

## Example

```
MCONVERT MATRIX=OUT(CORMTX) /APPEND.
```

## Overview

MCONVERT converts covariance matrix materials to correlation matrix materials, or vice versa. For MCONVERT to convert a correlation matrix, the matrix data must contain *CORR* values (Pearson correlation coefficients) and a vector of standard deviations (*STDDEV*). For MCONVERT to convert a covariance matrix, only *COV* values are required in the data.

## Options

**Matrix Files.** MCONVERT can read matrix materials from an external matrix data file, and it can write converted matrix materials to an external file.

**Matrix Materials.** MCONVERT can write the converted matrix only or both the converted matrix and the original matrix to the resulting matrix data file.

## Basic Specification

The minimum specification is the command itself. By default, MCONVERT reads the original matrix from the active dataset and then replaces it with the converted matrix.

## Syntax Rules

- The keywords *IN* and *OUT* cannot specify the same external file.
- The *APPEND* and *REPLACE* subcommands cannot be specified on the same MCONVERT command.

## Operations

- If the data are covariance matrix materials, MCONVERT converts them to a correlation matrix plus a vector of standard deviations.
- If the data are a correlation matrix and vector of standard deviations, MCONVERT converts them to a covariance matrix.
- If there are multiple *CORR* or *COV* matrices (for example, one for each grouping (factor) or one for each split variable), each will be converted to a separate matrix, preserving the values of any factor or split variables.



**Example**

```
GET FILE=COVMTX.  
MCONVERT MATRIX=OUT(CORMTX).
```

- GET retrieves the matrix data file *COVMTX*. *COVMTX* becomes the active dataset.
- By default, MCONVERT reads the original matrix from the active dataset. IN(\*) can be specified to make the default explicit.
- The keyword OUT on MATRIX writes the converted matrix to file *CORMTX*.

**REPLACE and APPEND Subcommands**

By default, MCONVERT writes only the converted matrix to the resulting matrix file. Use APPEND to copy both the original matrix and the converted matrix.

- The only specification is the keyword REPLACE or APPEND.
- REPLACE and APPEND are alternatives.
- REPLACE and APPEND affect the resulting matrix file only. The original matrix materials, whether in the active dataset or in an external file, remain intact.

**APPEND**

*Write the original matrix followed by the converted matrix to the matrix file. If there are multiple sets of matrix materials, APPEND appends each converted matrix to the end of a copy of its original matrix.*

**REPLACE**

*Write the original matrix followed by the covariance matrix to the matrix file.*

**Example**

```
MCONVERT MATRIX=OUT(COVMTX) /APPEND.
```

- MCONVERT reads matrix materials from the active dataset.
- The APPEND subcommand copies original matrix materials, appends each converted matrix to the end of the copy of its original matrix, and writes both sets to the file *COVMTX*.

# MEANS

MEANS is available in the Statistics Base option.

```
MEANS [TABLES={varlist} [BY varlist] [BY...] [/varlist...]  
      {ALL      }  
  
[/MISSING={TABLE      }]  
         {INCLUDE  }  
         {DEPENDENT}  
  
[/CELLS= [MEAN** ] [COUNT** ] [STDDEV**]  
         [MEDIAN] [GMEDIAN] [SEMEAN] [SUM ]  
         [MIN] [MAX] [RANGE] [VARIANCE]  
         [KURT] [SEKURT] [SKEW] [SESKEW]  
         [FIRST] [LAST]  
         [NPCT] [SPCT] [NPCT(var)] [SPCT(var)]  
         [HARMONIC] [GEOMETRIC]  
         [DEFAULT]  
         [ALL] ]  
  
[/STATISTICS=[ANOVA] [{LINEARITY}] [NONE**]  
              {ALL      }]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
MEANS TABLES=V1 TO V5 BY GROUP.
```

## Overview

By default, MEANS (alias BREAKDOWN) displays means, standard deviations, and group counts for a numeric dependent variable and group counts for a string variable within groups defined by one or more control (independent) variables. Other procedures that display univariate statistics are SUMMARIZE, FREQUENCIES, and DESCRIPTIVES.

### Options

**Cell Contents.** By default, MEANS displays means, standard deviations, and cell counts for a dependent variable across groups defined by one or more control variables. You can also display sums and variances using the CELLS subcommand.

**Statistics.** In addition to the statistics displayed for each cell of the table, you can obtain a one-way analysis of variance and test of linearity using the STATISTICS subcommand.

### Basic Specification

The basic specification is TABLES with a table list. The actual keyword TABLES can be omitted.

- The minimum table list specifies a dependent variable.
- By default, MEANS displays means, standard deviations, and number of cases.

### **Subcommand Order**

The table list must be first if the keyword TABLES is omitted. If the keyword TABLES is explicitly used, subcommands can be specified in any order.

### **Operations**

- MEANS displays the number and percentage of the processed and missing cases in the Case Process Summary table.
- MEANS displays univariate statistics for the population as a whole and for each value of each successive control variable defined by the BY keyword on the TABLE subcommand in the Group Statistics table.
- ANOVA and linearity statistics, if requested, are displayed in the ANOVA and Measures of Association tables.
- If a string variable is specified as a dependent variable on any table lists, the MEANS procedure produces limited statistics (COUNT, FIRST, and LAST).

### **Limitations**

- Each TABLES subcommand can contain a maximum of 10 BY variable lists.
- There is a maximum of 30 TABLES subcommands for each MEANS command.

## **Examples**

### **Specifying a Range of Dependent Variables**

```
MEANS TABLES=V1 TO V5 BY GROUP
  /STATISTICS=ANOVA.
```

- TABLES specifies that *V1* through *V5* are the dependent variables. *GROUP* is the control variable.
- Assuming that variables *V2*, *V3*, and *V4* lie between *V1* and *V5* in the active dataset, five tables are produced: *V1* by *GROUP*, *V2* by *GROUP*, *V3* by *GROUP*, and so on.
- STATISTICS requests one-way analysis-of-variance tables of *V1* through *V5* by *GROUP*.

### **Creating Analyses for Two Separate Sets of Dependent Variables**

```
MEANS VARA BY VARB BY VARC/V1 V2 BY V3 V4 BY V5.
```

- This command contains two TABLES subcommands that omit the optional TABLES keyword.
- The first table list produces a Group Statistics table for *VARA* within groups defined by each combination of values as well as the totals of *VARB* and *VARC*.
- The second table list produces a Group Statistics table displaying statistics for *V1* by *V3* by *V5*, *V1* by *V4* by *V5*, *V2* by *V3* by *V5*, and *V2* by *V4* by *V5*.

**ANOVA and tests of linearity**

```
MEANS
  TABLES=cigtry BY avcig
  /CELLS MEAN COUNT STDDEV
  /STATISTICS ANOVA LINEARITY .
```

- The procedure specifies *cigtry* as the target variable and *avcig* as a factor.
- `CELLS` requests the cell means, counts, and standard deviations.
- `STATISTICS` requests an analysis of variance and tests of linearity.

**TABLES Subcommand**

`TABLES` specifies the table list.

- You can specify multiple `TABLES` subcommands on a single `MEANS` command (up to a maximum of 30). The slash between the subcommands is required. You can also name multiple table lists separated by slashes on one `TABLES` subcommand.
- The dependent variable is specified first. If the dependent variable is a string variable, `MEANS` produces only limited statistics (`COUNT`, `FIRST`, and `LAST`). The control (independent) variables follow the `BY` keyword and can be numeric (integer or noninteger) or string.
- Each use of the keyword `BY` in a table list adds a dimension to the table requested. Statistics are displayed for each dependent variable by each combination of values and the totals of the control variables across dimensions. There is a maximum of 10 `BY` variable lists for each `TABLES` subcommand.
- The order in which control variables are displayed is the same as the order in which they are specified on `TABLES`. The values of the first control variable defined for the table appear in the leftmost column of the table and change the most slowly in the definition of groups.
- More than one dependent variable can be specified in a table list, and more than one control variable can be specified in each dimension of a table list.

**CELLS Subcommand**

By default, `MEANS` displays the means, standard deviations, and cell counts in each cell. Use `CELLS` to modify cell information.

- If `CELLS` is specified without keywords, `MEANS` displays the default statistics.
- If any keywords are specified on `CELLS`, only the requested information is displayed.
- `MEDIAN` and `GMEDIAN` are expensive in terms of computer resources and time. Requesting these statistics (via these keywords or `ALL`) may slow down performance.

<b>DEFAULT</b>	<i>Means, standard deviations, and cell counts.</i> This is the default if <code>CELLS</code> is omitted.
<b>MEAN</b>	<i>Cell means.</i>
<b>STDDEV</b>	<i>Cell standard deviations.</i>
<b>COUNT</b>	<i>Cell counts.</i>
<b>MEDIAN</b>	<i>Cell median.</i>



<b>GMEDIAN</b>	<i>Grouped median.</i>
<b>SEMEAN</b>	<i>Standard error of cell mean.</i>
<b>SUM</b>	<i>Cell sums.</i>
<b>MIN</b>	<i>Cell minimum.</i>
<b>MAX</b>	<i>Cell maximum.</i>
<b>RANGE</b>	<i>Cell range.</i>
<b>VARIANCE</b>	<i>Variances.</i>
<b>KURT</b>	<i>Cell kurtosis.</i>
<b>SEKURT</b>	<i>Standard error of cell kurtosis.</i>
<b>SKEW</b>	<i>Cell skewness.</i>
<b>SESKEW</b>	<i>Standard error of cell skewness.</i>
<b>FIRST</b>	<i>First value.</i>
<b>LAST</b>	<i>Last value.</i>
<b>NPCT</b>	<i>Percentage of the total number of cases.</i>
<b>SPCT</b>	<i>Percentage of the total sum.</i>
<b>NPCT(var)</b>	<i>Percentage of the total number of cases within the specified variable. The specified variable must be one of the control variables.</i>
<b>SPCT(var)</b>	<i>Percentage of the total sum within the specified variable. The specified variable must be one of the control variables.</i>
<b>HARMONIC</b>	<i>Harmonic mean.</i>
<b>GEOMETRIC</b>	<i>Geometric mean.</i>
<b>ALL</b>	<i>All cell information.</i>

## **STATISTICS Subcommand**

Use **STATISTICS** to request a one-way analysis of variance and a test of linearity for each **TABLE** list.

- Statistics requested on **STATISTICS** are computed in addition to the statistics displayed in the Group Statistics table.
- If **STATISTICS** is specified without keywords, **MEANS** computes **ANOVA**.
- If two or more dimensions are specified, the second and subsequent dimensions are ignored in the analysis-of-variance table. To obtain a two-way and higher analysis of variance, use the **ANOVA** or **MANOVA** procedure. The **ONEWAY** procedure calculates a one-way analysis of variance with multiple comparison tests.

<b>ANOVA</b>	<i>Analysis of variance.</i> <b>ANOVA</b> displays a standard analysis-of-variance table and calculates eta and eta squared (displayed in the Measures of Association table). This is the default if <b>STATISTICS</b> is specified without keywords.
<b>LINEARITY</b>	<i>Test of linearity.</i> <b>LINEARITY</b> (alias <b>ALL</b> ) displays additional statistics to the tables created by the <b>ANOVA</b> keyword: the sums of squares, degrees of freedom, and mean square associated with linear and nonlinear components, the <i>F</i> ratio, and significance level for the <b>ANOVA</b> table and Pearson's <i>r</i> and <i>r</i> <sup>2</sup> for the Measures of Association table. <b>LINEARITY</b> is ignored if the control variable is a string.
<b>NONE</b>	<i>No additional statistics.</i> This is the default if <b>STATISTICS</b> is omitted.

**Example**

```
MEANS TABLES=INCOME BY SEX BY RACE  
  /STATISTICS=ANOVA.
```

- MEANS produces a Group Statistics table of *INCOME* by *RACE* within *SEX* and computes an analysis of variance only for *INCOME* by *SEX*.

**MISSING Subcommand**

MISSING controls the treatment of missing values. If no MISSING subcommand is specified, each combination of a dependent variable and control variables is handled separately.

<b>TABLE</b>	<i>Delete cases with missing values on a tablewise basis. A case with a missing value for any variable specified for a table is not used. Thus, every case contained in a table has a complete set of nonmissing values for all variables in that table. When you separate table requests with a slash, missing values are handled separately for each list. Any MISSING specification will result in tablewise treatment of missing values.</i>
<b>INCLUDE</b>	<i>Include user-missing values. This option treats user-missing values as valid values.</i>
<b>DEPENDENT</b>	<i>Exclude user-missing values for dependent variables only. DEPENDENT treats user-missing values for all control variables as valid.</i>

**References**

Hays, W. L. 1981. *Statistics for the social sciences*, 3rd ed. New York: Holt, Rinehart, and Winston.

# MISSING VALUES

```
MISSING VALUES {varlist}(value list) [[/]{varlist} ...]
                {ALL      }                {ALL      }
```

*Keywords for numeric value lists:*

LO, LOWEST, HI, HIGHEST, THRU

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
MISSING VALUES V1 (8,9) V2 V3 (0) V4 ('X') V5 TO V9 (' ').
```

## **Overview**

MISSING VALUES declares values user-missing. These values can then receive special treatment in data transformations, statistical calculations, and case selection. By default, user-missing values are treated the same as the system-missing values. System-missing values are automatically assigned by the program when no legal value can be produced, such as when an alphabetical character is encountered in the data for a numeric variable, or when an illegal calculation, such as division by 0, is requested in a data transformation.

### **Basic Specification**

The basic specification is a single variable followed by the user-missing value or values in parentheses. Each specified value for the variable is treated as user-missing for any analysis.

### **Syntax Rules**

- Each variable can have a maximum of three individual user-missing values. A space or comma must separate each value. For numeric variables, you can also specify a range of missing values. [For more information, see the topic Specifying Ranges of Missing Values on p. 1199.](#)
- The missing-value specification must correspond to the variable type (numeric or string).
- The same values can be declared missing for more than one variable by specifying a variable list followed by the values in parentheses. Variable lists must have either all numeric or all string variables.
- Different values can be declared missing for different variables by specifying separate values for each variable. An optional slash can be used to separate specifications.
- Missing values for string variables must be enclosed in single or double quotes. The value specifications must include any leading or trailing blanks. [For more information, see the topic String Values in Command Specifications on p. 40.](#)

- For date format variables (for example, `DATE`, `ADATE`), missing values expressed in date formats must be enclosed in single or double quotes, and values must be expressed in the same date format as the defined date format for the variable.
- A variable list followed by an empty set of parentheses ( ) deletes any user-missing specifications for those variables.
- The keyword `ALL` can be used to refer to all user-defined variables in the active dataset, provided the variables are either all numeric or all string. `ALL` can refer to both numeric and string variables if it is followed by an empty set of parentheses. This will delete all user-missing specifications in the active dataset.
- More than one `MISSING VALUES` command can be specified per session.

### **Operations**

- Unlike most transformations, `MISSING VALUES` takes effect as soon as it is encountered. Special attention should be paid to its position among commands. [For more information, see the topic Command Order on p. 41.](#)
- Missing-value specifications can be changed between procedures. New specifications replace previous ones. If a variable is mentioned more than once on one or more `MISSING VALUES` commands before a procedure, only the last specification is used.
- Missing-value specifications are saved in IBM® SPSS® Statistics data files (see `SAVE`) and portable files (see `EXPORT`).

### **Limitations**

Missing values for string variables cannot exceed 8 bytes. (There is no limit on the defined width of the string variable, but defined missing values cannot exceed 8 bytes.)

## **Examples**

### **Declaring Missing Values for Multiple Variables**

```
MISSING VALUES V1 (8,9) V2 V3 (0) V4 ('X') V5 TO V9 (' ' ' ').
```

- The values 8 and 9 are declared missing for the numeric variable *V1*.
- The value 0 is declared missing for the numeric variables *V2* and *V3*.
- The value X is declared missing for the string variable *V4*.
- Blanks are declared missing for the string variables between and including *V5* and *V9*. All of these variables must have a width of four columns.

### **Clearing Missing Values for Selected Variables**

```
MISSING VALUES V1 ( ).
```

- Any previously declared missing values for *V1* are deleted.

### ***Declaring Missing Values for All Variables***

```
MISSING VALUES ALL (9) .
```

- The value 9 is declared missing for all variables in the active dataset; the variables must all be numeric. All previous user-missing specifications are overridden.

### ***Clearing Missing Values for All Variables***

```
MISSING VALUES ALL ( ) .
```

- All previously declared user-missing values for all variables in the active dataset are deleted. The variables in the active dataset can be both numeric and string.

## ***Specifying Ranges of Missing Values***

A range of values can be specified as missing for numeric variables but *not* for string variables.

- The keyword `THRU` indicates an inclusive list of values. Values must be separated from `THRU` by at least one blank space.
- The keywords `HIGHEST` and `LOWEST` with `THRU` indicate the highest and lowest values of a variable. `HIGHEST` and `LOWEST` can be abbreviated to `HI` and `LO`.
- Only one `THRU` specification can be used for each variable or variable list. Each `THRU` specification can be combined with one additional missing value.

### ***Example***

```
MISSING VALUES V1 (LOWEST THRU 0) .
```

- All negative values and 0 are declared missing for the variable *V1*.

### ***Example***

```
MISSING VALUES V1 (0 THRU 1.5) .
```

- Values from 0 through and including 1.5 are declared missing.

### ***Example***

```
MISSING VALUES V1 (LO THRU 0, 999) .
```

- All negative values, 0, and 999 are declared missing for the variable *V1*.

# MIXED

MIXED is available in the Advanced Statistics option.

```
MIXED dependent varname [BY factor list] [WITH covariate list]

[/CRITERIA = [CIN({95** })] [HCONVERGE({0** } {ABSOLUTE**})
              {value} {RELATIVE }

              [LCONVERGE({0** } {ABSOLUTE**})] [MXITER({100**})]
              {value} {RELATIVE } {n }

              [MXSTEP({5**})] [PCONVERGE({1E-6**}, {ABSOLUTE**})] [SCORING({1**})]
              {n } {value } {RELATIVE } {n }

              [SINGULAR({1E-12**})] ]

[/EMMEANS = TABLES ({OVERALL })]
              {factor }
              {factor*factor ...}
              [WITH (covariate=value [covariate = value ...])
              [COMPARE [({factor})] [REFCAT({value})] [ADJ({LSD** })] ]
              {FIRST} {BONFERRONI}
              {LAST } {SIDAK }

[/FIXED = [effect [effect ...]] [| [NOINT] [SSTYPE({1 })] ] ]
              {3**}

[/METHOD = {ML }
           {REML**}

[/MISSING = {EXCLUDE**}]
           {INCLUDE }

[/PRINT = [CORB] [COVB] [CPS] [DESCRIPTIVES] [G] [HISTORY(1**)] [LMATRIX] [R]
           (n )
           [SOLUTION] [TESTCOV]]

[/RANDOM = effect [effect ...]
          [| [SUBJECT(varname[*varname[*...]])] [COVTYPE({VC** })]]]
           {covstruct+}

[/REGWGT = varname]

[/REPEATED = varname[*varname[*...]] | SUBJECT(varname[*varname[*...]])
            [COVTYPE({DIAG** })]]
            {covstruct†}

[/SAVE = [tempvar [(name)] [tempvar [(name)]] ...]

[/TEST[(valuelist)] =
      ['label'] effect valuelist ... [| effect valuelist ...] [divisor=value]]
      [; effect valuelist ... [| effect valuelist ...] [divisor=value]]

[/TEST[(valuelist)] = ['label'] ALL list [| list] [divisor=value]]
      [; ALL list [| list] [divisor=value]]
```

\*\* Default if the subcommand is omitted.

† covstruct can take the following values: AD1, AR1, ARH1, ARMA11, CS, CSH, CSR, DIAG, FA1, FAH1, HF, ID, TP, TPH, UN, UNR, VC.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Example**

```
MIXED Y.
```

## **Overview**

The MIXED procedure fits a variety of mixed linear models. The mixed linear model expands the general linear model used in the GLM procedure in that the data are permitted to exhibit correlation and non-constant variability. The mixed linear model, therefore, provides the flexibility of modeling not only the means of the data but also their variances and covariances.

The MIXED procedure is also a flexible tool for fitting other models that can be formulated as mixed linear models. Such models include multilevel models, hierarchical linear models, and random coefficient models.

### **Important Changes to MIXED Compared to Previous Versions**

**Independence of random effects.** Prior to version 11.5, random effects were assumed to be independent. If you are using MIXED syntax jobs from a version prior to 11.5, be aware that the interpretation of the covariance structure may have changed. [For more information, see the topic Interpretation of Random Effect Covariance Structures on p. 1220.](#)

**Default covariance structures.** Prior to version 11.5, the default covariance structure for random effects was ID, and the default covariance structure for repeated effects was VC.

**Interpretation of VC covariance structure.** Prior to version 11.5, the variance components (VC) structure was a diagonal matrix with heterogeneous variances. Now, when the variance components structure is specified on a RANDOM subcommand, a scaled identity (ID) structure is assigned to each of the effects specified on the subcommand. If the variance components structure is specified on the REPEATED subcommand, it will be replaced by the diagonal (DIAG) structure. Note that the diagonal structure has the same interpretation as the variance components structure in versions prior to 11.5.

### **Basic Features**

**Covariance structures.** Various structures are available. Use multiple RANDOM subcommands to model a different covariance structure for each random effect.

**Standard errors.** Appropriate standard errors will be automatically calculated for all hypothesis tests on the fixed effects, and specified estimable linear combinations of fixed and random effects.

**Subject blocking.** Complete independence can be assumed across subject blocks.

**Choice of estimation method.** Two estimation methods for the covariance parameters are available.

**Tuning the algorithm.** You can control the values of algorithm-tuning parameters with the CRITERIA subcommand.

**Optional output.** You can request additional output through the PRINT subcommand. The SAVE subcommand allows you to save various casewise statistics back to the active dataset.

**Basic Specification**

- The basic specification is a variable list identifying the dependent variable, the factors (if any) and the covariates (if any).
- By default, MIXED adopts the model that consists of the intercept term as the only fixed effect and the residual term as the only random effect.

**Subcommand Order**

- The variable list must be specified first.
- Subcommands can be specified in any order.

**Syntax Rules**

- For many analyses, the MIXED variable list, the FIXED subcommand, and the RANDOM subcommand are the only specifications needed.
- A dependent variable must be specified.
- Empty subcommands are silently ignored.
- Multiple RANDOM subcommands are allowed. However, if an effect with the same subject specification appears in multiple RANDOM subcommands, only the last specification will be used.
- Multiple TEST subcommands are allowed.
- All subcommands, except the RANDOM and the TEST subcommands, should be specified only once. If a subcommand is repeated, only the last specification will be used.
- The following words are reserved as keywords in the MIXED procedure: BY, WITH, and WITHIN.

**Examples**

The following are examples of models that can be specified using MIXED:

**Model 1: Fixed-Effects ANOVA Model**

Suppose that *TREAT* is the treatment factor and *BLOCK* is the blocking factor.

```
MIXED Y BY TREAT BLOCK  
  /FIXED = TREAT BLOCK.
```

**Model 2: Randomized Complete Blocks Design**

Suppose that *TREAT* is the treatment factor and *BLOCK* is the blocking factor.

```
MIXED Y BY TREAT BLOCK  
  /FIXED = TREAT  
  /RANDOM = BLOCK.
```



**Model 3: Split-Plot Design**

An experiment consists of two factors,  $A$  and  $B$ . The experiment unit with respect to  $A$  is  $C$ . The experiment unit with respect to  $B$  is the individual subject, a subdivision of the factor  $C$ . Thus,  $C$  is the whole-plot unit, and the individual subject is the split-plot unit.

```
MIXED Y BY A B C
  /FIXED = A B A*B
  /RANDOM = C(A) .
```

**Model 4: Purely Random-Effects Model**

Suppose that  $A$ ,  $B$ , and  $C$  are random factors.

```
MIXED Y BY A B C
  /FIXED = | NOINT
  /RANDOM = INTERCEPT A B C A*B A*C B*C | COVTYPE(CS) .
```

The MIXED procedure allows effects specified on the same RANDOM subcommand to be correlated. Thus, in the model above, the parameters of a compound symmetry covariance matrix are computed across all levels of the random effects. In order to specify independent random effects, you need to specify separate RANDOM subcommands. For example:

```
MIXED Y BY A B C
  /FIXED = | NOINT
  /RANDOM = INTERCEPT | COVTYPE(ID)
  /RANDOM = A | COVTYPE(CS)
  /RANDOM = B | COVTYPE(CS)
  /RANDOM = C | COVTYPE(CS)
  /RANDOM = A*B | COVTYPE(CS)
  /RANDOM = A*C | COVTYPE(CS)
  /RANDOM = B*C | COVTYPE(CS) .
```

Here, the parameters of compound symmetry matrices are computed separately for each random effect.

**Model 5: Random Coefficient Model**

```
MIXED
  weight WITH time
  /CRITERIA = CIN(95) MXITER(100) MXSTEP(5) SCORING(1)
  SINGULAR(0.000000000001) HCONVERGE(0, ABSOLUTE)
  LCONVERGE(0, ABSOLUTE) PCONVERGE(0.000001, ABSOLUTE)
  /FIXED = time | SSTYPE(3)
  /METHOD = REML
  /PRINT = G SOLUTION TESTCOV
  /RANDOM INTERCEPT time | SUBJECT(patid) COVTYPE(UN) .
```

- The procedure fits a model for *weight*. You believe that there is a linear relationship between weight and time, but there are also patient-specific effects on the size of the relationship that can be accounted for by random effects.
- The fixed-effects portion of the model includes *time*.

- The random-effects portion of the model includes the intercept and *time* with subjects identified by *patid* and an unstructured covariance matrix structure. You are, in essence, fitting a linear regression to each patient in which the regression coefficients are random effects; hence it is a “random coefficients” model.
- The PRINT subcommand requests the estimates of fixed effects, the random effects covariance (G) matrix, and tests of the covariance parameters.
- All other options are set to their default values.

### **Model 6: Multilevel Analysis**

Suppose that *SCORE* is the score of a particular achievement test given over *TIME*. *STUDENT* is nested within *CLASS*, and *CLASS* is nested within *SCHOOL*.

```
MIXED SCORE WITH TIME
  /FIXED = TIME
  /RANDOM = INTERCEPT TIME | SUBJECT(SCHOOL) COVTYPE(ID)
  /RANDOM = INTERCEPT TIME | SUBJECT(SCHOOL*CLASS) COVTYPE(ID)
  /RANDOM = INTERCEPT TIME | SUBJECT(SCHOOL*CLASS*STUDENT) COVTYPE(ID) .
```

### **Model 7: Unconditional Linear Growth Model**

Suppose that *SUBJ* is the individual’s identification and *Y* is the response of an individual observed over *TIME*. The covariance structure is unspecified.

```
MIXED Y WITH TIME
  /FIXED = TIME
  /RANDOM = INTERCEPT TIME | SUBJECT(SUBJ) COVTYPE(ID) .
```

### **Model 8: Linear Growth Model with a Person-Level Covariate**

Suppose that *PCOVAR* is the person-level covariate.

```
MIXED Y WITH TIME PCOVAR
  /FIXED = TIME PCOVAR TIME*PCOVAR
  /RANDOM = INTERCEPT TIME | SUBJECT(SUBJ) COVTYPE(ID) .
```

### **Model 9: Repeated Measures Analysis**

Suppose that *SUBJ* is the individual’s identification and *Y* is the response of an individual observed over several *STAGES*. The covariance structure is compound symmetry.

```
MIXED Y BY STAGE
  /FIXED = STAGE
  /REPEATED = STAGE | SUBJECT(SUBJ) COVTYPE(CS) .
```

### **Model 10: Repeated Measures Analysis with Time-Dependent Covariate**

Suppose that *SUBJ* is the individual’s identification and *Y* is the response of an individual observed over several *STAGES*. *X* is an individual-level covariate that also measures over several *STAGES*. The residual covariance matrix structure is *AR(1)*.

```
MIXED Y BY STAGE WITH X
  /FIXED = X STAGE
```

```
/REPEATED = STAGE | SUBJECT(SUBJ) COVTYPE(AR1) .
```

### Model 11: Repeated Measures and Random Effects

```
MIXED
  sales BY marketid mktsize promo locid
  /CRITERIA = CIN(95) MXITER(100) MXSTEP(10) SCORING(1)
  SINGULAR(0.000000000001) HCONVERGE(0, ABSOLUTE) LCONVERGE(0, ABSOLUTE)
  PCONVERGE(0.000001, ABSOLUTE)
  /FIXED = promo | SSTYPE(3)
  /METHOD = REML
  /PRINT = G R SOLUTION TESTCOV
  /RANDOM = INTERCEPT | SUBJECT(marketid) COVTYPE(ID)
  /REPEATED = week | SUBJECT(locid marketid ) COVTYPE(AR1) .
```

- The procedure fits a model for *sales*. The factor of interest is *promo* (the promotion offered); other factors are used to account for location and market effects.
- The fixed-effects portion of the model includes *promo*.
- The random-effects portion of the model includes the intercept with “subjects” identified by *marketid* and a scaled identity covariance matrix structure. This is equivalent to, but more computationally efficient than, specifying `/RANDOM marketid | COVTYPE(ID)`.
- The repeated-effects portion of the model includes *week* with subjects identified by the unique combinations of *locid* and *marketid*, and a first order autoregressive covariance matrix structure.
- The `PRINT` subcommand requests the estimates of fixed effects, the random effects covariance (G) matrix, the residual covariance (R) matrix, and tests of the covariance parameters.
- All other options are set to their default values.

### Model 12: Crossover Trial

A crossover trial is a study where subjects are exposed to a series of treatments. The end goal is to establish treatment effects, but the model must also account for the effects of the sequencing of treatments. There are a number of ways to specify the model; for example, by using a random effects model of the customers within sequence:

```
MIXED
  amtspent BY coupval seq carry week custid
  /CRITERIA = CIN(95) MXITER(100) MXSTEP(10) SCORING(1)
  SINGULAR(0.000000000001) HCONVERGE(0, ABSOLUTE) LCONVERGE(0, ABSOLUTE)
  PCONVERGE(0.000001, ABSOLUTE)
  /FIXED = coupval seq carry week | SSTYPE(3)
  /METHOD = REML
  /PRINT = G SOLUTION TESTCOV
  /RANDOM = custid(seq) | COVTYPE(ID) .
```

- The procedure fits a model for a carryover trial for *amtspent* (the amount spent by the customer). The factor of interest is *coupval* (the value of the offer to the customer); other factors are used to account for the effects of the sequencing of offers.
- The fixed-effects portion of the model includes *coupval*, *seq* (the sequence in which the customer received offers), *carry* (an indicator to account for carryover effects), and *week* (the week of the trial in which the observation is taken).

- The random-effects portion of the model includes *custid(seq)* (customers within sequence) with a scaled identity covariance matrix structure.
- The `PRINT` subcommand requests the estimates of fixed effects, the random effects covariance (G) matrix, and tests of the covariance parameters.
- All other options are set to their default values.

You can also use a repeated effects model on the time periods:

```
MIXED
  amtspent BY coupval carry week seq
  /CRITERIA = CIN(95) MXITER(100) MXSTEP(100) SCORING(1)
  SINGULAR(0.000000000001) HCONVERGE(0, ABSOLUTE) LCONVERGE(0, ABSOLUTE)
  PCONVERGE(0.000001, ABSOLUTE)
  /FIXED = coupval seq carry week | SSTYPE(3)
  /METHOD = REML
  /PRINT = R SOLUTION TESTCOV
  /REPEATED = week | SUBJECT(custid ) COVTYPE(AR1) .
```

- The procedure fits a model for a carryover trial for *amtspent* (the amount spent by the customer). The factor of interest is *coupval* (the value of the offer to the customer); other factors are used to account for the effects of the sequencing of offers.
- The fixed-effects portion of the model includes *coupval*, *seq* (the sequence in which the customer received offers), *carry* (an indicator to account for carryover effects), and *week* (the week of the trial in which the observation is taken).
- The repeated-effects portion of the model includes *week* with subjects identified by *custid* and a first-order autoregressive covariance matrix structure.
- The `PRINT` subcommand requests the estimates of fixed effects, the residual covariance (R) matrix, and tests of the covariance parameters.
- All other options are set to their default values.

Lastly, you can use a repeated effects model on the treatments:

```
MIXED
  amtspent BY coupval carry week seq
  /CRITERIA = CIN(95) MXITER(100) MXSTEP(10) SCORING(1)
  SINGULAR(0.000000000001) HCONVERGE(0, ABSOLUTE) LCONVERGE(0, ABSOLUTE)
  PCONVERGE(0.000001, ABSOLUTE)
  /FIXED = coupval seq carry week | SSTYPE(3)
  /METHOD = REML
  /PRINT = R SOLUTION TESTCOV
  /REPEATED = coupval | SUBJECT(custid ) COVTYPE(UN) .
```

- The procedure fits a model for a carryover trial for *amtspent* (the amount spent by the customer). The factor of interest is *coupval* (the value of the offer to the customer); other factors are used to account for the effects of the sequencing of offers.
- The fixed-effects portion of the model includes *coupval*, *seq* (the sequence in which the customer received offers), *carry* (an indicator to account for carryover effects), and *week* (the week of the trial in which the observation is taken).
- The repeated-effects portion of the model includes *coupval* with subjects identified by *custid* and an unstructured covariance matrix structure.

- The `PRINT` subcommand requests the estimates of fixed effects, the residual covariance (R) matrix, and tests of the covariance parameters.
- All other options are set to their default values.

## Case Frequency

- If a `WEIGHT` variable is specified, its values are used as frequency weights by the `MIXED` procedure.
- Cases with missing weights or weights less than 0.5 are not used in the analyses.
- The weight values are rounded to the nearest whole numbers before use. For example, 0.5 is rounded to 1, and 2.4 is rounded to 2.

## Covariance Structure List

The following is the list of covariance structures being offered by the `MIXED` procedure. Unless otherwise implied or stated, the structures are not constrained to be non-negative definite in order to avoid nonlinear constraints and to reduce the optimization complexity. However, the variances are restricted to be non-negative.

- Separate covariance matrices are computed for each random effect; that is, while levels of a given random effect are allowed to co-vary, they are considered independent of the levels of other random effects.

**AD1** *First-order ante-dependence.* The constraint  $|\rho_k| \leq 1$  is imposed for stationarity.

$$\sigma^2 \begin{bmatrix} \sigma_1^2 & \sigma_2\sigma_1\rho_1 & \sigma_3\sigma_1\rho_1\rho_2 & \sigma_4\sigma_1\rho_1\rho_2\rho_3 \\ \sigma_2\sigma_1\rho_1 & \sigma_2^2 & \sigma_3\sigma_2\rho_2 & \sigma_4\sigma_2\rho_2\rho_3 \\ \sigma_3\sigma_1\rho_1\rho_2 & \sigma_3\sigma_2\rho_2 & \sigma_3^2 & \sigma_4\sigma_3\rho_3 \\ \sigma_4\sigma_1\rho_1\rho_2\rho_3 & \sigma_4\sigma_2\rho_2\rho_3 & \sigma_4\sigma_3\rho_3 & \sigma_4^2 \end{bmatrix}$$

**AR1** *First-order autoregressive.* The constraint  $|\rho| \leq 1$  is imposed for stationarity.

$$\sigma^2 \begin{bmatrix} 1 & \rho & \rho^2 & \rho^3 \\ \rho & 1 & \rho & \rho^2 \\ \rho^2 & \rho & 1 & \rho \\ \rho^3 & \rho^2 & \rho & 1 \end{bmatrix}$$

**ARH1** *Heterogenous first-order autoregressive.* The constraint  $|\rho_k| \leq 1$  is imposed for stationarity.

$$\sigma^2 \begin{bmatrix} \sigma_1^2 & \sigma_2\sigma_1\rho & \sigma_3\sigma_1\rho^2 & \sigma_4\sigma_1\rho^3 \\ \sigma_2\sigma_1\rho & \sigma_2^2 & \sigma_3\sigma_2\rho & \sigma_4\sigma_2\rho^2 \\ \sigma_3\sigma_1\rho^2 & \sigma_3\sigma_2\rho & \sigma_3^2 & \sigma_4\sigma_3\rho \\ \sigma_4\sigma_1\rho^3 & \sigma_4\sigma_2\rho^2 & \sigma_4\sigma_3\rho & \sigma_4^2 \end{bmatrix}$$

**ARMA11** *Autoregressive moving average (1,1).* The constraints  $|\phi| \leq 1$  and  $|\rho| \leq 1$  are imposed for stationarity.

$$\sigma^2 \begin{bmatrix} 1 & \phi\rho & \phi\rho^2 & \phi\rho^3 \\ \phi\rho & 1 & \phi\rho & \phi\rho^2 \\ \phi\rho^2 & \phi\rho & 1 & \phi\rho \\ \phi\rho^3 & \phi\rho^2 & \phi\rho & 1 \end{bmatrix}$$

**CS** *Compound symmetry.* This structure has constant variance and constant covariance.

$$\begin{bmatrix} \sigma^2 + \sigma_1 & \sigma_1 & \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma^2 + \sigma_1 & \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma_1 & \sigma^2 + \sigma_1 & \sigma_1 \\ \sigma_1 & \sigma_1 & \sigma_1 & \sigma^2 + \sigma_1 \end{bmatrix}$$

**CSH** *Heterogenous compound symmetry.* This structure has non-constant variance and constant correlation.

$$\begin{bmatrix} \sigma_1^2 & \sigma_2\sigma_1\rho & \sigma_3\sigma_1\rho & \sigma_4\sigma_1\rho \\ \sigma_2\sigma_1\rho & \sigma_2^2 & \sigma_3\sigma_2\rho & \sigma_4\sigma_2\rho \\ \sigma_3\sigma_1\rho & \sigma_3\sigma_2\rho & \sigma_3^2 & \sigma_4\sigma_3\rho \\ \sigma_4\sigma_1\rho & \sigma_4\sigma_2\rho & \sigma_4\sigma_3\rho & \sigma_4^2 \end{bmatrix}$$

**CSR** *Compound symmetry with correlation parameterization.* This structure has constant variance and constant covariance.

$$\sigma^2 \begin{bmatrix} 1 & \rho & \rho & \rho \\ \rho & 1 & \rho & \rho \\ \rho & \rho & 1 & \rho \\ \rho & \rho & \rho & 1 \end{bmatrix}$$

**DIAG** *Diagonal.* This is a diagonal structure with heterogenous variance. This is the default covariance structure for repeated effects.

$$\begin{bmatrix} \sigma_1^2 & 0 & 0 & 0 \\ 0 & \sigma_2^2 & 0 & 0 \\ 0 & 0 & \sigma_3^2 & 0 \\ 0 & 0 & 0 & \sigma_4^2 \end{bmatrix}$$

**FA1** *First-order factor analytic with constant diagonal offset ( $d \geq 0$ ).*

$$\begin{bmatrix} \lambda_1^2 + d & \lambda_2\lambda_1 & \lambda_3\lambda_1 & \lambda_4\lambda_1 \\ \lambda_2\lambda_1 & \lambda_2^2 + d & \lambda_3\lambda_2 & \lambda_4\lambda_2 \\ \lambda_3\lambda_1 & \lambda_3\lambda_2 & \lambda_3^2 + d & \lambda_4\lambda_3 \\ \lambda_4\lambda_1 & \lambda_4\lambda_2 & \lambda_4\lambda_3 & \lambda_4^2 + d \end{bmatrix}$$

**FAH1** *First-order factor analytic with heterogenous diagonal offsets ( $d_k \geq 0$ ).*

$$\begin{bmatrix} \lambda_1^2 + d_1 & \lambda_2\lambda_1 & \lambda_3\lambda_1 & \lambda_4\lambda_1 \\ \lambda_2\lambda_1 & \lambda_2^2 + d_2 & \lambda_3\lambda_2 & \lambda_4\lambda_2 \\ \lambda_3\lambda_1 & \lambda_3\lambda_2 & \lambda_3^2 + d_3 & \lambda_4\lambda_3 \\ \lambda_4\lambda_1 & \lambda_4\lambda_2 & \lambda_4\lambda_3 & \lambda_4^2 + d_4 \end{bmatrix}$$

**HF** *Huynh-Feldt.* This is a circular matrix that satisfies the condition  $\sigma_i^2 + \sigma_j^2 - 2\sigma_{ij} = 2\lambda$ .

$$\begin{bmatrix} \sigma_1^2 & \frac{\sigma_1^2 + \sigma_2^2}{2} - \lambda & \frac{\sigma_1^2 + \sigma_3^2}{2} - \lambda & \frac{\sigma_1^2 + \sigma_4^2}{2} - \lambda \\ \frac{\sigma_1^2 + \sigma_2^2}{2} - \lambda & \sigma_2^2 & \frac{\sigma_2^2 + \sigma_3^2}{2} - \lambda & \frac{\sigma_2^2 + \sigma_4^2}{2} - \lambda \\ \frac{\sigma_1^2 + \sigma_3^2}{2} - \lambda & \frac{\sigma_2^2 + \sigma_3^2}{2} - \lambda & \sigma_3^2 & \frac{\sigma_3^2 + \sigma_4^2}{2} - \lambda \\ \frac{\sigma_1^2 + \sigma_4^2}{2} - \lambda & \frac{\sigma_2^2 + \sigma_4^2}{2} - \lambda & \frac{\sigma_3^2 + \sigma_4^2}{2} - \lambda & \sigma_4^2 \end{bmatrix}$$

**ID** *Identity.* This is a scaled identity matrix.

$$\sigma^2 \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix}$$

**TP** *Toeplitz* ( $|\rho_k| \leq 1$ ).

$$\sigma^2 \begin{bmatrix} 1 & \rho_1 & \rho_2 & \rho_3 \\ \rho_1 & 1 & \rho_1 & \rho_2 \\ \rho_2 & \rho_1 & 1 & \rho_1 \\ \rho_3 & \rho_2 & \rho_1 & 1 \end{bmatrix}$$

**TPH** *Heterogenous Toeplitz* ( $|\rho_k| \leq 1$ ).

$$\begin{bmatrix} \sigma_1^2 & \sigma_2\sigma_1\rho_1 & \sigma_3\sigma_1\rho_2 & \sigma_4\sigma_1\rho_3 \\ \sigma_2\sigma_1\rho_1 & \sigma_2^2 & \sigma_3\sigma_2\rho_1 & \sigma_4\sigma_2\rho_2 \\ \sigma_3\sigma_1\rho_2 & \sigma_3\sigma_2\rho_1 & \sigma_3^2 & \sigma_4\sigma_3\rho_1 \\ \sigma_4\sigma_1\rho_3 & \sigma_4\sigma_2\rho_2 & \sigma_4\sigma_3\rho_1 & \sigma_4^2 \end{bmatrix}$$

**UN** *Unstructured.* This is a completely general covariance matrix.

$$\begin{bmatrix} \sigma_1^2 & \sigma_{21} & \sigma_{31} & \sigma_{41} \\ \sigma_{21} & \sigma_2^2 & \sigma_{32} & \sigma_{42} \\ \sigma_{31} & \sigma_{32} & \sigma_3^2 & \sigma_{43} \\ \sigma_{41} & \sigma_{42} & \sigma_{43} & \sigma_4^2 \end{bmatrix}$$

**UNR** *Unstructured correlations* ( $|\rho_{jk}| \leq 1$ ).

$$\begin{bmatrix} \sigma_1^2 & \sigma_2\sigma_1\rho_{21} & \sigma_3\sigma_1\rho_{31} & \sigma_4\sigma_1\rho_{41} \\ \sigma_2\sigma_1\rho_{21} & \sigma_2^2 & \sigma_3\sigma_2\rho_{32} & \sigma_4\sigma_2\rho_{42} \\ \sigma_3\sigma_1\rho_{31} & \sigma_3\sigma_2\rho_{32} & \sigma_3^2 & \sigma_4\sigma_3\rho_{43} \\ \sigma_4\sigma_1\rho_{41} & \sigma_4\sigma_2\rho_{42} & \sigma_4\sigma_3\rho_{43} & \sigma_4^2 \end{bmatrix}$$

**VC** *Variance components.* This is the default covariance structure for random effects. When the variance components structure is specified on a `RANDOM` subcommand, a scaled identity (`ID`) structure is assigned to each of the effects specified on the subcommand. If the variance components structure is specified on the `REPEATED` subcommand, it is replaced by the diagonal (`DIAG`) structure.

## Variable List

The variable list specifies the dependent variable, the factors, and the covariates in the model.

- The dependent variable must be the first specification on `MIXED`.
- The names of the factors, if any, must be preceded by the keyword `BY`.
- The names of the covariates, if any, must be preceded by the keyword `WITH`.
- The dependent variable and the covariates must be numeric.
- The factor variables can be of any type (numeric and string).
- Only cases with no missing values in all of the variables specified will be used.

## CRITERIA Subcommand

The CRITERIA subcommand controls the iterative algorithm used in the estimation and specifies numerical tolerance for checking singularity.

<b>CIN(value)</b>	<i>Confidence interval level.</i> This value is used whenever a confidence interval is constructed. Specify a value greater than or equal to 0 and less than 100. The default value is 95.
<b>HCONVERGE(value, type)</b>	<i>Hessian convergence criterion.</i> Convergence is assumed if $g'_k H_k^{-1} g_k$ is less than a multiplier of <i>value</i> . The multiplier is 1 for ABSOLUTE type and is the absolute value of the current log-likelihood function for RELATIVE type. The criterion is not used if <i>value</i> equals 0. This criterion is not used by default. Specify a non-negative value and a measure type of convergence.
<b>LCONVERGE(value, type)</b>	<i>Log-likelihood function convergence criterion.</i> Convergence is assumed if the ABSOLUTE or RELATIVE change in the log-likelihood function is less than <i>value</i> . The criterion is not used if <i>a</i> equals 0. This criterion is not used by default. Specify a non-negative value and a measure type of convergence.
<b>MXITER(n)</b>	<i>Maximum number of iterations.</i> Specify a non-negative integer. The default value is 100.
<b>PCONVERGE(value, type)</b>	<i>Parameter estimates convergence criterion.</i> Convergence is assumed if the maximum ABSOLUTE or maximum RELATIVE change in the parameter estimates is less than <i>value</i> . The criterion is not used if <i>a</i> equals 0. Specify a non-negative value and a measure type of convergence. The default value for <i>a</i> is $10^{-6}$ .
<b>MXSTEP(n)</b>	<i>Maximum step-halving allowed.</i> At each iteration, the step size is reduced by a factor of 0.5 until the log-likelihood increases or maximum step-halving is reached. Specify a positive integer. The default value is 10.
<b>SCORING(n)</b>	<i>Apply scoring algorithm.</i> Requests to use the Fisher scoring algorithm up to iteration number <i>n</i> . Specify a positive integer. The default is 1.
<b>SINGULAR(value)</b>	<i>Value used as tolerance in checking singularity.</i> Specify a positive value. The default value is $10^{-12}$ .

### Example

```
MIXED SCORE BY SCHOOL CLASS WITH AGE
/CRITERIA = CIN(90) LCONVERGE(0) MXITER(50) PCONVERGE(1E-5 RELATIVE)
/FIXED = AGE
/RANDOM = SCHOOL CLASS.
```

- The CRITERIA subcommand requests that a 90% confidence interval be calculated whenever appropriate.
- The log-likelihood convergence criterion is not used. Convergence is attained when the maximum relative change in parameter estimates is less than 0.00001 and number of iterations is less than 50.

### Example

```
MIXED SCORE BY SCHOOL CLASS WITH AGE
/CRITERIA = MXITER(100) SCORING(100)
/FIXED = AGE
```



/RANDOM = SCHOOL CLASS.

- The Fisher scoring algorithm is used for all iterations.

## **EMMEANS Subcommand**

EMMEANS displays estimated marginal means of the dependent variable in the cells and their standard errors for the specified factors. Note that these are predicted, not observed, means.

- The TABLES keyword, followed by an option in parentheses, is required. COMPARE is optional; if specified, it must follow TABLES.
- Multiple EMMEANS subcommands are allowed. Each is treated independently.
- If identical EMMEANS subcommands are specified, only the last identical subcommand is in effect. EMMEANS subcommands that are redundant but not identical (for example, crossed factor combinations such as  $A*B$  and  $B*A$ ) are all processed.

### **TABLES(option)**

*Table specification.* Valid options are the keyword OVERALL, factors appearing on the factor list, and crossed factors constructed of factors on the factor list. Crossed factors can be specified by using an asterisk (\*) or the keyword BY. All factors in a crossed factor specification must be unique.

If OVERALL is specified, the estimated marginal means of the dependent variable are displayed, collapsing over all factors.

If a factor, or a crossing factor, is specified on the TABLES keyword, MIXED will compute the estimated marginal mean for each level combination of the specified factor(s), collapsing over all other factors not specified with TABLES.

### **WITH (option)**

*Covariate values.* Valid options are covariates appearing on the covariate list on the VARIABLES subcommand. Each covariate must be followed by a numeric value or the keyword MEAN.

If a numeric value is used, the estimated marginal mean will be computed by holding the specified covariate at the supplied value.

When the keyword MEAN is used, the estimated marginal mean will be computed by holding the covariate at its overall mean. If a covariate is not specified in the WITH option, its overall mean will be used in estimated marginal mean calculations.

### **COMPARE(factor) REFCAT(value) ADJ(method)**

*Main- or simple-main-effects omnibus tests and pairwise comparisons of the dependent variable.* This option gives the mean difference, standard error, degrees of freedom, significance, and confidence intervals for each pair of levels for the effect specified in the COMPARE keyword, and an omnibus test for that effect. If only one factor is specified on TABLES, COMPARE can be specified by itself; otherwise, the factor specification is required. In this case, levels of the specified factor are compared with each other for each level of the other factors in the interaction.

The optional `ADJ` keyword allows you to apply an adjustment to the confidence intervals and significance values to account for multiple comparisons. Methods available are `LSD` (no adjustment), `BONFERRONI`, or `SIDAK`.

By default, all pairwise comparisons of the specified factor will be constructed. Optionally, comparisons can be made to a reference category by specifying the value of that category after the `REFCAT` keyword. If the compare factor is a string variable, the category value must be a quoted string. If the compare factor is a numeric variable, the category value should be specified as an unquoted numeric value. Alternatively, the keywords `FIRST` or `LAST` can be used to specify whether the first or the last category will be used as a reference category.

### Example

```
MIXED Y BY A B WITH X
  /FIXED A B X
  /EMMEANS TABLES(A*B) WITH(X=0.23) COMPARE(A) ADJ(SIDAK)
  /EMMEANS TABLES(A*B) WITH(X=MEAN) COMPARE(A) REFCAT(LAST) ADJ(LSD) .
```

- In the example, the first `EMMEANS` subcommand will compute estimated marginal means for all level combinations of  $A*B$  by fixing the covariate  $X$  at 0.23. Then for each level of  $B$ , all pairwise comparisons on  $A$  will be performed using `SIDAK` adjustment.
- In the second `EMMEANS` subcommand, the estimated marginal means will be computed by fixing the covariate  $X$  at its mean. Since `REFCAT(LAST)` is specified, comparison will be made to the last category of factor  $A$  using `LSD` adjustment.

## FIXED Subcommand

The `FIXED` subcommand specifies the fixed effects in the mixed model.

- Specify a list of terms to be included in the model, separated by commas or spaces.
- The intercept term is included by default.
- The default model is generated if the `FIXED` subcommand is omitted or empty. The default model consists of only the intercept term (if included).
- To explicitly include the intercept term, specify the keyword `INTERCEPT` on the `FIXED` subcommand. The `INTERCEPT` term must be specified first on the `FIXED` subcommand.
- To include a main-effect term, enter the name of the factor on the `FIXED` subcommand.
- To include an interaction-effect term among factors, use the keyword `BY` or the asterisk (\*) to connect factors involved in the interaction. For example,  $A*B*C$  means a three-way interaction effect of the factors  $A$ ,  $B$ , and  $C$ . The expression `A BY B BY C` is equivalent to  $A*B*C$ . Factors inside an interaction effect must be distinct. Expressions such as  $A*C*A$  and  $A*A$  are invalid.
- To include a nested-effect term, use the keyword `WITHIN` or a pair of parentheses on the `FIXED` subcommand. For example,  $A(B)$  means that  $A$  is nested within  $B$ , where  $A$  and  $B$  are factors. The expression `A WITHIN B` is equivalent to  $A(B)$ . Factors inside a nested effect must be distinct. Expressions such as  $A(A)$  and  $A(B*A)$  are invalid.

- Multiple-level nesting is supported. For example,  $A(B(C))$  means that  $B$  is nested within  $C$ , and  $A$  is nested within  $B(C)$ . When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus,  $A(B)(C)$  is invalid.
- Nesting within an interaction effect is valid. For example,  $A(B*C)$  means that  $A$  is nested within  $B*C$ .
- Interactions among nested effects are allowed. The correct syntax is the interaction followed by the common nested effect inside the parentheses. For example, the interaction between  $A$  and  $B$  within levels of  $C$  should be specified as  $A*B(C)$  instead of  $A(C)*B(C)$ .
- To include a covariate term in the model, enter the name of the covariate on the `FIXED` subcommand.
- Covariates can be connected using the keyword `BY` or the asterisk (\*). For example,  $X*X$  is the product of  $X$  and itself. This is equivalent to entering a covariate whose values are the squared values of  $X$ .
- Factor and covariate effects can be connected in many ways. Suppose that  $A$  and  $B$  are factors and  $X$  and  $Y$  are covariates. Examples of valid combinations of factor and covariate effects are  $A*X$ ,  $A*B*X$ ,  $X(A)$ ,  $X(A*B)$ ,  $X*A(B)$ ,  $X*Y(A*B)$ , and  $A*B*X*Y$ .
- No effects can be nested within a covariate effect. Suppose that  $A$  and  $B$  are factors and  $X$  and  $Y$  are covariates. The effects  $A(X)$ ,  $A(B*Y)$ ,  $X(Y)$ , and  $X(B*Y)$  are invalid.
- The following options, which are specific for the fixed effects, can be entered after the effects. Use the vertical bar (|) to precede the options.

**NOINT** *No intercept.* The intercept terms are excluded from the fixed effects.

**SSTYPE(n)** *Type of sum of squares.* Specify the methods for partitioning the sums of squares. Specify  $n = 1$  for Type I sum of squares or  $n = 3$  for Type III sum of squares. The default is Type III sum of squares.

### Example

```
MIXED SCORE BY SCHOOL CLASS WITH AGE PRETEST
  /FIXED = AGE(SCHOOL) AGE*PRETEST(SCHOOL)
  /RANDOM = CLASS.
```

- In this example, the fixed-effects design consists of the default `INTERCEPT`, a nested effect  $AGE$  within  $SCHOOL$ , and another nested effect of the product of  $AGE$  and  $PRETEST$  within  $SCHOOL$ .

### Example

```
MIXED SCORE BY SCHOOL CLASS
  /FIXED = | NOINT
  /RANDOM = SCHOOL CLASS.
```

- In this example, a purely random-effects model is fitted. The random effects are  $SCHOOL$  and  $CLASS$ . The fixed-effects design is empty because the implicit intercept term is removed by the `NOINT` keyword.
- You can explicitly insert the `INTERCEPT` effect as `/FIXED = INTERCEPT | NOINT`. But the specification will be identical to `/FIXED = | NOINT`.

## **METHOD Subcommand**

The `METHOD` subcommand specifies the estimation method.

- If this subcommand is not specified, the default is `REML`.
- The keywords `ML` and `REML` are mutually exclusive. Only one of them can be specified once.

**ML**                      *Maximum likelihood.*  
**REML**                    *Restricted maximum likelihood.* This is the default.

## **MISSING Subcommand**

The `MISSING` subcommand specifies the way to handle cases with user-missing values.

- If this subcommand is not specified, the default is `EXCLUDE`.
- Cases, which contain system-missing values in one of the variables, are always deleted.
- The keywords `EXCLUDE` and `INCLUDE` are mutually exclusive. Only one of them can be specified at once.

**EXCLUDE**                *Exclude both user-missing and system-missing values.* This is the default.  
**INCLUDE**                *User-missing values are treated as valid.* System-missing values cannot be included in the analysis.

## **PRINT Subcommand**

The `PRINT` subcommand specifies additional output. If no `PRINT` subcommand is specified, the default output includes:

- A model dimension summary table
- A covariance parameter estimates table
- A model fit summary table
- A test of fixed effects table

**CORB**                    *Asymptotic correlation matrix of the fixed-effects parameter estimates.*  
**COVB**                    *Asymptotic covariance matrix of the fixed-effects parameter estimates.*  
**CPS**                     *Case processing summary.* Displays the sorted values of the factors, the repeated measure variables, the repeated measure subjects, the random-effects subjects, and their frequencies.  
**DESCRIPTIVES**        *Descriptive statistics.* Displays the sample sizes, the means, and the standard deviations of the dependent variable, and covariates (if specified). These statistics are displayed for each distinct combination of the factors.  
**G**                         *Estimated covariance matrix of random effects.* This keyword is accepted only when at least one `RANDOM` subcommand is specified. Otherwise, it will be ignored. If a `SUBJECT` variable is specified for a random effect, then the common block is displayed.  
**HISTORY(n)**            *Iteration history.* The table contains the log-likelihood function value and parameter estimates for every  $n$  iterations beginning with the  $0$ th iteration (the initial estimates). The default is to print every iteration ( $n = 1$ ). If `HISTORY` is specified, the last iteration is always printed regardless of the value of  $n$ .

<b>LMATRIX</b>	<i>Estimable functions.</i> Displays the estimable functions used for testing the fixed effects and for testing the custom hypothesis.
<b>R</b>	<i>Estimated covariance matrix of residual.</i> This keyword is accepted only when a REPEATED subcommand is specified. Otherwise, it will be ignored. If a SUBJECT variable is specified, the common block is displayed.
<b>SOLUTION</b>	<i>A solution for the fixed-effects and the random-effects parameters.</i> The fixed-effects and the random-effects parameter estimates are displayed. Their approximate standard errors are also displayed.
<b>TESTCOV</b>	<i>Tests for the covariance parameters.</i> Displays the asymptotic standard errors and Wald tests for the covariance parameters.

## ***RANDOM Subcommand***

The RANDOM subcommand specifies the random effects in the mixed model.

- Depending on the covariance type specified, random effects specified in one RANDOM subcommand may be correlated.
- One covariance G matrix will be constructed for each RANDOM subcommand. The dimension of the random effect covariance G matrix is equal to the sum of the levels of all random effects in the subcommand.
- When the variance components (VC) structure is specified, a scaled identity (ID) structure will be assigned to each of the effects specified. This is the default covariance type for the RANDOM subcommand.
- Note that the RANDOM subcommand in the MIXED procedure is different in syntax from the RANDOM subcommand in the GLM and VARCOMP procedures.
- Use a separate RANDOM subcommand when a different covariance structure is assumed for a list of random effects. If the same effect is listed on more than one RANDOM subcommand, it must be associated with a different SUBJECT combination.
- Specify a list of terms to be included in the model, separated by commas or spaces.
- No random effects are included in the mixed model unless a RANDOM subcommand is specified correctly.
- Specify the keyword INTERCEPT to include the intercept as a random effect. The MIXED procedure does not include the intercept in the RANDOM subcommand by default. The INTERCEPT term must be specified first on the RANDOM subcommand.
- To include a main-effect term, enter the name of the factor on the RANDOM subcommand.
- To include an interaction-effect term among factors, use the keyword BY or the asterisk (\*) to join factors involved in the interaction. For example, A\*B\*C means a three-way interaction effect of A, B, and C, where A, B, and C are factors. The expression A BY B BY C is equivalent to A\*B\*C. Factors inside an interaction effect must be distinct. Expressions such as A\*C\*A and A\*A are invalid.
- To include a nested-effect term, use the keyword WITHIN or a pair of parentheses on the RANDOM subcommand. For example, A(B) means that A is nested within B, where A and B are factors. The expression A WITHIN B is equivalent to A(B). Factors inside a nested effect must be distinct. Expressions such as A(A) and A(B\*A) are invalid.

- Multiple-level nesting is supported. For example,  $A(B(C))$  means that  $B$  is nested within  $C$ , and  $A$  is nested within  $B(C)$ . When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus,  $A(B)(C)$  is invalid.
- Nesting within an interaction effect is valid. For example,  $A(B*C)$  means that  $A$  is nested within  $B*C$ .
- Interactions among nested effects are allowed. The correct syntax is the interaction followed by the common nested effect inside the parentheses. For example, the interaction between  $A$  and  $B$  within levels of  $C$  should be specified as  $A*B(C)$  instead of  $A(C)*B(C)$ .
- To include a covariate term in the model, enter the name of the covariate on the `FIXED` subcommand.
- Covariates can be connected using the keyword `BY` or the asterisk (\*). For example,  $X*X$  is the product of  $X$  and itself. This is equivalent to entering a covariate whose values are the squared values of  $X$ .
- Factor and covariate effects can be connected in many ways. Suppose that  $A$  and  $B$  are factors and  $X$  and  $Y$  are covariates. Examples of valid combinations of factor and covariate effects are  $A*X$ ,  $A*B*X$ ,  $X(A)$ ,  $X(A*B)$ ,  $X*A(B)$ ,  $X*Y(A*B)$ , and  $A*B*X*Y$ .
- No effects can be nested within a covariate effect. Suppose that  $A$  and  $B$  are factors and  $X$  and  $Y$  are covariates. The effects  $A(X)$ ,  $A(B*Y)$ ,  $X(Y)$ , and  $X(B*Y)$  are invalid.
- The following options, which are specific for the random effects, can be entered after the effects. Use the vertical bar (|) to precede the options.

**SUBJECT(varname\*varname\*...)** *Identify the subjects.* Complete independence is assumed across subjects, thus producing a block-diagonal structure in the covariance matrix of the random effect with identical blocks. Specify a list of variable names (of any type) connected by asterisks. The number of subjects is equal to the number of distinct combinations of values of the variables. A case will not be used if it contains a missing value on any of the subject variables.

**COVTYPE(type)** *Covariance structure.* Specify the covariance structure of the identical blocks for the random effects (see [Covariance Structure List on p. 1207](#)). The default covariance structure for random effects is `VC`.

- If the `REPEATED` subcommand is specified, the variables in the `RANDOM` subject list must be a subset of the variables in the `REPEATED` subject list.
- Random effects are considered independent of each other, and a separate covariance matrix is computed for each effect.

### **Example**

```
MIXED SCORE BY SCHOOL CLASS
  /RANDOM = INTERCEPT SCHOOL CLASS.
```

## **REGWGT Subcommand**

The `REGWGT` subcommand specifies the name of a variable containing the regression weights.

- Specify a numeric variable name following the `REGWGT` subcommand.
- Cases with missing or non-positive weights are not used in the analyses.
- The regression weights will be applied only to the covariance matrix of the residual term.

## ***REPEATED Subcommand***

The `REPEATED` subcommand specifies the residual covariance matrix in the mixed-effects model. If no `REPEATED` subcommand is specified, the residual covariance matrix assumes the form of a scaled identity matrix with the scale being the usual residual variance.

- Specify a list of variable names (of any type) connected by asterisks (repeated measure) following the `REPEATED` subcommand.
- Distinct combinations of values of the variables are used simply to identify the repeated observations. Order of the values will determine the order of occurrence of the repeated observations. Therefore, the lowest values of the variables associate with the first repeated observation, and the highest values associate with the last repeated observation.
- The `VC` covariance structure is obsolete in the `REPEATED` subcommand. If it is specified, it will be replaced with the `DIAG` covariance structure. An annotation will be made in the output to indicate this change.
- The default covariance type for repeated effects is `DIAG`.
- The following keywords, which are specific for the `REPEATED` subcommand, can be entered after the effects. Use the vertical bar (`|`) to precede the options.

**SUBJECT(varname\*varname\*...)** *Identify the subjects.* Complete independence is assumed across subjects, thus producing a block-diagonal structure in the residual covariance matrix with identical blocks. The number of subjects is equal to the number of distinct combinations of values of the variables. A case will not be used if it contains a missing value on any of the subject variables.

**COVTYPE(type)** *Covariance structure.* Specify the covariance structure of the identical blocks for the residual covariance matrix (see [Covariance Structure List on p. 1207](#)). The default structure for repeated effects is `DIAG`.

- The `SUBJECT` keyword must be specified to identify the subjects in a repeated measurement analysis. The analysis will not be performed if this keyword is omitted.
- The list of subject variables must contain all of the subject variables specified in all `RANDOM` subcommands.
- Any variable used in the repeated measure list must not be used in the repeated subject specification.

### ***Example***

```
MIXED SCORE BY CLASS
  /RANDOM = CLASS | SUBJECT(SCHOOL)
  /REPEATED = FLOOR | SUBJECT(SCHOOL*STUDENT) .
```

However, the syntax in each of the following examples is invalid:

```
MIXED SCORE BY CLASS
  /RANDOM = CLASS | SUBJECT(SCHOOL)
  /REPEATED = FLOOR | SUBJECT(STUDENT) .
```

```
MIXED SCORE BY CLASS
  /RANDOM = CLASS | SUBJECT(SCHOOL*STUDENT)
  /REPEATED = FLOOR | SUBJECT(STUDENT) .
```

```
MIXED SCORE BY CLASS
  /RANDOM = CLASS | SUBJECT(SCHOOL)
  /REPEATED = STUDENT | SUBJECT(STUDENT*SCHOOL) .
```

- In the first two examples, the RANDOM subject list contains a variable not on the REPEATED subject list.
- In the third example, the REPEATED subject list contains a variable on the REPEATED variable list.

## SAVE Subcommand

Use the SAVE subcommand to save one or more casewise statistics to the active dataset.

- Specify one or more temporary variables, each followed by an optional new name in parentheses.
- If new names are not specified, default names are generated.

<b>FIXPRED</b>	<i>Fixed predicted values.</i> The regression means without the random effects.
<b>PRED</b>	<i>Predicted values.</i> The model fitted value.
<b>RESID</b>	<i>Residuals.</i> The data value minus the predicted value.
<b>SEFIXP</b>	<i>Standard error of fixed predicted values.</i> These are the standard error estimates for the fixed effects predicted values obtained by the keyword FIXPRED.
<b>SEPREP</b>	<i>Standard error of predicted values.</i> These are the standard error estimates for the overall predicted values obtained by the keyword PRED.
<b>DFFIXP</b>	<i>Degrees of freedom of fixed predicted values.</i> These are the Satterthwaite degrees of freedom for the fixed effects predicted values obtained by the keyword FIXPRED.
<b>DFPREP</b>	<i>Degrees of freedom of predicted values.</i> These are the Satterthwaite degrees of freedom for the fixed effects predicted values obtained by the keyword PRED.

### Example

```
MIXED SCORE BY SCHOOL CLASS WITH AGE
  /FIXED = AGE
  /RANDOM = SCHOOL CLASS(SCHOOL)
  /SAVE = FIXPRED(BLUE) PRED(BLUP) SEFIXP(SEBLUE) SEPREP(SEBLUP) .
```

- The SAVE subcommand appends four variables to the active dataset: *BLUE*, containing the fixed predicted values, *BLUP*, containing the predicted values, *SEBLUE*, containing the standard error of *BLUE*, and *SEBLUP*, containing the standard error of *BLUP*.



## **TEST Subcommand**

The `TEST` subcommand allows you to customize your hypotheses tests by directly specifying null hypotheses as linear combinations of parameters.

- Multiple `TEST` subcommands are allowed. Each is handled independently.
- The basic format for the `TEST` subcommand is an optional list of values enclosed in a pair of parentheses, an optional label in quotes, an effect name or the keyword `ALL`, and a list of values.
- When multiple linear combinations are specified within the same `TEST` subcommand, a semicolon (;) terminates each linear combination except the last one.
- At the end of a contrast coefficients row, you can use the option `DIVISOR=value` to specify a denominator for coefficients in that row. When specified, the contrast coefficients in that row will be divided by the given value. Note that the equals sign is required.
- The value list preceding the first effect or the keyword `ALL` contains the constants, to which the linear combinations are equated under the null hypotheses. If this value list is omitted, the constants are assumed to be zeros.
- The optional label is a string with a maximum length of 255 bytes. Only one label per `TEST` subcommand can be specified.
- The effect list is divided into two parts. The first part is for the fixed effects, and the second part is for the random effects. Both parts have the same syntax structure.
- Effects specified in the fixed-effect list should have already been specified or implied on the `FIXED` subcommand.
- Effects specified in the random-effect list should have already been specified on the `RANDOM` subcommand.
- To specify the coefficient for the intercept, use the keyword `INTERCEPT`. Only one value is expected to follow `INTERCEPT`.
- The number of values following an effect name must be equal to the number of parameters (including the redundant ones) corresponding to that effect. For example, if the effect  $A*B$  takes up to six parameters, then exactly six values must follow  $A*B$ .
- A number can be specified as a fraction with a positive denominator. For example,  $1/3$  or  $-1/3$  are valid, but  $1/-3$  is invalid.
- When `ALL` is specified, only a list of values can follow. The number of values must be equal to the number of parameters (including the redundant ones) in the model.
- Effects appearing or implied on the `FIXED` and `RANDOM` subcommands but not specified on `TEST` are assumed to take the value 0 for all of their parameters.
- If `ALL` is specified for the first row in a `TEST` matrix, then all subsequent rows should begin with the `ALL` keyword.
- If effects are specified for the first row in a `TEST` matrix, then all subsequent rows should use the effect name (thus `ALL` is not allowed).
- When `SUBJECT( )` is specified on a `RANDOM` subcommand, the coefficients given in the `TEST` subcommand will be divided by the number of subjects of that random effect automatically.

**Example**

```

MIXED Y BY A B C
  /FIX = A
  /RANDOM = B C
  /TEST = 'Contrasts of A' A 1/3 1/3 1/3; A 1 -1 0; A 1 -1/2 -1/2
  /TEST(1) = 'Contrast of B' | B 1 -1
  /TEST = 'BLUP at First Level of A'
          ALL 0 1 0 0 | 1 0 1 0;
          ALL          | 1 0 0 1;
          ALL 0 1 0 0;
          ALL 0 1 0 0 | 0 1 0 1.

```

Suppose that factor  $A$  has three levels and factors  $B$  and  $C$  each have two levels.

- The first TEST is labeled *Contrasts of A*. It performs three contrasts among levels of  $A$ . The first is technically not a contrast but the mean of level 1, level 2, and level 3 of  $A$ , the second is between level 1 and level 2 of  $A$ , and the third is between level 1 and the mean of level 2 and level 3 of  $A$ .
- The second TEST is labeled *Contrast of B*. Coefficients for  $B$  are preceded by the vertical bar (|) because  $B$  is a random effect. This contrast computes the difference between level 1 and level 2 of  $B$ , and tests if the difference equals 1.
- The third TEST is labeled *BLUP at First Level of A*. There are four parameters for the fixed effects (intercept and  $A$ ), and there are four parameters for the random effects ( $B$  and  $C$ ). Coefficients for the fixed-effect parameters are separated from those for the random-effect parameters by the vertical bar (|). The coefficients correspond to the parameter estimates in the order in which the parameter estimates are listed in the output.

**Example**

Suppose that factor  $A$  has three levels and factor  $B$  has four levels.

```

MIXED Y BY A B
  /FIXED = A B
  /TEST = 'test example' A 1 -1 0 DIVISOR=3;
          B 0 0 1 -1 DIVISOR=4.

```

- For effect  $A$ , all contrast coefficients will be divided by 3; therefore, the actual coefficients are  $(1/3, -1/3, 0)$ .
- For effect  $B$ , all contrast coefficients will be divided by 4; therefore, the actual coefficients are  $(0, 0, 1/4, -1/4)$ .

**Interpretation of Random Effect Covariance Structures**

This section is intended to provide some insight into the specification random effects and how their covariance structures differ from versions prior to 11.5. Throughout the examples, let  $A$  and  $B$  be factors with three levels, and let  $X$  and  $Y$  be covariates.

**Example (Variance Component Models)**

Random effect covariance matrix of  $A$ :

$$G_A = \sigma_A^2 I_3$$

Random effect covariance matrix of  $B$ :

$$G_B = \sigma_B^2 I_3$$

Overall random effect covariance matrix:

$$G = \begin{bmatrix} G_A & 0 \\ 0 & G_B \end{bmatrix}_{6 \times 6}$$

Prior to version 11.5, this model could be specified by:

```
/RANDOM = A B | COVTYPE(ID)
```

*or*

```
/RANDOM = A | COVTYPE(ID)
/RANDOM = B | COVTYPE(ID)
```

with or without the explicit specification of the covariance structure.

As of version 11.5, this model could be specified by:

```
/RANDOM = A B | COVTYPE(VC)
```

*or*

```
/RANDOM = A | COVTYPE(VC)
/RANDOM = B | COVTYPE(VC)
```

with or without the explicit specification of the covariance structure.

*or*

```
/RANDOM = A | COVTYPE(ID)
/RANDOM = B | COVTYPE(ID)
```

with the explicit specification of the covariance structure.

**Example (Independent Random Effects with Heterogeneous Variances)**

Random effect covariance matrix of  $A$ :

$$G_A = \begin{bmatrix} \sigma_{A1}^2 & 0 & 0 \\ 0 & \sigma_{A2}^2 & 0 \\ 0 & 0 & \sigma_{A3}^2 \end{bmatrix}$$

Random effect covariance matrix of  $B$ :

$$G_B = \begin{bmatrix} \sigma_{B1}^2 & 0 & 0 \\ 0 & \sigma_{B2}^2 & 0 \\ 0 & 0 & \sigma_{B3}^2 \end{bmatrix}$$

Overall random effect covariance matrix:

$$G = \begin{bmatrix} G_A & 0 \\ 0 & G_B \end{bmatrix}_{6 \times 6}$$

Prior to version 11.5, this model could be specified by:

```
/RANDOM = A B | COVTYPE(VC)
```

*or*

```
/RANDOM = A | COVTYPE(VC)
/RANDOM = B | COVTYPE(VC)
```

As of version 11.5, this model could be specified by:

```
/RANDOM = A B | COVTYPE(DIAG)
```

*or*

```
/RANDOM = A | COVTYPE(DIAG)
/RANDOM = B | COVTYPE(DIAG)
```

### **Example (Correlated Random Effects)**

Overall random effect covariance matrix; one column belongs to  $X$  and one column belongs to  $Y$ .

$$G = \begin{bmatrix} \sigma^2 & \sigma^2 \rho \\ \sigma^2 \rho & \sigma^2 \end{bmatrix}$$

Prior to version 11.5, it was impossible to specify this model.

As of version 11.5, this model could be specified by:

```
/RANDOM = A B | COVTYPE(CSR)
```

# MLP

MLP is available in the Neural Networks option.

```
MLP dependent variable [(MLEVEL = {S})] [dependent variable...]  
                        {O}  
                        {N}  
  
    [BY factor list] [WITH covariate list]  
  
[/EXCEPT VARIABLES = varlist]  
  
[/RESCALE [COVARIATE = {STANDARDIZED**}]  
          {NORMALIZED }  
          {ADJNORMALIZED }  
          {NONE }  
  
          [DEPENDENT = {STANDARDIZED }]]  
          {NORMALIZED [(CORRECTION = {0.02**})]}  
          {number}  
          {ADJNORMALIZED [(CORRECTION = {0.02**})]}  
          {number}  
          {NONE }  
  
[/PARTITION {TRAINING = {70** } TESTING = {30** } HOLDOUT = {0** }]]  
            {integer}           {integer}           {integer}  
            {VARIABLE = varname }  
  
[/ARCHITECTURE [AUTOMATIC = {YES**} [(MINUNITS = {1** }, MAXUNITS = {50** } )]]  
              {NO }  
              {integer}           {integer}  
  
              [HIDDENLAYERS = {1** [(NUMUNITS = {AUTO** } )]}  
                {integer}  
                {2 [(NUMUNITS = {AUTO** } )]}  
                {integer, integer}  
              [HIDDENFUNCTION = {TANH** }]  
                {SIGMOID}           [OUTPUTFUNCTION = {IDENTITY}]]  
                {SIGMOID }  
                {SOFTMAX }  
                {TANH }  
  
[/CRITERIA [TRAINING = {BATCH** } ] [MINIBATCHSIZE = {AUTO** }]  
           {ONLINE }  
           {MINIBATCH}           {integer}  
  
           [MEMSIZE = {1000** } ] [OPTIMIZATION = {GRADIENTDESCENT}]  
           {integer}           {SCALEDCONJUGATE}  
  
           [LEARNINGINITIAL = {0.4** } ] [LEARNINGLOWER = {0.001**}]  
           {number}           {number }  
  
           [MOMENTUM = {0.9** } ] [LEARNINGEPOCHS = {10** }]  
           {number}           {integer}  
  
           [LAMBDAINITIAL = {0.000005**}] [SIGMAINITIAL = {0.0005**}]  
           {number }           {number }  
  
           [INTERVALCENTER = {0** } ] [INTERVALOFFSET = {0.5** }]]  
           {number}           {number}  
  
[/STOPPINGRULES [ERRORSTEPS = {1** } [(DATA = {AUTO** } )]]  
               {integer}           {BOTH }  
  
               [TRAININGTIMER = {ON**} [(MAXTIME = {15** } )]]  
               {OFF }           {number}  
  
               [MAXEPOCHS = {AUTO** }  
                 {integer}]
```

```

                [ERRORCHANGE = {0.0001**}]    [ERRORRATIO = {0.001**}]
                        {number }                {number }

[/MISSING USERMISSING = {EXCLUDE**}]
                        {INCLUDE }

[/PRINT [CPS**] [NETWORKINFO**] [SUMMARY**] [CLASSIFICATION**]
        [SOLUTION] [IMPORTANCE] [NONE]]

[/PLOT [NETWORK**] [PREDICTED] [RESIDUAL] [ROC]
       [GAIN] [LIFT] [NONE]]

[/SAVE [PREDDVAL[(varname [varname...])]]
       [PSEUDOPROB[(rootname[:{25 }][rootname...])]]]
                        {integer}

[/OUTFILE [MODEL = 'file' ['file'...]]]

```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 16.0

- Command introduced.

### **Example**

```
MLP dep_var BY A B C WITH X Y Z.
```

## **Overview**

Neural networks are a data mining tool for finding unknown patterns in databases. Neural networks can be used to make business decisions by forecasting demand for a product as a function of price and other variables or by categorizing customers based on buying habits and demographic characteristics. The MLP procedure fits a particular kind of neural network called a multilayer perceptron. The multilayer perceptron uses a feedforward architecture and can have multiple hidden layers. It is one of the most commonly used neural network architectures.

### **Options**

**Prediction or classification.** One or more dependent variables may be specified, and they may be scale, categorical, or a combination. If a dependent variable has a scale measurement level, then the neural network predicts continuous values that approximate the “true” value of some continuous function of the input data. If a dependent variable is categorical, then the neural network is used to classify cases into the “best” category based on the input predictors.

**Rescaling.** MLP optionally rescales covariates or scale dependent variables before training the neural network. There are three rescaling options: standardization, normalization, and adjusted normalization.

**Training, testing, and holdout data.** MLP optionally divides the dataset into training, testing, and holdout data. The neural network is trained using the training data. The training data or testing data, or both, can be used to track errors across steps and determine when to stop training. The holdout data is completely excluded from the training process and is used for independent assessment of the final network.

**Architecture selection.** MLP can perform automatic architecture selection, or it can build a neural network based on user specifications. Automatic architecture selection creates a neural network with one hidden layer and finds the “best” number of hidden units. Alternatively, you can specify one or two hidden layers and define the number of hidden units in each layer.

**Activation functions.** Units in the hidden layers can use the hyperbolic or sigmoid activation functions. Units in the output layer can use the hyperbolic, sigmoid, identity, or softmax activation functions.

**Training methods.** The neural network can be built using batch, online, or mini-batch training. Gradient descent and scaled conjugate gradient optimization algorithms are available.

**Missing Values.** The MLP procedure has an option for treating user-missing values of categorical variables as valid. User-missing values of scale variables are always treated as invalid.

**Output.** MLP displays pivot table output but offers an option for suppressing most such output. Graphical output includes a network diagram (default) and a number of optional charts: predicted-by-observed values, residual-by-predicted values, ROC (Receiver Operating Characteristic) curves, cumulative gains, lift, and independent variable importance. The procedure also optionally saves predicted values in the active dataset. Synaptic weight estimates can be saved in IBM® SPSS® Statistics or XML files also.

### ***Basic Specification***

The basic specification is the MLP command followed by one or more dependent variables, the BY keyword and one or more factors, and the WITH keyword and one or more covariates. By default, the MLP procedure standardizes covariates and selects a training sample before training the neural network. Automatic architecture selection is used to find the “best” neural network architecture. User-missing values are excluded, and default pivot table output is displayed.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### ***Syntax Rules***

- All subcommands are optional.
- Subcommands may be specified in any order.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.

- Empty subcommands are not allowed.
- Any split variable defined on the `SPLIT FILE` command may not be used as a dependent variable, factor, covariate, or partition variable.

### **Limitations**

The `WEIGHT` setting is ignored with a warning by the `MLP` procedure.

### **Categorical Variables**

Although the `MLP` procedure accepts categorical variables as predictors or as the dependent variable, the user should be cautious when using a categorical variable with a very large number of categories.

The `MLP` procedure temporarily recodes categorical predictors and dependent variables using one-of- $c$  coding for the duration of the procedure. If there are  $c$  categories of a variable, then the variable is stored as  $c$  vectors, with the first category denoted  $(1,0,\dots,0)$ , the next category  $(0,1,0,\dots,0)$ , ..., and the final category  $(0,0,\dots,0,1)$ .

This coding scheme increases the number of synaptic weights. In particular, the total number of input units is the number of scale predictors plus the number of categories across all categorical predictors. As a result, this coding scheme can lead to slower training, but more “compact” coding methods usually lead to poorly fit neural networks. If your network training is proceeding very slowly, you might try reducing the number of categories in your categorical predictors by combining similar categories or dropping cases that have extremely rare categories before running the `MLP` procedure.

All one-of- $c$  coding is based on the training data, even if a testing or holdout sample is defined (see [PARTITION Subcommand](#)). Thus, if the testing or holdout samples contain cases with predictor categories that are not present in the training data, then those cases are not used by the procedure or in scoring. If the testing or holdout samples contain cases with dependent variable categories that are not present in the training data, then those cases are not used by the procedure but they may be scored.

### **Replicating results**

The `MLP` procedure uses random number generation during random assignment of partitions, random subsampling for initialization of synaptic weights, random subsampling for automatic architecture selection, and the simulated annealing algorithm used in weight initialization and automatic architecture selection. To reproduce the same randomized results in the future, use the `SET` command to set the initialization value for the random number generator before each run of the `MLP` procedure.

`MLP` results are also dependent on data order. The online and mini-batch training methods are explicitly dependent upon data order; however, even batch training is data-order dependent due to the weight initialization algorithm. For all three training methods, initialization of synaptic weights uses simulated annealing, which takes a random sample from the entire data set and randomly splits it into training (70%) and testing (30%) samples. The size of the random sample is  $N = \min(1000, memsize)$ , where *memsize* is the user-controlled maximum number of cases



stored in memory. (See the `/CRITERIA MEMSIZE` keyword.) If the entire dataset has less than  $N$  cases, then all  $N$  cases are used.

To minimize data order effects, randomly order the cases before running the MLP procedure. To verify the stability of a given solution, you may want to obtain several different solutions with cases sorted in different random orders. In situations with extremely large file sizes, multiple runs can be performed with a sample of cases sorted in different random orders.

Finally, MLP results may be influenced by the variable order on the command line due to the different pattern of initial values assigned when the command line variable order is changed. As with data order effects, you might try different command line variable orders to assess the stability of a given solution.

In summary, if you want to exactly replicate MLP results in the future, use the same initialization value for the random number generator, the same data order, and the same command line variable order, in addition to using the same MLP procedure settings.

## Examples

### **Basic specification with default neural network settings**

```
MLP DepVar BY A B C WITH X Y Z.
```

- The MLP procedure treats *DepVar* as the dependent variable.
- Predictors *A*, *B*, and *C* are factors, and *X*, *Y*, and *Z* are covariates.
- By default, covariates are standardized before training. Also, the active dataset is partitioned into training and testing data samples, with 70% going to the training data and 30% to the testing data sample.
- Automatic architecture selection is used to find the “best” neural network architecture.
- User-missing values are excluded and default output is displayed.

### **User-specified neural network with two hidden layers**

```
MLP DepVar BY A B C WITH X Y Z
  /PARTITION
    TRAINING=100 TESTING=0
  /ARCHITECTURE
    AUTOMATIC=NO
    HIDDENLAYERS=2 (NUMUNITS=25,10)
    OUTPUTFUNCTION=SIGMOID.
```

- The MLP procedure treats *DepVar* as the dependent variable. Predictors *A*, *B*, and *C* are factors, and *X*, *Y*, and *Z* are covariates.
- By default, covariates are standardized before training. The `PARTITION` subcommand overrides the default partitioning of the active dataset into training and testing data and treats all cases as training data.

- The ARCHITECTURE subcommand turns off automatic architecture selection (AUTOMATIC = NO) and specifies a neural network with two hidden layers. There are 25 hidden units in the first hidden layer and 10 hidden units in the second hidden layer. The sigmoid activation function is used for units in the output layer.
- User-missing values are excluded and default output is displayed.

#### **Automatic architecture with partitions specified by variable**

```
*Multilayer Perceptron Network.
MLP default (MLEVEL=N) BY ed WITH age employ address income debtinc
  creddebt othdebt
  /RESCALE COVARIATE=STANDARDIZED
  /PARTITION VARIABLE=partition
  /ARCHITECTURE AUTOMATIC=YES (MINUNITS=1 MAXUNITS=50)
  /CRITERIA TRAINING=BATCH OPTIMIZATION=SCALEDCONJUGATE LAMBDAINITIAL=0.0000005
    SIGMAINITIAL=0.00005 INTERVALCENTER=0 INTERVALOFFSET=0.5
  /PRINT CPS NETWORKINFO SUMMARY CLASSIFICATION IMPORTANCE
  /PLOT ROC GAIN LIFT PREDICTED
  /STOPPINGRULES ERRORSTEPS= 1 (DATA=AUTO) TRAININGTIMER=ON (MAXTIME=15)
    MAXEPOCHS=AUTO ERRORCHANGE=1.0E-4 ERRORRATIO=0.0010
  /MISSING USERMISSING=EXCLUDE .
```

- The procedure builds a network for the nominal-level variable *default*, based upon the factor *ed* and covariates *age* through *othdebt*.
- Cases are assigned to training, testing, and holdout samples based on the values of *partition*.
- In addition to the default tabular output, a sensitivity analysis to compute the importance of each predictor is requested.
- The default graphical output (the network diagram) is not requested, but an ROC curve, cumulative gains chart, lift chart, and predicted by observed chart will be produced.
- All other options are set to their default values.

#### **Multiple dependent variables; two hidden layers with automatic numbers of units selection**

```
*Multilayer Perceptron Network.
MLP los (MLEVEL=S) cost (MLEVEL=S) BY agecat gender diabetes bp smoker choles
  active obesity angina mi nitro anticlot time doa ekg cpk tropt clotsolv
  bleed magnes digi betablk der proc comp
  /RESCALE DEPENDENT=ADJNORMALIZED (CORRECTION=0.02)
  /PARTITION TRAINING=7 TESTING=2 HOLDOUT=1
  /ARCHITECTURE AUTOMATIC=NO HIDDENLAYERS=2 (NUMUNITS=AUTO)
    HIDDENFUNCTION=TANH OUTPUTFUNCTION=TANH
  /CRITERIA TRAINING=ONLINE OPTIMIZATION=GRADIENTDESCENT LEARNINGINITIAL= 0.4
    LEARNINGLOWER= 0.001 LEARNINGEPOCHS= 10 MOMENTUM= 0.9 INTERVALCENTER=0
    INTERVALOFFSET=0.5 MEMSIZE=1000
  /PRINT CPS NETWORKINFO SUMMARY IMPORTANCE
  /PLOT PREDICTED RESIDUAL
  /SAVE PREDVAL
  /STOPPINGRULES ERRORSTEPS= 1 (DATA=AUTO) TRAININGTIMER=ON (MAXTIME=15)
    MAXEPOCHS=AUTO ERRORCHANGE=1.0E-4 ERRORRATIO=0.0010
  /MISSING USERMISSING=INCLUDE .
```

- The procedure has fit a network to the scale variables *los* and *cost*, using *agecat* through *anticlot* and *time* through *comp* as factors.
- The RESCALE subcommand specifies that dependent variables are rescaled using the adjusted normalized method.

- The `PARTITION` subcommand requests that cases be assigned to the training, testing, and holdout samples in a 7:2:1 ratio.
- The `ARCHITECTURE` subcommand specifies a custom architecture with two hidden layers and the hyperbolic tangent as the activation function for the output layer.
- The `CRITERIA` subcommand specifies that online training will be used to estimate the network parameters, using the default settings for the gradient descent algorithm.
- The `PRINT` subcommand requests a sensitivity analysis to compute the importance of each predictor, in addition to the default output.
- The `PLOT` subcommand does not request default graphical output (the network diagram), but predicted-by-observed and residuals-by-predicted charts will be produced.
- The `SAVE` subcommand requests that predicted values be saved to the active dataset.
- The `MISSING` subcommand specifies that user-missing values of factors and categorical dependents be included in the analysis.
- All other options are set to their default values.

## ***Variable Lists***

The command line variable lists specify the dependent variables, any categorical predictors (also known as factors), and any scale predictors (also known as covariates).

### ***Dependent Variables***

- A list of one or more dependent variables must be the first specification on the `MLP` command.
- Each dependent variable may be followed by the measurement level specification, which contains, in parentheses, the `MLEVEL` keyword followed by an equals sign and then `S` for scale, `O` for ordinal, or `N` for nominal. `MLP` treats ordinal and nominal dependent variables equivalently as categorical.
- If a measurement level is specified, then it temporarily overrides a dependent variable's setting in the data dictionary.
- If no measurement level is specified, then `MLP` defaults to the dictionary setting.
- If a measurement level is not specified and no setting is recorded in the data dictionary, then a numeric variable is treated as scale and a string variable is treated as categorical.
- Dependent variables can be numeric or string.
- A string variable may be defined as ordinal or nominal only.

### ***Predictor Variables***

- The names of the factors, if any, must be preceded by the keyword `BY`.
- If keyword `BY` is specified with no factors, then a warning is issued and `BY` is ignored.
- The names of the covariates, if any, must be preceded by the keyword `WITH`.
- If keyword `WITH` is specified with no covariates, then a warning is issued and `WITH` is ignored.
- A dependent variable may not be specified within a factor or covariate list. If a dependent variable is specified within one of these lists, then an error is issued.

- All variables specified within a factor or covariate list must be unique. If duplicate variables are specified within a list, then the duplicates are ignored.
- If duplicate variables are specified across the factor and covariate lists, then an error is issued.
- The universal keywords `TO` and `ALL` may be specified in the factor and covariate lists.
- Factor variables can be numeric or string.
- Covariates must be numeric.
- If no predictors at all are specified, then the procedure fits an input layer containing only the bias unit—that is, the constant-only input layer.
- At least one predictor must be specified.

## ***EXCEPT Subcommand***

The `EXCEPT` subcommand lists any variables that the `MLP` procedure should exclude from the factor or covariate lists on the command line. This subcommand is useful if the factor or covariate lists contain a large number of variables—specified using the `TO` or `ALL` keyword, for example—but there are a few variables (for example, Case ID) that should be excluded.

The `EXCEPT` subcommand is introduced strictly for the purpose of simplifying syntax. Missing values on factors or covariates specified on `EXCEPT` do not affect whether a case is included in the analysis. For example, the following two `MLP` commands are equivalent. In both commands, listwise deletion is based on the dependent variable and factors *A*, *B*, and *C*.

```
MLP DepVar BY A B C.
```

```
MLP DepVar BY A B C D /EXCEPT VARIABLES=D.
```

- The `EXCEPT` subcommand ignores duplicate variables and variables that are not specified on the command line's factor or covariate lists.
- There is no default variable list on the `EXCEPT` subcommand.

## ***RESCALE Subcommand***

The `RESCALE` subcommand is used to rescale covariates or scale dependent variables.

All rescaling is performed based on the training data, even if a testing or holdout sample is defined (see [PARTITION Subcommand](#)). That is, depending on the type of rescaling, the mean, standard deviation, minimum value, or maximum value of a covariate or dependent variable are computed using only the training data. It is important that these covariates or dependent variables have similar distributions across the training, testing, and holdout samples. If the data are partitioned by specifying percentages on the `PARTITION` subcommand, then the `MLP` procedure attempts to ensure this similarity by random assignment. However, if you use the `PARTITION` subcommand `VARIABLE` keyword to assign cases to the training, testing, and holdout samples, then we recommend that you confirm the distributions are similar across samples before running the `MLP` procedure.

**COVARIATE Keyword**

The COVARIATE keyword specifies the rescaling method to use for covariates specified following WITH on the command line. If no covariates are specified on the command line, then the COVARIATE keyword is ignored.

<b>STANDARDIZED</b>	<i>Subtract the mean and divide by the standard deviation, <math>(x - \text{mean})/s</math>. This is the default rescaling method for covariates.</i>
<b>NORMALIZED</b>	<i>Subtract the minimum and divide by the range, <math>(x - \text{min})/(\text{max} - \text{min})</math>.</i>
<b>ADJNORMALIZED</b>	<i>Adjusted version of subtract the minimum and divide by the range, <math>[2 * (x - \text{min})/(\text{max} - \text{min})] - 1</math>.</i>
<b>NONE</b>	<i>No rescaling of covariates.</i>

**DEPENDENT Keyword**

The DEPENDENT keyword specifies the rescaling method to use for scale dependent variables.

- This keyword applies only to scale dependent variables—that is, either MLEVEL=S is specified on the command line or the variable has a scale measurement level based on its data dictionary setting. If a dependent variable is not scale, then the DEPENDENT keyword is ignored for that variable.
- The availability of these rescaling methods for scale dependent variables depends on the output layer activation function in effect.
- If the identity activation function is in effect, then any of the rescaling methods, including NONE, may be requested. If the sigmoid activation function is in effect, then NORMALIZED is required. If the hyperbolic tangent activation function is in effect, then ADJNORMALIZED is required.
- If automatic architecture selection is in effect (/ARCHITECTURE AUTOMATIC=YES), then the default output layer activation function (identity if there are any scale dependent variables) is always used. In this case, the default rescaling method (STANDARDIZED) is also used and the DEPENDENT keyword is ignored.

<b>STANDARDIZED</b>	<i>Subtract the mean and divide by the standard deviation, <math>(x - \text{mean})/s</math>. This is the default rescaling method for scale dependent variables if the output layer uses the identity activation function. This rescaling method may not be specified if the output layer uses the sigmoid or hyperbolic tangent activation function.</i>
<b>NORMALIZED</b>	<i>Subtract the minimum and divide by the range, <math>(x - \text{min})/(\text{max} - \text{min})</math>. This is the required rescaling method for scale dependent variables if the output layer uses the sigmoid activation function. This rescaling method may not be specified if the output layer uses the hyperbolic tangent activation function.</i>

The NORMALIZED keyword may be followed by the CORRECTION option, which specifies a number  $\epsilon$  that is applied as a correction to the rescaling formula. In particular, the corrected formula is  $[x - (\text{min} - \epsilon)] / [(\text{max} + \epsilon) - (\text{min} - \epsilon)]$ . This correction ensures that all rescaled dependent variable values will be within the range of the activation function. A real number greater than or equal to 0 must be specified. The default is 0.02.

- ADJNORMALIZED** *Adjusted version of subtract the minimum and divide by the range,  $[2*(x-min)/(max-min)]-1$ . This is the required rescaling method for scale dependent variables if the output layer uses the hyperbolic tangent activation function. This rescaling method may not be specified if the output layer uses the sigmoid activation function.*
- The ADJNORMALIZED keyword may be followed by the CORRECTION option, which specifies a number  $\epsilon$  that is applied as a correction to the rescaling formula. In particular, the corrected formula is  $\{2*[(x-(min-\epsilon))/((max+\epsilon)-(min-\epsilon))]-1$ . This correction ensures that all rescaled dependent variable values will be within the range of the activation function. A real number greater than or equal to 0 must be specified. The default is 0.02.
- NONE** *No rescaling of scale dependent variables.*

## ***PARTITION Subcommand***

The PARTITION subcommand specifies the method of partitioning the active dataset into training, testing, and holdout samples. The training sample comprises the data records used to train the neural network. The testing sample is an independent set of data records used to track prediction error during training in order to prevent overtraining. The holdout sample is another independent set of data records used to assess the final neural network.

- The partition can be defined by specifying the ratio of cases randomly assigned to each sample (training, testing, and holdout) or by a variable that assigns each case to the training, testing, or holdout sample.
- If the PARTITION subcommand is not specified, then the default partition randomly assigns 70% of the cases to the training sample, 30% to the testing sample, and 0% to the holdout sample. If you want to specify a different random assignment, then you must specify new values for the TRAINING, TESTING, and HOLDOUT keywords. The value specified on each keyword gives the relative number of cases in the active dataset to assign to each sample. For example, /PARTITION TRAINING = 50 TESTING = 30 HOLDOUT = 20 is equivalent to /PARTITION TRAINING = 5 TESTING = 3 HOLDOUT = 2; both subcommands randomly assign 50% of the cases to the training sample, 30% to the testing sample, and 20% to the holdout sample.
- If you want to be able to reproduce results based on the TRAINING, TESTING, and HOLDOUT keywords later, use the SET command to set the initialization value for the random number generator before running the MLP procedure.
- Be aware of the relationship between rescaling and partitioning. [For more information, see the topic RESCALE Subcommand on p. 1230.](#)
- All partitioning is performed after listwise deletion of any cases with invalid data for any variable used by the procedure. See [MISSING Subcommand](#) for details about valid and invalid data.

### ***TRAINING Keyword***

The TRAINING keyword specifies the relative number of cases in the active dataset to randomly assign to the training sample. The value must be an integer greater than 0. The default (if the PARTITION subcommand is not specified) is 70.

**TESTING Keyword**

The `TESTING` keyword specifies the relative number of cases in the active dataset to randomly assign to the testing sample. The value must be an integer greater than 0. The default (if the `PARTITION` subcommand is not specified) is 30.

**HOLDOUT Keyword**

The `HOLDOUT` keyword specifies the relative number of cases in the active dataset to randomly assign to the holdout sample. The value must be an integer greater than 0. The default (if the `PARTITION` subcommand is not specified) is 0.

**VARIABLE Keyword**

The `VARIABLE` keyword specifies a variable that assigns each case in the active dataset to the training, testing, or holdout sample. Cases with a positive value on the variable are assigned to the training sample, cases with a value of 0 to the testing sample, and cases with a negative value to the holdout sample. Cases with a system-missing value are excluded from the analysis. (Any user-missing values for the partition variable are always treated as valid.)

The variable may not be the dependent variable or any variable specified on the command line factor or covariate lists. The variable must be numeric.

**ARCHITECTURE Subcommand**

The `ARCHITECTURE` subcommand is used to specify the neural network architecture. By default, automatic architecture selection is used to build the network. However, you have the option of overriding automatic architecture selection and building a more specific structure.

**AUTOMATIC Keyword**

The `AUTOMATIC` keyword indicates whether to use automatic architecture selection to build the neural network. Automatic architecture selection builds a network with one hidden layer. Using a prespecified range defining the minimum and maximum number of hidden units, automatic architecture selection computes the “best” number of units in the hidden layer. Automatic architecture selection uses the default activation functions for the hidden and output layers.

If automatic architecture selection is used, then a random sample from the total dataset (excluding any data records included in the holdout sample as defined on the `PARTITION` subcommand) is taken and split into training (70%) and testing (30%) samples. This random sample is used to find the architecture and fit the network. Then, the network is retrained using the entire dataset (taking into account the training, testing, and holdout samples defined on the `PARTITION` subcommand), with the synaptic weights obtained from the random sample used as the initial weights.

The size of the random sample  $N = \min(1000, \text{memsize})$ , where `memsize` is the user-specified maximum number of cases to store in memory (see the `MEMSIZE` keyword in [CRITERIA Subcommand](#)). If the total dataset (excluding holdout cases) has less than  $N$  cases, then all cases (excluding holdout cases) are used. If you want to be able to reproduce results based on the

AUTOMATIC keyword later, use the SET command to set the initialization value for the random number generator before running the MLP procedure.

- YES**                    *Use automatic architecture selection to build the network.* This is the default.
- The YES keyword may be followed by parentheses containing the MINUNITS and MAXUNITS options, which specify the minimum and maximum number of units, respectively, that automatic architecture selection will consider in determining the “best” number of units. It is invalid to specify only one option; you must specify both or neither. The options may be specified in any order and must be separated by a comma or space character. Both numbers must be integers greater than 0, with MINUNITS less than MAXUNITS. The defaults are MINUNITS=1, MAXUNITS=50.
- If AUTOMATIC=YES is specified, then all other ARCHITECTURE subcommand keywords are invalid.
- NO**                    *Do not use automatic architecture selection to build the network.* All other ARCHITECTURE subcommand keywords are valid only if AUTOMATIC=NO is specified.

### **HIDDENLAYERS Keyword**

The HIDDENLAYERS keyword specifies the number of hidden layers in the neural network.

This keyword is honored only if automatic architecture selection is not used—that is, if AUTOMATIC=NO. If automatic architecture selection is in effect, then the HIDDENLAYERS keyword is ignored.

- 1**                    *One hidden layer.* This is the default. The HIDDENLAYERS=1 specification may be followed by the NUMUNITS option, which gives the number of units in the first hidden layer, excluding the bias unit. Specify AUTO to automatically compute the number of units based on the number of input and output units. Alternatively, specify an integer greater than or equal to 1 to request a particular number of hidden units. The default is AUTO.
- 2**                    *Two hidden layers.* The HIDDENLAYERS=2 specification may be followed by the NUMUNITS option, which gives the number of units in the first and second hidden layers, excluding the bias unit in each layer. Specify AUTO to automatically compute the numbers of units based on the number of input and output units. Alternatively, specify two integers greater than or equal to 1 to request particular numbers of hidden units in the first and second hidden layers, respectively. The default is AUTO.



**HIDDENFUNCTION Keyword**

The HIDDENFUNCTION keyword specifies the activation function to use for all units in the hidden layers. This keyword is honored only if automatic architecture selection is not used—that is, if AUTOMATIC=NO. If automatic architecture selection is in effect, then the HIDDENFUNCTION keyword is ignored.

- TANH**            *Hyperbolic tangent.* This function has form:  $\gamma(c) = \tanh(c) = (e^c - e^{-c}) / (e^c + e^{-c})$ . It takes real-valued arguments and transforms them to the range (-1, 1). This is the default activation function for all units in the hidden layers.
- SIGMOID**        *Sigmoid.* This function has form:  $\gamma(c) = 1 / (1 + e^{-c})$ . It takes real-valued arguments and transforms them to the range (0, 1).

**OUTPUTFUNCTION Keyword**

The OUTPUTFUNCTION keyword specifies the activation function to use for all units in the output layer.

The activation function used in the output layer has a special relationship with the error function, which is the measure that the neural network is trying to minimize. In particular, the error function is automatically assigned based on the activation function for the output layer. Sum-of-squares error, the sum of the squared deviations between the observed dependent variable values and the model-predicted values, is used when the identity, sigmoid, or hyperbolic tangent activation function is applied to the output layer. Cross-entropy error is used when the softmax activation function is applied to the output layer.

The OUTPUTFUNCTION keyword is honored only if automatic architecture selection is not used—that is, if AUTOMATIC=NO. If automatic architecture selection is in effect, then OUTPUTFUNCTION is ignored.

- IDENTITY**        *Identity.* This function has form:  $\gamma(c) = c$ . It takes real-valued arguments and returns them unchanged. This is the default activation function for units in the output layer if there are any scale dependent variables.
- SIGMOID**        *Sigmoid.* This function has form:  $\gamma(c) = 1 / (1 + e^{-c})$ . It takes real-valued arguments and transforms them to the range (0, 1).
- TANH**            *Hyperbolic tangent.* This function has form:  $\gamma(c) = \tanh(c) = (e^c - e^{-c}) / (e^c + e^{-c})$ . It takes real-valued arguments and transforms them to the range (-1, 1).
- SOFTMAX**        *Softmax.* This function has form:  $\gamma(c_k) = \exp(c_k) / \sum_j \exp(c_j)$ . It takes a vector of real-valued arguments and transforms it to a vector whose elements fall in the range (0, 1) and sum to 1. Softmax is available only if all dependent variables are categorical; if SOFTMAX is specified and there are any scale dependent variables, then an error is issued. This is the default activation function for units in the output layer if all dependent variables are categorical.

**CRITERIA Subcommand**

The CRITERIA subcommand specifies computational and resource settings for the MLP procedure.

**TRAINING Keyword**

The `TRAINING` keyword specifies the training type, which determines how the neural network processes training data records.

The online and mini-batch training methods are explicitly dependent upon data order; however, even batch training is dependent upon data order because initialization of synaptic weights involves subsampling from the dataset. To minimize data order effects, randomly order the cases before running the `MLP` procedure.

<b>BATCH</b>	<i>Batch training.</i> Updates the synaptic weights only after passing all training data records—that is, batch training uses information from all records in the training dataset. Batch training is often preferred because it directly minimizes the total prediction error. However, batch training may need to update the weights many times until one of the stopping rules is met and, hence, may need many data passes. It is most useful for smaller datasets. This is the default training type.
<b>ONLINE</b>	<i>Online training.</i> Updates the synaptic weights after every single training data record—that is, online training uses information from one record at a time. Online training continuously gets a record and updates the weights until one of the stopping rules is met. If all the records are used once and none of the stopping rules is met, then the process continues by recycling the data records. Online training is superior to batch only for larger datasets with associated predictors. If there are many records and many inputs, and their values are not independent of each other, then online training can more quickly obtain a reasonable answer than batch training.
<b>MINIBATCH</b>	<i>Mini-batch training.</i> Divides the training data records into $K$ groups of approximately equal size, then updates the synaptic weights after passing one group—that is, mini-batch training uses information from a group of records. The process then recycles the data group if necessary. The number of training records per mini-batch is determined by the <code>MINIBATCHSIZE</code> keyword. Mini-batch training offers a compromise between batch and online training, and it may be best for “medium-size” datasets.

**MINIBATCHSIZE Keyword**

The `MINIBATCHSIZE` keyword specifies the number of training records per mini-batch.

- Specify `AUTO` to automatically compute the number of records per mini-batch as  $R = \min(\max(M/10, 2), memsize)$ , where  $M$  is the number of training records and  $memsize$  is the maximum number of cases to store in memory (see the `MEMSIZE` keyword below). If the remainder of  $M/R$  is  $r$ , then when the end of the data is reached, the process places the final  $r$  records in the same mini-batch with the first  $R-r$  records of the next data pass. This “wrapping” of mini-batches will place different cases in the mini-batches with each data pass unless  $R$  divides  $M$  with no remainder.
- Alternatively, specify an integer greater than or equal to 2 and less than or equal to  $memsize$  to request a particular number of records. If the number of training records turns out to be less than the specified `MINIBATCHSIZE`, the number of training records is used instead.
- The default is `AUTO`.
- This keyword is ignored if `TRAINING = MINIBATCH` is not in effect.

**MEMSIZE Keyword**

The MEMSIZE keyword specifies the maximum number of cases to store in memory when synaptic weight initialization, automatic architecture selection, and/or mini-batch training is in effect.

- Specify an integer greater than or equal to 2. The default is 1000.

**OPTIMIZATION Keyword**

The OPTIMIZATION keyword specifies the optimization algorithm used to determine the synaptic weights.

**GRADIENTDESCENT**

*Gradient descent.* Gradient descent is the required optimization algorithm for online and mini-batch training. It is optional for batch training.

When gradient descent is used with online and mini-batch training, the algorithm's user-specified parameters are the initial learning rate, lower bound for the learning rate, momentum, and number of data passes (see the LEARNINGINITIAL, LEARNINGLOWER, MOMENTUM, and LEARNINGEPOCHS keywords, respectively). With batch training, the user-specified parameters are the initial learning rate and the momentum.

**SCALEDCONJUGATE**

*Scaled conjugate gradient.* Scaled conjugate gradient is the default for batch training. The assumptions that justify the use of conjugate gradient methods do not apply to the online and mini-batch training, so this method may not be used if TRAINING = ONLINE or MINIBATCH. The user-specified parameters are the initial lambda and sigma (see the LAMBDAINITIAL and SIGMAINITIAL keywords).

**LEARNINGINITIAL Keyword**

The LEARNINGINITIAL keyword specifies the initial learning rate  $\eta_0$  for the gradient descent optimization algorithm.

- Specify a number greater than 0. The default is 0.4.
- This keyword is ignored if OPTIMIZATION = GRADIENTDESCENT is not in effect.

**LEARNINGLOWER Keyword**

The LEARNINGLOWER keyword specifies the lower boundary for the learning rate  $\eta_{\text{low}}$  when gradient descent is used with online or mini-batch training.

- Specify a number greater than 0 and less than the initial learning rate (see the LEARNINGINITIAL keyword). The default is 0.001.
- This keyword is ignored if TRAINING = ONLINE or MINIBATCH and OPTIMIZATION = GRADIENTDESCENT are not in effect.

**MOMENTUM Keyword**

The MOMENTUM keyword specifies the initial momentum rate  $\alpha$  for the gradient descent optimization algorithm.

- Specify a number greater than 0. The default is 0.9.
- This keyword is ignored if `OPTIMIZATION = GRADIENTDESCENT` is not in effect.

### ***LEARNINGEPOCHS Keyword***

The `LEARNINGEPOCHS` keyword specifies the number of epochs (data passes of the training set)  $p$  to reduce the learning rate when gradient descent is used with online or mini-batch training. You can control the learning rate decay factor  $\beta$  by specifying the number of epochs it takes for the learning rate to decrease from  $\eta_0$  to  $\eta_{low}$ . This corresponds to  $\beta = (1/pK) * \ln(\eta_0/\eta_{low})$ , where  $K$  is the total number of mini-batches in the training dataset. For online training,  $K = M$ , where  $M$  is the number of training records.

- Specify an integer greater than 0. The default is 10.
- This keyword is ignored if `TRAINING = ONLINE` or `MINIBATCH` and `OPTIMIZATION = GRADIENTDESCENT` are not in effect.

### ***LAMBDAINITIAL Keyword***

The `LAMBDAINITIAL` keyword specifies the initial lambda,  $\lambda_0$ , for the scaled conjugate gradient optimization algorithm.

- Specify a number greater than 0 and less than  $10^{-6}$ . The default is 0.0000005.
- This keyword is ignored if `OPTIMIZATION = SCALEDCONJUGATE` is not in effect.

### ***SIGMAINITIAL Keyword***

The `SIGMAINITIAL` keyword specifies the initial sigma,  $\sigma_0$ , for the scaled conjugate gradient optimization algorithm.

- Specify a number greater than 0 and less than  $10^{-4}$ . The default is 0.00005.
- This keyword is ignored if `OPTIMIZATION = SCALEDCONJUGATE` is not in effect.

### ***INTERVALCENTER and INTERVALOFFSET Keywords***

The `INTERVALCENTER` and `INTERVALOFFSET` keywords specify the interval  $[a_0 - a, a_0 + a]$  in which weight vectors are randomly generated when simulated annealing is used. `INTERVALCENTER` corresponds to  $a_0$  and `INTERVALOFFSET` corresponds to  $a$ .

- Simulated annealing is used to break out of a local minimum, with the goal of finding the global minimum, during the optimization algorithm. This approach is used in weight initialization and automatic architecture selection.
- Specify a number for `INTERVALCENTER`. The `INTERVALCENTER` default is 0. Specify a number greater than 0 for `INTERVALOFFSET`. The `INTERVALOFFSET` default is 0.5. The default interval is  $[-0.5, 0.5]$ .

## **STOPPINGRULES Subcommand**

The `STOPPINGRULES` subcommand specifies the rules that determine when to stop training the neural network. Training proceeds through at least one data pass. Training can then be stopped according to the following criteria, which are listed as `STOPPINGRULES` keywords. Stopping rules are checked in the listed order. (In the keyword descriptions, a step is a data pass for the online and mini-batch methods, an iteration for the batch method.)

*Note:* After each complete data pass, online and mini-batch training require an extra data pass in order to compute the training error. This extra data pass can slow training considerably, so if you use online or mini-batch training, we recommend specifying a testing dataset. Then, if you use only the testing set in the `ERRORSTEPS` criterion, the `ERRORCHANGE` and `ERRORRATIO` criteria will not be checked.

### **ERRORSTEPS Keyword**

The `ERRORSTEPS` keyword specifies the number of steps,  $n$ , to allow before checking for a decrease in error. If there is no decrease in error after  $n$  steps, then training stops.

- Any integer greater than or equal to 1 may be specified. The default is 1.
- The `DATA` option following `ERRORSTEPS` specifies how to compute error.

#### **(DATA=AUTO)**

*Compute error using the testing sample if it exists or using the training sample otherwise.* If the error at any step does not decrease below the current minimum error (based on preceding steps) over the next  $n$  steps, then training stops.

For online and mini-batch training, if there is no testing sample, then the procedure computes error using the training sample. Batch training, on the other hand, guarantees a decrease in the training sample error after each data pass, thus this option is ignored if batch training is in effect and there is no testing sample.

`DATA = AUTO` is the default option.

#### **(DATA=BOTH)**

*Compute errors using the testing sample and the training sample.* If neither the testing sample error nor the training sample error decreases below its current minimum error over the next  $n$  steps, then training stops.

For batch training, which guarantees a decrease in the training sample error after each data pass, this option is the same as `DATA= AUTO`.

`DATA = BOTH` may be specified only if testing data are defined—that is, `/PARTITION TESTING` is specified with a number greater than zero or `/PARTITION VARIABLE` is used. If `DATA = BOTH` is specified when `/PARTITION TESTING = 0`, or when `/PARTITION VARIABLE` is used but no testing data exist in the active dataset, then an error is issued.

### **TRAININGTIMER Keyword**

The `TRAININGTIMER` keyword specifies whether the training timer is turned on or off.

- If `TRAININGTIMER = ON`, then the `MAXTIME` option gives the maximum number of minutes allowed for training. Training stops if the algorithm exceeds the maximum allotted time.
- If `TRAININGTIMER = OFF`, then the `MAXTIME` option is ignored.

- `TRAININGTIMER` may be specified with keyword `ON` or `OFF`. The default is `ON`.
- The `MAXTIME` option may be specified with any number greater than 0. The default is 15.

### **MAXEPOCHS Keyword**

The `MAXEPOCHS` keyword specifies the maximum number of epochs (data passes) allowed for the training data. If the maximum number of epochs is exceeded, then training stops.

- Specify `AUTO` to automatically compute the maximum number of epochs as  $\max(2N+1, 100)$ , where  $N$  is the number of synaptic weights in the neural network.
- Alternatively, specify an integer greater than 0 to request a particular maximum number of epochs.
- The default is `AUTO`.

### **ERRORCHANGE Keyword**

The `ERRORCHANGE` keyword specifies the relative change in training error criterion. Training stops if the relative change in the training error compared to the previous step is less than the criterion value.

- Any number greater than 0 may be specified. The default is 0.0001.
- For online and mini-batch training, this criterion is ignored if the `ERRORSTEPS` criterion uses only testing data.

### **ERRORRATIO Keyword**

The `ERRORRATIO` keyword specifies the training error ratio criterion. Training stops if the ratio of the training error to the error of the null model is less than the criterion value. The null model predicts the average value for all dependent variables.

- Any number greater than 0 may be specified. The default is 0.001.
- For online and mini-batch training, this criterion is ignored if the `ERRORSTEPS` criterion uses only testing data.

## **MISSING Subcommand**

The `MISSING` subcommand is used to control whether user-missing values for categorical variables—that is, factors and categorical dependent variables—are treated as valid values. By default, user-missing values for categorical variables are treated as invalid.

- User-missing values for scale variables are always treated as invalid.
- System-missing values for any variables are always treated as invalid.

#### **USERMISSING=EXCLUDE**

*User-missing values for categorical variables are treated as invalid.* This is the default.

#### **USERMISSING=INCLUDE**

*User-missing values for categorical variables are treated as valid values.*

---

## ***PRINT Subcommand***

The `PRINT` subcommand indicates the tabular output to display and can be used to request a sensitivity analysis. If `PRINT` is not specified, then the default tables are displayed. If `PRINT` is specified, then only the requested `PRINT` output is displayed.

### ***CPS Keyword***

The `CPS` keyword displays the case processing summary table, which summarizes the number of cases included and excluded in the analysis, in total and by training, testing, and holdout samples. This table is shown by default.

### ***NETWORKINFO Keyword***

The `NETWORKINFO` keyword displays information about the neural network, including the dependent variables, number of input and output units, number of hidden layers and units, and activation functions. This table is shown by default.

### ***SUMMARY Keyword***

The `SUMMARY` keyword displays a summary of the neural network results, including the error, the relative error or percent of incorrect predictions, the stopping rule used to stop training, and the training time.

- The error is the sum-of-squares error when the identity, sigmoid, or hyperbolic tangent activation function is applied to the output layer. It is the cross-entropy error when the softmax activation function is applied to the output layer.
- In addition, relative errors or percents of incorrect predictions are displayed, depending on the dependent variable measurement levels. If any dependent variable has a scale measurement level, then the average overall relative error (relative to the mean model) is displayed. If all dependent variables are categorical, then the average percent of incorrect predictions is displayed. Relative errors or percents of incorrect predictions are also displayed for individual dependent variables.
- Summary results are given for the training data and for testing and hold-out data if they exist.
- This table is shown by default.

### ***CLASSIFICATION Keyword***

The `CLASSIFICATION` keyword displays a classification table for each categorical dependent variable. The table gives the number of cases classified correctly and incorrectly for each dependent variable category.

- In addition to classification tables, the `CLASSIFICATION` keyword reports the percent of the total cases that were correctly classified. A case is correctly classified if its highest predicted probabilities correspond to the observed categories for that case.
- Classification results are given for the training data and for testing and holdout data if they exist.

- Classification results are shown by default.
- The CLASSIFICATION keyword is ignored for scale dependent variables.

**SOLUTION Keyword**

The SOLUTION keyword displays the synaptic weights—that is, the coefficient estimates, from layer  $i-1$  unit  $j$  to layer  $i$  unit  $k$ . The synaptic weights are based on the training sample even if the active dataset is partitioned into training, testing, and holdout data.

This table is not shown by default because the number of synaptic weights may be extremely large, and these weights are generally not used for interpreting network results.

**IMPORTANCE Keyword**

The IMPORTANCE keyword performs a sensitivity analysis, which computes the importance of each predictor in determining the neural network. The analysis is based on the combined training and testing samples or only the training sample if there is no testing sample. This keyword creates a table and a chart displaying importance and normalized importance for each predictor.

Sensitivity analysis is not performed by default because it is computationally expensive and time-consuming if there are a large number of predictors or cases.

**NONE Keyword**

The NONE keyword suppresses all PRINT output except the Notes table and any warnings. This keyword may not be specified with any other PRINT keywords.

**PLOT Subcommand**

The PLOT subcommand indicates the chart output to display. If PLOT is not specified, then the default chart (the network diagram) is displayed. If PLOT is specified, then only the requested PLOT output is displayed.

**NETWORK Keyword**

The NETWORK keyword displays the network diagram. This chart is shown by default.

**PREDICTED Keyword**

The PREDICTED keyword displays a predicted-by-observed value chart for each dependent variable. For categorical dependent variables, a boxplot of predicted pseudo-probabilities is displayed. For scale dependent variables, a scatterplot is displayed.

- Predicted-by-observed value charts are based on the combined training and testing samples or only the training sample if there is no testing sample.



**RESIDUAL Keyword**

The `RESIDUAL` keyword displays a residual-by-predicted value chart for each scale dependent variable. This chart is available only for scale dependent variables. The `RESIDUAL` keyword is ignored for categorical dependent variables.

- Residual-by-predicted value charts are based on the combined training and testing samples or only the training sample if there is no testing sample.

**ROC Keyword**

The `ROC` keyword displays an ROC (Receiver Operating Characteristic) chart for each categorical dependent variable. It also displays a table giving the area under each curve in the chart.

- For a given dependent variable, the ROC chart displays one curve for each category. If the dependent variable has two categories, then each curve treats the category at issue as the positive state versus the other category. If the dependent variable has more than two categories, then each curve treats the category at issue as the positive state versus the aggregate of all other categories.
- This chart is available only for categorical dependent variables. The `ROC` keyword is ignored for scale dependent variables.
- ROC charts and area computations are based on the combined training and testing samples or only the training sample if there is no testing sample.

**GAIN Keyword**

The `GAIN` keyword displays a cumulative gains chart for each categorical dependent variable.

- The display of one curve for each dependent variable category is the same as for the `ROC` keyword.
- This chart is available only for categorical dependent variables. The `GAIN` keyword is ignored for scale dependent variables.
- Cumulative gains charts are based on the combined training and testing samples or only the training sample if there is no testing sample.

**LIFT Keyword**

The `LIFT` keyword displays a lift chart for each categorical dependent variable.

- The display of one curve for each dependent variable category is the same as for the `ROC` keyword.
- This chart is available only for categorical dependent variables. The `LIFT` keyword is ignored for scale dependent variables.
- Lift charts are based on the combined training and testing samples or only the training sample if there is no testing sample.

**NONE Keyword**

The NONE keyword suppresses all PLOT output. This keyword may not be specified with any other PLOT keywords.

**SAVE Subcommand**

The SAVE subcommand writes optional temporary variables to the active dataset.

**PREDVAL(varname varname...)**

*Predicted value or category.* This saves the predicted value for scale dependent variables and the predicted category for categorical dependent variables.

Specify one or more unique, valid variable names. There should be as many variable names specified as there are dependent variables, and the names should be listed in the order of the dependent variables on the command line. If you do not specify enough variable names, then default names are used for any remaining variables. If you specify too many variable names, then any remaining names are ignored.

If there is only one dependent variable, then the default variable name is *MLP\_PredictedValue*. If there are multiple dependent variables, then the default variable names are *MLP\_PredictedValue\_1*, *MLP\_PredictedValue\_2*, etc., corresponding to the order of the dependent variables on the command line.

**PSEUDOPROB(rootname:n rootname...)**

*Predicted pseudo-probability.* If a dependent variable is categorical, then this keyword saves the predicted pseudo-probabilities of the first *n* categories of that dependent variable.

Specify one or more unique, valid variable names. There should be as many variable names specified as there are categorical dependent variables, and the names should be listed in the order of the categorical dependent variables on the command line. The specified names are treated as rootnames. Suffixes are added to each rootname to get a group of variable names corresponding to the categories for a given dependent variable. If you do not specify enough variable names, then default names are used for any remaining categorical dependent variables. If you specify too many variable names, then any remaining names are ignored.

A colon and a positive integer giving the number of probabilities to save for a dependent variable can follow the rootname.

If there is only one dependent variable, then the default rootname is *MLP\_PseudoProbability*. If there are multiple dependent variables, then the default rootnames are *MLP\_PseudoProbability\_1*, *MLP\_PseudoProbability\_2*, etc., corresponding to the order of the categorical dependent variables on the command line and taking into account the position of any scale dependent variables.

The default *n* is 25.

This keyword is ignored for scale dependent variables.

**Probabilities and Pseudo-probabilities**

Categorical dependent variables with softmax activation and cross-entropy error will have a predicted value for each category, where each predicted value is the probability that the case belongs to the category.

Categorical dependent variables with sum-of-squares error will have a predicted value for each category, but the predicted values cannot be interpreted as probabilities. The `SAVE` subcommand saves these predicted pseudo-probabilities even if any are less than zero or greater than one or the sum for a given dependent variable is not 1.

The ROC, cumulative gains, and lift charts (see `/PLOT ROC, GAIN, and LIFT`, respectively) are created based on pseudo-probabilities. In the event that any of the pseudo-probabilities are less than zero or greater than one or the sum for a given variable is not 1, they are first rescaled to be between zero and one and to sum to 1. The `SAVE` subcommand saves the original pseudo-probabilities, but the charts are based on rescaled pseudo-probabilities.

Pseudo-probabilities are rescaled by dividing by their sum. For example, if a case has predicted pseudo-probabilities of 0.50, 0.60, and 0.40 for a three-category dependent variable, then each pseudo-probability is divided by the sum 1.50 to get 0.33, 0.40, and 0.27.

If any of the pseudo-probabilities are negative, then the absolute value of the lowest is added to all pseudo-probabilities before the above rescaling. For example, if the pseudo-probabilities are -0.30, 0.50, and 1.30, then first add 0.30 to each value to get 0.00, 0.80, and 1.60. Next, divide each new value by the sum 2.40 to get 0.00, 0.33, and 0.67.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand saves XML-format (PMML) files containing the synaptic weights. You can use this model file to apply the model information to other data files for scoring purposes. For more information, see the topic [Scoring expressions on p. 116](#).

- Filenames must be specified in full. `MLP` does not supply extensions.
- The `MODEL` keyword is not honored if split-file processing is in effect (see [SPLIT FILE](#)). If this keyword is specified when split-file processing is on, a warning is displayed.

**MODEL = 'file' 'file'...**

*Writes the synaptic weights to XML (PMML) files.* Specify one or more unique, valid filenames. There should be as many filenames as there are dependent variables, and the names should be listed in the order of the dependent variables on the command line. If you do not specify enough filenames, then an error is issued. If you specify too many filenames, then any remaining names are ignored.

If any 'file' specification refers to an existing file, then the file is overwritten. If any 'file' specifications refer to the same file, then only the last instance of this 'file' specification is honored.

# ***MODEL CLOSE***

```
MODEL CLOSE NAME={handlelist}
                  {ALL      }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Release history***

Release 13.0

- Command introduced.

## ***Example***

```
MODEL CLOSE NAME=discrimmod1 twostep1.
MODEL CLOSE NAME=ALL.
```

## ***Overview***

MODEL CLOSE is used to discard cached models and their associated model handle names (see MODEL HANDLE on p. 1247).

## ***Basic specification***

The basic specification is NAME followed by a list of model handles. Each model handle name should match the name specified on the MODEL HANDLE command. The keyword ALL specifies that all model handles are to be closed.

# MODEL HANDLE

```
MODEL HANDLE NAME=handle FILE='file specification'  
  [/OPTIONS [MISSING={SUBSTITUTE**}] ]  
           {SYSMIS      }  
  
  [/MAP  VARIABLES=varlist MODELVARIABLES=varlist  ]
```

\*\*Default if the keyword is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release history**

Release 13.0

- Command introduced.

## **Example**

```
MODEL HANDLE NAME=discrimmod1 FILE='/modelfiles/discrim1.mml'.
```

## **Overview**

MODEL HANDLE reads an external XML file or ZIP archive containing specifications for a predictive model. It caches the model specifications and associates a unique name (handle) with the cached model. The model can then be used by the APPLYMODEL and STRAPPLYMODEL transformation functions to calculate scores and other results (see Scoring expressions on p. 116). The MODEL CLOSE command is used to discard a cached model from memory.

Different models can be applied to the same data by using separate MODEL HANDLE commands for each of the models.

## **Options**

**Variable Mapping.** You can map any or all of the variables in the original model to different variables in the current active dataset. By default, the model is applied to variables in the current active dataset with the same names as the variables in the original model.

**Handling Missing Values.** You can choose how to handle cases with missing values. By default, an attempt is made to substitute a sensible value for a missing value, but you can choose to treat missing values as system-missing.

## **Basic specification**

The basic specification is NAME and FILE. NAME specifies the model handle name to be used when referring to this model. FILE specifies the external file containing the model specifications.

**Subcommand order**

- Subcommands can be specified in any order.

**Syntax rules**

- When using the MAP subcommand, you must specify both the VARIABLES and MODELVARIABLES keywords.
- Multiple MAP subcommands are allowed. Each MAP subcommand should provide the mappings for a distinct subset of the variables. Subsequent mappings of a given variable override any previous mappings of that same variable.

**Operations**

- A model handle is used only during the current working session. The handle is not saved with the data file.
- Issuing a SET LOCALE command that changes the computer's code page requires closing any existing model handles (using MODEL CLOSE) and reopening the models (using MODEL HANDLE) before proceeding with scoring.

**Models Supported for Scoring**

IBM® SPSS® Statistics can score models created by SPSS Statistics, IBM® SPSS® Modeler, and IBM® SPSS® AnswerTree®.

The following table lists the set of scoring functions available for each type of model that supports scoring. The function type denoted as PROBABILITY (category) refers to specification of a particular category (the optional third parameter) for the PROBABILITY function.

Table 141-1  
Supported functions by model type

Model type	Supported functions
Tree (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE, NODEID
Tree (scale target)	PREDICT, NODEID, STDDEV
Boosted Tree (C5.0)	PREDICT, CONFIDENCE
Linear Regression	PREDICT, STDDEV
Automatic Linear Models	PREDICT
Binary Logistic Regression	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Conditional Logistic Regression	PREDICT
Multinomial Logistic Regression	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
General Linear Model	PREDICT, STDDEV
Discriminant	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
TwoStep Cluster	PREDICT
K-Means Cluster	PREDICT, CONFIDENCE
Kohonen	PREDICT
Neural Net (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE

Model type	Supported functions
Neural Net (scale target)	PREDICT
Naive Bayes	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Anomaly Detection	PREDICT
Ruleset	PREDICT, CONFIDENCE
Generalized Linear Model (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Generalized Linear Model (scale target)	PREDICT, STDDEV
Generalized Linear Mixed Model (categorical target)	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Generalized Linear Mixed Model (scale target)	PREDICT
Ordinal Multinomial Regression	PREDICT, PROBABILITY, PROBABILITY (category), CONFIDENCE
Cox Regression	PREDICT, CUMHAZARD
Nearest Neighbor (scale target)	PREDICT, STDDEV, NEIGHBOR, NEIGHBOR(K), DISTANCE, DISTANCE(K)
Nearest Neighbor (categorical target)	PREDICT, PROBABILITY, CONFIDENCE, NEIGHBOR, NEIGHBOR(K), DISTANCE, DISTANCE(K)

- For the Binary Logistic Regression, Multinomial Logistic Regression, and Naive Bayes models, the value returned by the CONFIDENCE function is identical to that returned by the PROBABILITY function.
- For the K-Means model, the value returned by the CONFIDENCE function is the least distance.
- For tree and ruleset models, the confidence can be interpreted as an adjusted probability of the predicted category and is always less than the value given by PROBABILITY. For these models, the confidence value is more reliable than the value given by PROBABILITY.
- For neural network models, the confidence provides a measure of whether the predicted category is much more likely than the second-best predicted category.
- For Ordinal Multinomial Regression and Generalized Linear Model, the PROBABILITY function is supported when the target variable is binary.
- For nearest neighbor models without a target variable, the available functions are NEIGHBOR and DISTANCE.

For information on applying scoring functions from a model, see [Scoring expressions](#).

## **NAME Subcommand**

NAME specifies the model handle name. The rules for valid model handle names are the same as for IBM® SPSS® Statistics variable names with the addition of the \$ character as an allowed first character. The model handle name should be unique for each model.

## **FILE keyword**

The FILE keyword is used to specify the external model file that you want to refer to by the model handle.

- File specifications should be enclosed in quotation marks.
- Fully qualified paths are recommended to avoid ambiguity.

## ***OPTIONS subcommand***

Use `OPTIONS` to control the treatment of missing data.

## ***MISSING keyword***

The `MISSING` keyword controls the treatment of missing values, encountered during the scoring process, for the predictor variables defined in the model. A missing value in the context of scoring refers to one of the following:

- A predictor contains no value. For numeric fields (variables), this means the system-missing value. For string fields, this means a null string.
- The value has been defined as user-missing, in the model, for the given predictor. Values defined as user-missing in the active dataset, but not in the model, are not treated as missing values in the scoring process.
- The predictor is categorical and the value is not one of the categories defined in the model.

**SYSMIS** *Return the system-missing value when scoring a case with a missing value.*

**SUBSTITUTE** *Use value substitution when scoring cases with missing values. This is the default.*

The method for determining a value to substitute for a missing value depends on the type of predictive model:

- **IBM® SPSS® Statistics models.** For independent variables in linear regression (`REGRESSION` command) and discriminant (`DISCRIMINANT` command) models, if mean value substitution for missing values was specified when building and saving the model, then this mean value is used in place of the missing value in the scoring computation, and scoring proceeds. If the mean value is not available, then `APPLYMODEL` and `STRAPPLYMODEL` return the system-missing value.
- **IBM® SPSS® AnswerTree® models & TREE command models.** For the CHAID and Exhaustive CHAID algorithms, the biggest child node is selected for a missing split variable. The biggest child node is determined by the algorithm to be the one with the largest population among the child nodes using learning sample cases. For C&RT and QUEST algorithms, surrogate split variables (if any) are used first. (Surrogate splits are splits that attempt to match the original split as closely as possible using alternate predictors.) If no surrogate splits are specified or all surrogate split variables are missing, the biggest child node is used.
- **IBM® SPSS® Modeler models.** Linear regression models are handled as described under SPSS Statistics models. Logistic regression models are handled as described under Logistic Regression models. C&R Tree models are handled as described for C&RT models under AnswerTree models.
- **Logistic Regression models.** For covariates in logistic regression models, if a mean value of the predictor was included as part of the saved model, then this mean value is used in place of the missing value in the scoring computation, and scoring proceeds. If the predictor is categorical



(for example, a factor in a logistic regression model), or if the mean value is not available, then APPLYMODEL and STRAPPLYMODEL return the system-missing value.

### **Example**

```
MODEL HANDLE NAME=twostep1 FILE='twostep1.mml'  
  /OPTIONS MISSING=SYSMIS.
```

- In this example, missing values encountered during scoring give rise to system-missing results.

## **MAP subcommand**

Use MAP to map a set of variable names from the input model to a different set of variable names in the active dataset. Both the VARIABLES and MODELVARIABLES keywords must be included. MODELVARIABLES is used to specify the list of variable names from the model that are to be mapped. VARIABLES is used to specify the list of target variable names in the active dataset.

- Both variable lists must contain the same number of names.
- No validation is performed against the current active file dictionary when the MODEL HANDLE command is processed. Errors associated with incorrect target variable names or variable data type mismatch are signaled when an APPLYMODEL or STRAPPLYMODEL transformation is processed.

### **Example**

```
MODEL HANDLE NAME=creditmod1 FILE='credit1.mml'  
  /MAP VARIABLES=agecat curdebt  
    MODELVARIABLES=age debt.
```

- In this example, the variable *age* from the model file is mapped to the variable *agecat* in the active dataset. Likewise, the variable *debt* from the model file is mapped to the variable *curdebt* in the active dataset.

# ***MODEL LIST***

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Release History***

Release 13.0

- Command introduced.

## ***Example***

```
MODEL LIST.
```

## ***Overview***

MODEL LIST produces a list, in pivot table format, of the existing model handles (see MODEL HANDLE on p. 1247). The listing includes the handle name, the type of predictive model (for example, NOMREG) associated with the model handle, the external XML model file associated with the model handle, and the method (specified on the MODEL HANDLE command) for handling cases with missing values.

## ***Basic specification***

The basic specification is simply MODEL LIST. There are no additional specifications.

## ***Operations***

- The MODEL LIST command lists only the handles created in the current working session.

# MODEL NAME

```
MODEL NAME [model name] ['model label']
```

## Example

```
MODEL NAME PLOT1 'PLOT OF THE OBSERVED SERIES'.
```

## Overview

MODEL NAME specifies a model name and label for the next procedure in the session.

### Basic Specification

The specification on MODEL NAME is a name, a label, or both.

- The default model name is *MOD\_n*, where *n* increments by 1 each time an unnamed model is created. This default is in effect if it is not changed on the MODEL NAME command, or if the command is not specified. There is no default label.

### Syntax Rules

- If both a name and label are specified, the name must be specified first.
- Only one model name and label can be specified on the command.
- The model name must be unique. The name can contain up to eight characters and must begin with a letter (*A–Z*).
- The model label can contain up to 60 characters and must be specified in quotes.

### Operations

- MODEL NAME is executed at the next model-generating procedure.
- If the MODEL NAME command is used more than once before a procedure, the last command is in effect.
- If a duplicate model name is specified, the default *MOD\_n* name will be used instead.
- *MOD\_n* reinitializes at the start of every session and when the READ MODEL command is specified (see READ MODEL). If any models in the active dataset are already named *MOD\_n*, those numbers are skipped when new *MOD\_n* names are assigned.
- The following procedures can generate models that can be named with the MODEL NAME command: AREG, ARIMA, EXSMOOTH, SEASON, and SPECTRA in the Forecasting add-on module; ACF, CASEPLOT, CCF, CURVEFIT, PACF, PLOT, and TSLOT in the Core system; and WLS and 2SLS in the Regression add-on module.

## Example

```
MODEL NAME CURVE1 'First CURVEFIT model'.
```

---

*MODEL NAME*

```
CURVEFIT Y1 .  
CURVEFIT Y2 .  
CURVEFIT Y3  
    /APPLY 'CURVE1' .
```

- In this example, the model name *CURVE1* and the label *First CURVEFIT model* are assigned to the first *CURVEFIT* command.
- The second *CURVEFIT* command has no *MODEL NAME* command before it, so it is assigned the name *MOD\_n*, where *n* is the next unused integer in the sequence.
- The third *CURVEFIT* command applies the model named *CURVE1* to the series *Y3*. This model is named *MOD\_m*, where  $m = n + 1$ .

# MRSETS

```
MRSETS

/MDGROUP NAME= setname {LABEL= 'label'      }
                        {LABELSOURCE=VARLABEL}
CATEGORYLABELS={VARLABELS      }
                {COUNTEDVALUES}
VARIABLES= varlist
VALUE= {value  }
       {'chars' }

/MCGROUP NAME= setname LABEL= 'label'
            VARIABLES= varlist

/DELETE NAME= {[setlist]}
             {ALL      }

/DISPLAY NAME= {[setlist]}
              {ALL      }
```

The set name must begin with a \$ and follow IBM® SPSS® Statistics variable naming conventions.

Square brackets shown in the DELETE and DISPLAY subcommands are required if one or more set names is specified, but not with the keyword ALL.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- LABELSOURCE keyword introduced on MDGROUP subcommand.
- CATEGORYLABELS keyword introduced on MDGROUP subcommand.

## **Example**

```
MRSETS
/MDGROUP NAME=$mltnews LABEL='News sources'
VARIABLES=news5 news4 news3 news2 news1
VALUE=1
/DISPLAY NAME=[$mltnews].
```

```
MRSETS
/MCGROUP NAME=$mltcars
LABEL='Car maker, most recent car'
VARIABLES=car1 car2 car3
/DISPLAY NAME=[$mltcars].
```

## Overview

The MRSETS command defines and manages multiple response sets. The set definitions are saved in the IBM® SPSS® Statistics data file, so they are available whenever the file is in use. Multiple response sets can be used in the GGRAPH and CTABLES (Tables option) commands. Two types of multiple response sets can be defined:

- Multiple dichotomy (MD) groups combine variables so that each variable becomes a category in the group. For example, take five variables that ask for *yes/no* responses to the questions: Do you get news from the Internet? Do you get news from the radio? Do you get news from television? Do you get news from news magazines? Do you get news from newspapers? These variables are coded 1 for *yes* and 0 for *no*. A multiple dichotomy group combines the five variables into a single variable with five categories in which a respondent could be counted zero to five times, depending on how many of the five elementary variables contain a 1 for that respondent. It is not required that the elementary variables be dichotomous. If the five elementary variables had the values 1 for *regularly*, 2 for *occasionally*, and 3 for *never*, it would still be possible to create a multiple dichotomy group that counts the variables with 1's and ignores the other responses.
- Multiple category (MC) groups combine variables that have identical categories. For example, suppose that instead of having five *yes/no* questions for the five news sources, there are three variables, each coded 1 = *Internet*, 2 = *radio*, 3 = *television*, 4 = *magazines*, and 5 = *newspapers*. For each variable, a respondent could select one of these values. In a multiple category group based on these variables, a respondent could be counted zero to three times, once for each variable for which he or she selected a news source. For this sort of multiple response group, it is important that all of the source variables have the same set of values and value labels and the same missing values.

The MRSETS command also allows you to delete sets and to display information about the sets in the data file.

## Syntax Conventions

The following conventions apply to the MRSETS command:

- All subcommands are optional, but at least one must be specified.
- Subcommands can be issued more than once in any order.
- Within a subcommand, attributes can be specified in any order. If an attribute is specified more than once, the last instance is honored.
- Equals signs are required where shown in the syntax diagram.
- Square brackets are required where shown in the syntax diagram.
- The TO convention and the ALL keyword are honored in variable lists.

## MDGROUP Subcommand

```
/MDGROUP NAME= setname {LABEL= 'label'      }
                        {LABELSOURCE=VARLABEL}
CATEGORYLABELS={VARLABELS      }
                {COUNTEDVALUES}
VARIABLES= varlist
```

```
VALUE= {value }
        { 'chars' }
```

The MDGROUP subcommand defines or modifies a multiple dichotomy set. A name, variable list, and value must be specified. Optionally, you can control assignment of set and category labels.

<b>NAME</b>	<i>The name of the multiple dichotomy set.</i> The name must follow IBM® SPSS® Statistics variable naming conventions and begin with a \$. If the name refers to an existing set, the set definition is overwritten.
<b>LABEL</b>	<i>The label for the set.</i> The label must be quoted and cannot be wider than the limit for variable labels. By default, the set is unlabeled. LABEL and LABELSOURCE are mutually exclusive.
<b>LABELSOURCE</b>	<i>Use the variable label for the first variable in the set with a defined variable label as the set label.</i> If none of the variables in the set have defined variable labels, the name of the first variable in the set is used as the set label. LABELSOURCE is an alternative to LABEL and is only available with CATEGORYLABELS=COUNTEDVALUES.
<b>CATEGORYLABELS = [VARLABELS COUNTEDVALUES]</b>	<i>Use variable labels or value labels of the counted values as category labels for the set.</i> VARLABELS uses the defined variable labels (or variable names for variables without defined variable labels) as the set category labels. This is the default. COUNTEDVALUES uses the defined value labels of the counted values as the set category labels. The counted value for each variable must have a defined value label and the labels must be unique (the value label for the counted value must be different for each variable).
<b>VARIABLES</b>	<i>The list of elementary variables that define the set.</i> Variables must be of the same type (numeric or string). At least two variables must be specified.
<b>VALUE</b>	<i>The value that indicates presence of a response.</i> This is also referred to as the “counted” value. If the set type is numeric, the counted value must be an integer. If the set type is string, the counted value, after trimming trailing blanks, cannot be wider than the narrowest elementary variable.

Elementary variables need not have variable labels, but because variable labels are used as value labels for categories of the MD variable, a warning is issued if two or more variables of an MD set have the same variable label. A warning is also issued if two or more elementary variables use different labels for the counted value—for example, if it is labeled *Yes* for Q1 and *No* for Q2. When checking for label conflicts, case is ignored.

## **MCGROUP Subcommand**

```
/MCGROUP NAME= setname LABEL= 'label'
          VARIABLES= varlist
```

The MCGROUP subcommand defines or modifies a multiple category group. A name and variable list must be specified. Optionally, a label can be specified for the set.

<b>NAME</b>	<i>The name of the multiple category set.</i> The name must follow IBM® SPSS® Statistics variable naming conventions and begin with a \$. If the name refers to an existing set, the set definition is overwritten.
<b>LABEL</b>	<i>The label for the set.</i> The label must be quoted and cannot be wider than the limit for variable labels. By default, the set is unlabeled.
<b>VARIABLES</b>	<i>The list of elementary variables that define the set.</i> Variables must be of the same type (numeric or string). At least two variables must be specified.

The elementary variables need not have value labels, but a warning is issued if two or more elementary variables have different labels for the same value. When checking for label conflicts, case is ignored.

## ***DELETE Subcommand***

```
/DELETE NAME= {[setlist]}  
              {ALL }
```

The DELETE subcommand deletes one or more set definitions. If one or more set names is given, the list must be enclosed in square brackets. ALL can be used to delete all sets; it is not enclosed in brackets.

## ***DISPLAY Subcommand***

```
/DISPLAY NAME= {[setlist]}  
              {ALL }
```

The DISPLAY subcommand creates a table of information about one or more sets. If one or more set names is given, the list must be enclosed in square brackets. ALL can be used to refer to all sets; it is not enclosed in brackets.



# MULT RESPONSE

MULT RESPONSE is available in the Statistics Base option.

```
MULT RESPONSE†  
  
  /GROUPS=groupname['label'] (varlist ({value1,value2}))  
      {value  
      }  
  ...[groupname...]  
  
  /VARIABLES=varlist (min,max)  [varlist... ]  
  
  /FREQUENCIES=varlist  
      }  
  
  /TABLES=varlist BY varlist... [BY varlist] [(PAIRED)]  
      [/varlist BY...]  
  
  [/MISSING={TABLE**} [INCLUDE]]  
      {MDGROUP}  
      {MRGROUP}  
  
  [/FORMAT={LABELS**} {TABLE** } [DOUBLE]]  
      {NOLABELS} {CONDENSE}  
      {ONEPAGE }  
  
  [/BASE={CASES** }]  
      {RESPONSES}  
  
  [/CELLS={COUNT**} [ROW] [COLUMN] [TOTAL] [ALL]]
```

†A minimum of two subcommands must be used: at least one from the pair GROUPS or VARIABLES and one from the pair FREQUENCIES or TABLES.

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
MULT RESPONSE GROUPS=MAGS (TIME TO STONE (2))  
  /FREQUENCIES=MAGS.
```

## Overview

MULT RESPONSE displays frequencies and optional percentages for multiple-response items in univariate tables and multivariate crosstabulations. Another procedure that analyzes multiple-response items is TABLES, which has most, but not all, of the capabilities of MULT RESPONSE. TABLES has special formatting capabilities that make it useful for presentations.

Multiple-response items are questions that can have more than one value for each case. For example, the respondent may have been asked to circle all magazines read within the last month in a list of magazines. You can organize multiple-response data in one of two ways for use in the program. For each possible response, you can create a variable that can have one of two values, such as 1 for *no* and 2 for *yes*; this is the multiple-dichotomy method. Alternatively, you can

estimate the maximum number of possible answers from a respondent and create that number of variables, each of which can have a value representing one of the possible answers, such as 1 for *Time*, 2 for *Newsweek*, and 3 for *PC Week*. If an individual did not give the maximum number of answers, the extra variables receive a missing-value code. This is the multiple-response or multiple-category method of coding answers.

To analyze the data entered by either method, you combine variables into groups. The technique depends on whether you have defined multiple-dichotomy or multiple-response variables. When you create a multiple-dichotomy group, each component variable with at least one *yes* value across cases becomes a category of the group variable. When you create a multiple-response group, each value becomes a category and the program calculates the frequency for a particular value by adding the frequencies of all component variables with that value. Both multiple-dichotomy and multiple-response groups can be crosstabulated with other variables in `MULT RESPONSE`.

### **Options**

**Cell Counts and Percentages.** By default, crosstabulations include only counts and no percentages. You can request row, column, and total table percentages using the `CELLS` subcommand. You can also base percentages on responses instead of respondents using `BASE`.

**Format.** You can suppress the display of value labels and request condensed format for frequency tables using the `FORMAT` subcommand.

### **Basic Specification**

The subcommands required for the basic specification fall into two groups: `GROUPS` and `VARIABLES` name the elements to be included in the analysis; `FREQUENCIES` and `TABLES` specify the type of table display to be used for tabulation. The basic specification requires at least one subcommand from each group:

- `GROUPS` defines groups of multiple-response items to be analyzed and specifies how the component variables will be combined.
- `VARIABLES` identifies all individual variables to be analyzed.
- `FREQUENCIES` requests frequency tables for the groups and/or individual variables specified on `GROUPS` and `VARIABLES`.
- `TABLES` requests crosstabulations of groups and/or individual variables specified on `GROUPS` and `VARIABLES`.

### **Subcommand Order**

- The basic subcommands must be used in the following order: `GROUPS`, `VARIABLES`, `FREQUENCIES`, and `TABLES`. Only one set of basic subcommands can be specified.
- All basic subcommands must precede all optional subcommands. Optional subcommands can be used in any order.

### **Operations**

- Empty categories are not displayed in either frequency tables or crosstabulations.

- If you define a multiple-response group with a very wide range, the tables require substantial amounts of workspace. If the component variables are sparsely distributed, you should recode them to minimize the workspace required.
- MULT RESPONSE stores category labels in the workspace. If there is insufficient space to store the labels after the tables are built, the labels are not displayed.

### **Limitations**

- The component variables must have integer values. Non-integer values are truncated.
- A maximum of 100 existing variables named or implied by GROUPS and VARIABLES together.
- A maximum of 20 groups defined on GROUPS.
- A maximum of 32,767 categories for a multiple-response group or an individual variable.
- A maximum of 10 table lists on TABLES.
- A maximum of 5 dimensions per table.
- A maximum of 100 groups and variables named or implied on FREQUENCIES and TABLES together.
- A maximum of 200 non-empty rows and 200 non-empty columns in a single table.

## **GROUPS Subcommand**

GROUPS defines both multiple-dichotomy and multiple-response groups.

- Specify a name for the group and an optional label, followed by a list of the component variables and the value or values to be used in the tabulation.
- Enclose the variable list in parentheses and enclose the values in an inner set of parentheses following the last variable in the list.
- The label for the group is optional and can be up to 40 characters in length, including imbedded blanks. Quotes around the label are not required.
- To define a multiple-dichotomy group, specify only one tabulating value (the value that represents *yes*) following the variable list. Each component variable becomes a value of the group variable, and the number of cases that have the tabulating value becomes the frequency. If there are no cases with the tabulating value for a given component variable, that variable does not appear in the tabulation.
- To define a multiple-response group, specify two values following the variable list. These are the minimum and maximum values of the component variables. The group variable will have the same range of values. The frequency for each value is tabulated across all component variables in the list.
- You can use any valid variable name for the group except the name of an existing variable specified on the same MULT RESPONSE command. However, you can reuse a group name on another MULT RESPONSE command.
- The group names and labels exist only during MULT RESPONSE and disappear once MULT RESPONSE has been executed. If group names are referred to in other procedures, an error results.

- For a multiple-dichotomy group, the category labels come from the variable labels defined for the component variables.
- For a multiple-response group, the category labels come from the value labels for the first component variable in the group. If categories are missing for the first variable but are present for other variables in the group, you must define value labels for the missing categories. (You can use the `ADD VALUE LABELS` command to define extra value labels.)

**Example**

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
  /FREQUENCIES=MAGS.
```

- The `GROUPS` subcommand creates a multiple-dichotomy group named *MAGS*. The variables between and including *TIME* and *STONE* become categories of *MAGS*, and the frequencies are cases with the value 2 (indicating *yes, read the magazine*) for the component variables.
- The group label is *MAGAZINES READ*.

**Example**

```
MULT RESPONSE GROUPS=PROBS 'PERCEIVED NATIONAL PROBLEMS'
  (PROB1 TO PROB3 (1,9))
  /FREQUENCIES=PROBS.
```

- The `GROUPS` subcommand creates the multiple-response group *PROBS*. The component variables are the existing variables between and including *PROB1* and *PROB3*, and the frequencies are tabulated for the values 1 through 9.
- The frequency for a given value is the number of cases that have that value in any of the variables *PROB1* to *PROB3*.

## VARIABLES Subcommand

`VARIABLES` specifies existing variables to be used in frequency tables and crosstabulations. Each variable is followed by parentheses enclosing a minimum and a maximum value, which are used to allocate cells for the tables for that variable.

- You can specify any numeric variable on `VARIABLES`, but non-integer values are truncated.
- If `GROUPS` is also specified, `VARIABLES` follows `GROUPS`.
- To provide the same minimum and maximum for each of a set of variables, specify a variable list followed by a range specification.
- The component variables specified on `GROUPS` can be used in frequency tables and crosstabulations, but you must specify them again on `VARIABLES`, along with a range for the values. You do not have to respecify the component variables if they will not be used as individual variables in any tables.

**Example**

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
  /VARIABLES SEX(1,2) EDUC(1,3)
  /FREQUENCIES=MAGS SEX EDUC.
```

- The `VARIABLES` subcommand names the variables `SEX` and `EDUC` so that they can be used in a frequencies table.

### Example

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
/VARIABLES=EDUC (1,3) TIME (1,2).
/TABLES=MAGS BY EDUC TIME.
```

- The variable `TIME` is used in a group and also in a table.

## FREQUENCIES Subcommand

`FREQUENCIES` requests frequency tables for groups and individual variables. By default, a frequency table contains the count for each value, the percentage of responses, and the percentage of cases. For another method of producing frequency tables for individual variables, see the `FREQUENCIES` procedure.

- All groups must be created by `GROUPS`, and all individual variables to be tabulated must be named on `VARIABLES`.
- You can use the keyword `TO` to imply a set of group or individual variables. `TO` refers to the order in which variables are specified on the `GROUPS` or `VARIABLES` subcommand.

### Example

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
/FREQUENCIES=MAGS.
```

- The `FREQUENCIES` subcommand requests a frequency table for the multiple-dichotomy group `MAGS`, tabulating the frequency of the value 2 for each of the component variables `TIME` to `STONE`.

### Example

```
MULT RESPONSE
GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
PROBS 'PERCEIVED NATIONAL PROBLEMS' (PROB1 TO PROB3 (1,9))
MEMS 'SOCIAL ORGANIZATION MEMBERSHIPS' (VFW AMLEG ELKS (1))
/VARIABLES SEX(1,2) EDUC(1,3)
/FREQUENCIES=MAGS TO MEMS SEX EDUC.
```

- The `FREQUENCIES` subcommand requests frequency tables for `MAGS`, `PROBS`, `MEMS`, `SEX`, and `EDUC`.
- You cannot specify `MAGS TO EDUC` because `SEX` and `EDUC` are individual variables, and `MAGS`, `PROBS`, and `MEMS` are group variables.

## TABLES Subcommand

`TABLES` specifies the crosstabulations to be produced by `MULT RESPONSE`. Both individual variables and group variables can be tabulated together.

- The first list defines the rows of the tables; the next list (following `BY`) defines the columns. Subsequent lists following `BY` keywords define control variables, which produce subtables. Use the keyword `BY` to separate the dimensions. You can specify up to five dimensions (four `BY` keywords) for a table.
- To produce more than one table, name one or more variables for each dimension of the tables. You can also specify multiple table lists separated by a slash. If you use the keyword `TO` to imply a set of group or individual variables, `TO` refers to the order in which groups or variables are specified on the `GROUPS` or `VARIABLES` subcommand.
- If `FREQUENCIES` is also specified, `TABLES` follows `FREQUENCIES`.
- The value labels for columns are displayed on three lines with eight characters per line. To avoid splitting words, reverse the row and column variables, or redefine the variable or value labels (depending on whether the variables are multiple-dichotomy or multiple-response variables).

**Example**

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
/VARIABLES=EDUC (1,3)/TABLES=EDUC BY MAGS.
```

- The `TABLES` subcommand requests a crosstabulation of variable `EDUC` by the multiple-dichotomy group `MAGS`.

**Example**

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
MEMS 'SOCIAL ORGANIZATION MEMBERSHIPS' (VFW AMLEG ELKS (1))
/VARIABLES EDUC (1,3)/TABLES=MEMS MAGS BY EDUC.
```

- The `TABLES` subcommand specifies two crosstabulations—`MEMS` by `EDUC` and `MAGS` by `EDUC`.

**Example**

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
/VARIABLES SEX (1,2) EDUC (1,3)
/TABLES=MAGS BY EDUC SEX/EDUC BY SEX/MAGS BY EDUC BY SEX.
```

- The `TABLES` subcommand uses slashes to separate three table lists. It produces two tables from the first table list (`MAGS` by `EDUC` and `MAGS` by `SEX`) and one table from the second table list (`EDUC` by `SEX`). The third table list produces separate tables for each sex (`MAGS` by `EDUC` for male and for female).

**Example**

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
PROBS 'NATIONAL PROBLEMS MENTIONED' (PROB1 TO PROB3 (1,9))
/TABLES=MAGS BY PROBS.
```

- The `TABLES` subcommand requests a crosstabulation of the multiple-dichotomy group `MAGS` with the multiple-response group `PROBS`.

## PAIRED Keyword

When `MULT RESPONSE` crosstabulates two multiple-response groups, by default it tabulates each variable in the first group with each variable in the second group and sums the counts for each cell. Thus, some responses can appear more than once in the table. Use `PAIRED` to pair the first variable in the first group with the first variable in the second group, the second variable in the first group with the second variable in the second group, and so on.

- The keyword `PAIRED` is specified in parentheses on the `TABLES` subcommand following the last variable named for a specific table list.
- When you request paired crosstabulations, the order of the component variables on the `GROUPS` subcommand determines the construction of the table.
- Although the tables can contain individual variables and multiple-dichotomy groups in a paired table request, only variables within multiple-response groups are paired.
- `PAIRED` also applies to a multiple-response group used as a control variable in a three-way or higher-order table.
- Paired tables are identified in the output by the label *PAIRED GROUP*.
- Percentages in paired tables are always based on responses rather than cases.

### Example

```
MULT RESPONSE GROUPS=PSEX 'SEX OF CHILD' (P1SEX P2SEX P3SEX (1,2))
/PAGE 'AGE OF ONSET OF PREGNANCY' (P1AGE P2AGE P3AGE (1,4))
/TABLES=PSEX BY PAGE (PAIRED).
```

- The `PAIRED` keyword produces a paired crosstabulation of *PSEX* by *PAGE*, which is a combination of the tables *P1SEX* by *P1AGE*, *P2SEX* by *P2AGE*, and *P3SEX* by *P3AGE*.

### Example

```
MULT RESPONSE GROUPS=PSEX 'SEX OF CHILD' (P1SEX P2SEX P3SEX (1,2))
PAGE 'AGE OF ONSET OF PREGNANCY' (P1AGE P2AGE P3AGE (1,4))
/VARIABLES=EDUC (1,3)
/TABLES=PSEX BY PAGE BY EDUC (PAIRED).
```

- The `TABLES` subcommand pairs only *PSEX* with *PAGE*. *EDUC* is not paired because it is an individual variable, not a multiple-response group.

## CELLS Subcommand

By default, `MULT RESPONSE` displays cell counts but not percentages in crosstabulations. `CELLS` requests percentages for crosstabulations.

- If you specify one or more keywords on `CELLS`, `MULT RESPONSE` displays cell counts plus the percentages you request. The count cannot be eliminated from the table cells.

<b>COUNT</b>	<i>Cell counts.</i> This is the default if you omit the <code>CELLS</code> subcommand.
<b>ROW</b>	<i>Row percentages.</i>

<b>COLUMN</b>	<i>Column percentages.</i>
<b>TOTAL</b>	<i>Two-way table total percentages.</i>
<b>ALL</b>	<i>Cell counts, row percentages, column percentages, and two-way table total percentages. This is the default if you specify the CELLS subcommand without keywords.</i>

**Example**

```
MULT RESPONSE GROUPS=MAGS 'MAGAZINES READ' (TIME TO STONE (2))
/VARIABLES=SEX (1,2) (EDUC (1,3)
/TABLES=MAGS BY EDUC SEX
/CELLS=ROW COLUMN.
```

- The CELLS subcommand requests row and column percentages in addition to counts.

**BASE Subcommand**

BASE lets you obtain cell percentages and marginal frequencies based on responses rather than respondents. Specify one of two keywords:

<b>CASES</b>	<i>Base cell percentages on cases. This is the default if you omit the BASE subcommand and do not request paired tables. You cannot use this specification if you specify PAIRED on TABLE.</i>
<b>RESPONSES</b>	<i>Base cell percentages on responses. This is the default if you request paired tables.</i>

**Example**

```
MULT RESPONSE GROUPS=PROBS 'NATIONAL PROBLEMS MENTIONED'
(PROB1 TO PROB3 (1,9))/VARIABLES=EDUC (1,3)
/TABLES=EDUC BY PROBS
/CELLS=ROW COLUMN
/BASE=RESPONSES.
```

- The BASE subcommand requests marginal frequencies and cell percentages based on responses.

**MISSING Subcommand**

MISSING controls missing values. Its minimum specification is a single keyword.

- By default, MULT RESPONSE deletes cases with missing values on a table-by-table basis for both individual variables and groups. In addition, values falling outside the specified range are not tabulated and are included in the missing category. Thus, specifying a range that excludes missing values is equivalent to the default missing-value treatment.
- For a multiple-dichotomy group, a case is considered missing by default if none of the component variables contains the tabulating value for that case. The keyword MDGROUP overrides the default and specifies listwise deletion for multiple-dichotomy groups.
- For a multiple-response group, a case is considered missing by default if none of the components has valid values falling within the tabulating range for that case. Thus, cases with missing or excluded values on some (but not all) of the components of a group are included in



tabulations of the group variable. The keyword `MRGROUP` overrides the default and specifies listwise deletion for multiple-response groups.

- You can use `INCLUDE` with `MDGROUP`, `MRGROUP`, or `TABLE`. The user-missing value is tabulated if it is included in the range specification.

<b>TABLE</b>	<i>Exclude missing values on a table-by-table basis.</i> Missing values are excluded on a table-by-table basis for both component variables and groups. This is the default if you omit the <code>MISSING</code> subcommand.
<b>MDGROUP</b>	<i>Exclude missing values listwise for multiple-dichotomy groups.</i> Cases with missing values for any component dichotomy variable are excluded from the tabulation of the multiple-dichotomy group.
<b>MRGROUP</b>	<i>Exclude missing values listwise for multiple-response groups.</i> Cases with missing values for any component variable are excluded from the tabulation of the multiple-response group.
<b>INCLUDE</b>	<i>Include user-missing values.</i> User-missing values are treated as valid values if they are included in the range specification on the <code>GROUPS</code> or <code>VARIABLES</code> subcommands.

### Example

```
MULT RESPONSE GROUPS=FINANCL 'FINANCIAL PROBLEMS MENTIONED'
(FINPROB1 TO FINPROB3 (1,3))
SOCIAL 'SOCIAL PROBLEMS MENTIONED' (SOCPROB1 TO SOCPROB4 (4,9))
/VARIABLES=EDUC (1,3)
/TABLES=EDUC BY FINANCL SOCIAL
/MISSING=MRGROUP.
```

- The `MISSING` subcommand indicates that a case will be excluded from counts in the first table if any of the variables in the group `FINPROB1` to `FINPROB3` has a missing value or a value outside the range 1 to 3. A case is excluded from the second table if any of the variables in the group `SOCPROB1` to `SOCPROB4` has a missing value or value outside the range 4 to 9.

## FORMAT Subcommand

`FORMAT` controls table formats. The minimum specification on `FORMAT` is a single keyword.

Labels are controlled by two keywords:

<b>LABELS</b>	<i>Display value labels in frequency tables and crosstabulations.</i> This is the default.
<b>NOLABELS</b>	<i>Suppress value labels in frequency tables and crosstabulations for multiple-response variables and individual variables.</i> You cannot suppress the display of variable labels used as value labels for multiple-dichotomy groups.

The following keywords apply to the format of frequency tables:

<b>DOUBLE</b>	<i>Double spacing for frequency tables.</i> By default, <code>MULT RESPONSE</code> uses single spacing.
<b>TABLE</b>	<i>One-column format for frequency tables.</i> This is the default if you omit the <code>FORMAT</code> subcommand.

- CONDENSE**            *Condensed format for frequency tables.* This option uses a three-column condensed format for frequency tables for all multiple-response groups and individual variables. Labels are suppressed. This option does not apply to multiple-dichotomy groups.
- ONEPAGE**            *Conditional condensed format for frequency tables.* Three-column condensed format is used if the resulting table would not fit on a page. This option does not apply to multiple-dichotomy groups.

**Example**

```
MULT RESPONSE GROUPS=PROBS 'NATIONAL PROBLEMS MENTIONED'  
  (PROB1 TO PROB3 (1,9))/VARIABLES=EDUC (1,3)  
  /FREQUENCIES=EDUC PROBS  
  /FORMAT=CONDENSE.
```

- The `FORMAT` subcommand specifies condensed format, which eliminates category labels and displays the categories in three parallel sets of columns, each set containing one or more rows of categories (rather than displaying one set of columns aligned vertically down the page).

# MULTIPLE CORRESPONDENCE

MULTIPLE CORRESPONDENCE is available in the Categories option.

```
MULTIPLE CORRESPONDENCE [/VARIABLES =] varlist

  /ANALYSIS = varlist [[WEIGHT={1**}]
                    {n }

  /DISCRETIZATION = [varlist [[{GROUPING } ] [{NCAT={7} } ] [DISTR={NORMAL }]]]
                    {n } {UNIFORM}
                    {EQINTV={n}}
                    {RANKING }
                    {MULTIPLYING}

  /MISSING = [{varlist} [{PASSIVE**}] [{MODEIMPU**}]]
            {ALL* } {ACTIVE } {EXTRACAT }
            {MODEIMPU**}
            {EXTRACAT }
            {LISTWISE }

  /SUPPLEMENTARY = [OBJECT(objlist)] [VARIABLE(varlist)]

  /CONFIGURATION = [{INITIAL}] ('filename')
                  {FIXED }

  /DIMENSION = {2**}
              {n }

  /NORMALIZATION = {VPRINCIPAL**}
                  {OPRINCIPAL }
                  {SYMMETRICAL }
                  {INDEPENDENT }
                  {n }

  /MAXITER = {100**}
            {n }

  /CRITITER = {.00001**}
            {value }

  /PRINT = [CORR**] [DESCRIP**{(varlist)}] [HISTORY][DISCRIM**] [NONE]
          [OBJECT[({varname})varlist]] [OCORR] [QUANT{(varlist)}]

  /PLOT = [BIPLOT[({varlist})[(varlist)][(n)]]
         [CATEGORY (varlist)[(n)]]
         [JOINTCAT{(varlist)}[(n)]]
         [DISCRIM**{(varlist)}[(n)]]
         [NONE]
         [OBJECT**{(varlist)}[(n)]]
         [RESID(varlist[({1**})][(n)]]
         {n }
         [TRANS(varlist[({1**})][(n)]]
         {n }
         [NDIM(value,value)]

  /SAVE = [TRDATA[({TRA }[(n)]]] [OBJECT[({OBSCO }[(n)]]] ]
         {rootname} {rootname}

  /OUTFILE = {DISCRDATA('filename')} {OBJECT('filename')} {TRDATA('filename')}
```

\*\* Default if subcommand is omitted

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

**Release History**

Release 13.0

- Command introduced.

**Overview**

MULTIPLE CORRESPONDENCE (Multiple Correspondence Analysis; also known as homogeneity analysis) quantifies nominal (categorical) data by assigning numerical values to the cases (objects) and categories, such that in the low-dimensional representation of the data, objects within the same category are close together and objects in different categories are far apart. Each object is as close as possible to the category points of categories that apply to the object. In this way, the categories divide the objects into homogeneous subgroups. Variables are considered homogeneous when they classify objects in the same categories into the same subgroups.

**Basic Specification**

The basic specification is the command MULTIPLE CORRESPONDENCE with the VARIABLES and ANALYSIS subcommands.

**Syntax Rules**

- The VARIABLES and ANALYSIS subcommands always must appear.
- All subcommands can appear in any order.
- For the first subcommand after the procedure name, a slash is accepted, but not required.
- Variables specified in the ANALYSIS subcommand must be found in the VARIABLES subcommand.
- Variables specified in the SUPPLEMENTARY subcommand must be found in the ANALYSIS subcommand.

**Operations**

- If the same subcommand is repeated, it causes a syntax error and the procedure terminates.

**Limitations**

- MULTIPLE CORRESPONDENCE operates on category indicator variables. The category indicators should be positive integers. You can use the DISCRETIZATION subcommand to convert fractional value variables and string variables into positive integers. If DISCRETIZATION is not specified, fractional value variables are automatically converted into positive integers by grouping them into seven categories (or into the number of distinct values of the variable if this number is less than seven) with a close-to-normal distribution, and string variables are automatically converted into positive integers by ranking.
- In addition to system-missing values and user-defined missing values, MULTIPLE CORRESPONDENCE treats category indicator values less than 1 as missing. If one of the values of a categorical variable has been coded 0 or some negative value and you want to treat it as a valid category, use the COMPUTE command to add a constant to the values of that

variable such that the lowest value will be 1. You can also use the `RANKING` option of the `DISCRETIZATION` subcommand for this purpose, except for variables you want to treat as numerical, since the spacing of the categories will not be maintained.

- There must be at least three valid cases.
- Split-File has no implications for `MULTIPLE CORRESPONDENCE`.

## Example

```
MULTIPLE CORRESPONDENCE /VARIABLES = TEST1 TEST2 TEST3 TO TEST6 TEST7 TEST8
  /ANALYSIS = TEST1 TO TEST2 (WEIGHT=2)
              TEST3 TO TEST5
              TEST6 TEST7
              TEST8
  /DISCRETIZATION = TEST1 (GROUPING NCAT=5 DISTR=UNIFORM)
                  TEST6 (GROUPING) TEST8 (MULTIPLYING)
  /MISSING = TEST5 (ACTIVE) TEST6 (ACTIVE EXTRACAT) TEST8 (LISTWISE)
  /SUPPLEMENTARY = OBJECT(1 3) VARIABLE(TEST1)
  /CONFIGURATION = ('iniconf.sav')
  /DIMENSION = 2
  /NORMALIZATION = VPRINCIPAL
  /MAXITER = 150
  /CRITITER = .000001
  /PRINT = DESCRIP DISCRIM CORR QUANT(TEST1 TO TEST3) OBJECT
  /PLOT = TRANS(TEST2 TO TEST5) OBJECT(TEST2 TEST3)
  /SAVE = TRDATA OBJECT
  /OUTFILE = TRDATA('/data/trans.sav') OBJECT('/data/obs.sav').
```

- `VARIABLES` defines variables. The keyword `TO` refers to the order of the variables in the working data file.
- The `ANALYSIS` subcommand defines variables used in the analysis. It is specified that `TEST1` and `TEST2` have a weight of 2 (for the other variables, `WEIGHT` is not specified and thus they have the default weight value of 1).
- `DISCRETIZATION` specifies that `TEST6` and `TEST8`, which are fractional value variables, are discretized: `TEST6` by recoding into seven categories with a normal distribution (default because unspecified) and `TEST8` by “multiplying”. `TEST1`, which is a categorical variable, is recoded into five categories with a close to uniform distribution.
- `MISSING` specifies that objects with missing values on `TEST5` and `TEST6` are included in the analysis: missing values on `TEST5` are replaced with the mode (default if not specified) and missing values on `TEST6` are treated as an extra category. Objects with a missing value on `TEST8` are excluded from the analysis. For all other variables, the default is in effect; that is, missing values (not objects) are excluded from the analysis.
- `CONFIGURATION` specifies `iniconf.sav` as the file containing the coordinates of a configuration that is to be used as the initial configuration (default because unspecified).
- `DIMENSION` specifies the number of dimensions to be 2. This is the default, so this subcommand could be omitted here.
- The `NORMALIZATION` subcommand specifies optimization of the association between variables. This is the default, so this subcommand could be omitted here.
- `MAXITER` specifies the maximum number of iterations to be 150 instead of the default value of 100.
- `CRITITER` sets the convergence criterion to a value smaller than the default value.

- `PRINT` specifies descriptives, discrimination measures, and correlations (all default), and quantifications for `TEST1` to `TEST3`, and the object scores.
- `PLOT` is used to request transformation plots for the variables `TEST2` to `TEST5`, an object points plot labeled with the categories of `TEST2`, and an object points plot labeled with the categories of `TEST3`.
- The `SAVE` subcommand adds the transformed variables and the object scores to the working data file.
- The `OUTFILE` subcommand writes the transformed data to a data file called `trans.sav` and the object scores to a data file called `obs.sav`, both in the directory `/data`.

## Options

**Discretization.** You can use the `DISCRETIZATION` subcommand to discretize fractional value variables or to recode categorical variables.

**Missing data.** You can specify the treatment of missing data per variable with the `MISSING` subcommand.

**Supplementary objects and variables.** You can specify objects and variables that you want to treat as supplementary.

**Read configuration.** `MULTIPLE CORRESPONDENCE` can read a configuration from a file through the `CONFIGURATION` subcommand. This configuration can be used as the initial configuration or as a fixed configuration in which to fit variables.

**Number of dimensions.** You can specify how many dimensions `MULTIPLE CORRESPONDENCE` should compute.

**Normalization.** You can specify one of five different options for normalizing the objects and variables.

**Tuning the algorithm.** You can control the values of algorithm-tuning parameters with the `MAXITER` and `CRITITER` subcommands.

**Optional output.** You can request optional output through the `PRINT` subcommand.

**Optional plots.** You can request a plot of object points, transformation plots per variable, plots of category points per variable, or a joint plot of category points for specified variables. Other plot options include residuals plots, a biplot, and a plot of discrimination measures.

**Writing discretized data, transformed data, and object scores.** You can write the discretized data, the transformed data, and the object scores to outfiles for use in further analyses.

**Saving transformed data and object scores.** You can save the transformed variables and the object scores in the working data file.

## VARIABLES Subcommand

`VARIABLES` specifies the variables that may be analyzed in the current `MULTIPLE CORRESPONDENCE` procedure.

- The `VARIABLES` subcommand is required. The actual keyword `VARIABLES` can be omitted.

- At least two variables must be specified, except if the `CONFIGURATION` subcommand with the `FIXED` keyword is used.
- The keyword `TO` on the `VARIABLES` subcommand refers to the order of variables in the working data file. (Note that this behavior of `TO` is different from that in the `varlist` in the `ANALYSIS` subcommand.)

## ***ANALYSIS Subcommand***

`ANALYSIS` specifies the variables to be used in the computations, and the variable weight for each variable or variable list. `ANALYSIS` also specifies supplementary variables; no weight can be specified for supplementary variables.

- At least two variables must be specified, except if the `CONFIGURATION` subcommand with the `FIXED` keyword is used.
- All the variables on `ANALYSIS` must be specified on the `VARIABLES` subcommand.
- The `ANALYSIS` subcommand is required.
- The keyword `TO` in the variable list honors the order of variables in the `VARIABLES` subcommand.
- Variable weights are indicated by the keyword `WEIGHT` in parentheses following the variable or variable list.

**WEIGHT**                    *Specifies the variable weight.* The default value is 1. If `WEIGHT` is specified for supplementary variables, this is ignored (but with a syntax warning).

## ***DISCRETIZATION Subcommand***

`DISCRETIZATION` specifies fractional value variables you want to discretize. Also, you can use `DISCRETIZATION` for ranking or for two ways of recoding categorical variables.

- A string variable's values are always converted into positive integers, by assigning category indicators according to the ascending alphanumeric order. `DISCRETIZATION` for string variables applies to these integers.
- When the `DISCRETIZATION` subcommand is omitted, or when the `DISCRETIZATION` subcommand is used without a `varlist`, fractional value variables are converted into positive integers by grouping them into seven categories (or into the number of distinct values of the variable if this number is less than seven) with a close-to-normal distribution.
- When no specification is given for variables in a `varlist` following `DISCRETIZATION`, these variables are grouped into seven categories (or into the number of distinct values of the variable if this number is less than seven) with a close-to-normal distribution.
- In `MULTIPLE CORRESPONDENCE` a system-missing value, user-defined missing values, and values less than 1 are considered to be missing values (see next section). However, in discretizing a variable, values less than 1 are considered to be valid values and are thus

included in the discretization process. System-missing values and user-defined missing values are excluded.

<b>GROUPING</b>	<i>Recode into the specified number of categories or recode intervals of equal size into categories.</i>
<b>RANKING</b>	<i>Rank cases. Rank 1 is assigned to the case with the smallest value on the variable.</i>
<b>MULTIPLYING</b>	<i>Multiplying the standardized values (z-scores) of a fractional value variable by 10, rounding, and adding a value such that the lowest value is 1.</i>

### **GROUPING Keyword**

<b>NCAT</b>	<i>Recode into ncat categories. When NCAT is not specified, the number of categories is set to seven (or the number of distinct values of the variable if this number is less than seven).</i>
<b>EQINTV</b>	<i>Recode intervals of equal size into categories. The interval size must be specified (there is no default value). The resulting number of categories depends on the interval size.</i>

### **NCAT Keyword**

NCAT has the keyword **DISTR**, which has the following keywords:

<b>NORMAL</b>	<i>Normal distribution. This is the default when DISTR is not specified.</i>
<b>UNIFORM</b>	<i>Uniform distribution.</i>

### **MISSING Subcommand**

In **MULTIPLE CORRESPONDENCE**, system-missing values, user-defined missing values, and values less than 1 are treated as missing values. However, in discretizing a variable, values less than 1 are considered as valid values. The **MISSING** subcommand allows you to indicate how to handle missing values for each variable.

<b>PASSIVE</b>	<i>Exclude missing values on a variable from analysis. This is the default applicable to all variables, when the MISSING subcommand is omitted or specified without variable names or keywords. Also, any variable which is not included in the subcommand gets this specification. Passive treatment of missing values means that, in optimizing the quantification of a variable, only objects with non-missing values on the variable are involved and that only the non-missing values of variables contribute to the solution. Thus, when PASSIVE is specified, missing values do not affect the analysis. If an object has only missing values, and for all variables the MISSING option is passive, the object will be handled as a supplementary object. If on the PRINT subcommand, correlations are requested and passive treatment of missing values is specified for a variable, the missing values have to be imputed. For the correlations of the original variables, missing values on a variable are imputed with the most frequent category (mode) of the variable.</i>
<b>ACTIVE</b>	<i>Impute missing values. You can choose to use mode imputation, or to consider objects with missing values on a variable as belonging to the same category and impute missing values with an extra category indicator.</i>
<b>LISTWISE</b>	<i>Exclude cases with missing values on the specified variable(s). The cases used in the analysis are cases without missing values on the variable(s) specified. Also, any variable that is not included in the subcommand gets this specification.</i>



- The ALL keyword may be used to indicate all variables. If it is used, it must be the only variable specification.
- A mode or extracat imputation is done before listwise deletion.

### **PASSIVE Keyword**

<b>MODEIMPU</b>	<i>Impute missing values on a variable with the mode of the quantified variable. This is the default.</i>
<b>EXTRACAT</b>	<i>Impute missing values on a variable with the quantification of an extra category. This implies that objects with a missing value are considered to belong to the same (extra) category.</i>

*Note:* With passive treatment of missing values, imputation only applies to correlations and is done afterwards. Thus the imputation has no effect on the quantification or the solution.

### **ACTIVE Keyword**

<b>MODEIMPU</b>	<i>Impute missing values on a variable with the most frequent category (mode). When there are multiple modes, the smallest category indicator is used. This is the default.</i>
<b>EXTRACAT</b>	<i>Impute missing values on a variable with an extra category indicator. This implies that objects with a missing value are considered to belong to the same (extra) category.</i>

*Note:* With active treatment of missing values, imputation is done before the analysis starts, and thus will affect the quantification and the solution.

### **SUPPLEMENTARY Subcommand**

The SUPPLEMENTARY subcommand specifies the objects or/and variables that you want to treat as supplementary. Supplementary variables must be found in the ANALYSIS subcommand. You can not weight supplementary objects and variables (specified weights are ignored). For supplementary variables, all options on the MISSING subcommand can be specified except LISTWISE.

<b>OBJECT</b>	<i>Objects you want to treat as supplementary are indicated with an object number list in parentheses following OBJECT. The keyword TO is allowed. The OBJECT specification is not allowed when CONFIGURATION = FIXED.</i>
<b>VARIABLE</b>	<i>Variables you want to treat as supplementary are indicated with a variable list in parentheses following VARIABLE. The keyword TO is allowed and honors the order of variables in the VARIABLES subcommand. The VARIABLE specification is ignored when CONFIGURATION = FIXED, for in that case all the variables in the ANALYSIS subcommand are automatically treated as supplementary variables.</i>

## **CONFIGURATION Subcommand**

The CONFIGURATION subcommand allows you to read data from a file containing the coordinates of a configuration. The first variable in this file should contain the coordinates for the first dimension, the second variable should contain the coordinates for the second dimension, and so forth.

<b>INITIAL('filename')</b>	Use the configuration in the specified file as the starting point of the analysis.
<b>FIXED('filename')</b>	<i>Fit variables in the fixed configuration found in the specified file.</i> The variables to fit in should be specified on the ANALYSIS subcommand but will be treated as supplementary variables. The SUPPLEMENTARY subcommand will be ignored. Also, variable weights will be ignored.

## **DIMENSION Subcommand**

DIMENSION specifies the number of dimensions you want MULTIPLE CORRESPONDENCE to compute.

- If you do not specify the DIMENSION subcommand, MULTIPLE CORRESPONDENCE computes a two dimensional solution.
- DIMENSION is followed by an integer indicating the number of dimensions.
- The maximum number of dimensions is the smaller of a) the number of observations minus 1 and b) the total number of valid variable levels (categories) minus the number of variables if there are no variables with missing values to be treated as passive. If there are variables with missing values to be treated as passive, the maximum number of dimensions is the smaller of a) the number of observations minus 1 and b) the total number of valid variable levels (categories) minus the larger of c) 1 and d) the number of variables without missing values to be treated as passive.
- The maximum number of dimensions is the smaller of the number of observations minus 1 and the total number of valid variable levels (categories) minus the number of variables without missing values.
- MULTIPLE CORRESPONDENCE adjusts the number of dimensions to the maximum if the specified value is too large.
- The minimum number of dimensions is 1.

## **NORMALIZATION Subcommand**

The NORMALIZATION subcommand specifies one of five options for normalizing the object scores and the variables.

- Only one normalization method can be used in a given analysis.

<b>VPRINCIPAL</b>	<i>Optimize the association between variables.</i> With VPRINCIPAL, the categories are in the centroid of the objects in the particular categories. VPRINCIPAL is the default if the NORMALIZATION subcommand is
-------------------	--

	not specified. This is useful when you are primarily interested in the association between the variables.
<b>OPRINCIPAL</b>	<i>Optimize distances between objects.</i> This is useful when you are primarily interested in differences or similarities between the objects.
<b>SYMMETRICAL</b>	<i>Use this normalization option if you are primarily interested in the relation between objects and variables.</i>
<b>INDEPENDENT</b>	<i>Use this normalization option if you want to examine distances between objects and associations between variables separately.</i>

The fifth method allows the user to specify any real value in the closed interval  $[-1, 1]$ . A value of 1 is equal to the **OPRINCIPAL** method, a value of 0 is equal to the **SYMMETRICAL** method, and a value of  $-1$  is equal to the **VPRINCIPAL** method. By specifying a value greater than  $-1$  and less than 1, the user can spread the eigenvalue over both objects and variables. This method is useful for making a tailor-made biplot. If the user specifies a value outside of this interval, the procedure issues a syntax error message and terminates.

## **MAXITER Subcommand**

**MAXITER** specifies the maximum number of iterations **MULTIPLE CORRESPONDENCE** can go through in its computations.

- If **MAXITER** is not specified, the maximum number of iterations is 100.
- The specification on **MAXITER** is a positive integer indicating the maximum number of iterations. There is no uniquely predetermined (that is, hard-coded) maximum for the value that can be used.

## **CRITITER Subcommand**

**CRITITER** specifies a convergence criterion value. **MULTIPLE CORRESPONDENCE** stops iterating if the difference in fit between the last two iterations is less than the **CRITITER** value.

- If **CRITITER** is not specified, the convergence value is 0.00001.
- The specification on **CRITITER** is any positive value.

## **PRINT Subcommand**

The Model Summary statistics (Cronbach's alpha and the variance accounted for) and the **HISTORY** statistics (the variance accounted for, the loss, and the increase in variance accounted for) for the last iteration are always displayed. That is, they cannot be controlled by the **PRINT** subcommand. The **PRINT** subcommand controls the display of optional additional output. The output of the **MULTIPLE CORRESPONDENCE** procedure is always based on the transformed variables. However, the correlations of the original variables can be requested as well by the keyword **OCORR**.

The default keywords are **DESCRIP**, **DISCRIM**, and **CORR**. That is, the three keywords are in effect when the **PRINT** subcommand is omitted or when the **PRINT** subcommand is given without any keywords. Note that when some keywords are specified, the default is nullified and only the keywords specified become in effect. If a keyword that cannot be followed by a

`varlist` is duplicated or if a contradicting keyword is encountered, then the later one silently becomes effective (in case of a contradicting use of `NONE`, only the keywords following `NONE` are effective). For example,

```
/PRINT <=> /PRINT = DESCRIP DISCRIM CORR
```

```
/PRINT = DISCRIM DISCRIM <=> /PRINT = DISCRIM
```

```
/PRINT = DISCRIM NONE CORR <=> /PRINT = CORR
```

If a keyword that can be followed by a `varlist` is duplicated, it will cause a syntax error and the procedure will terminate. For example, `/PRINT = QUANT QUANT` is a syntax error.

The following keywords can be specified:

<b>DESCRIP(varlist)</b>	<i>Descriptive statistics (frequencies, missing values, and mode).</i> The variables in the <code>varlist</code> must be specified on the <code>VARIABLES</code> subcommand, but need not appear on the <code>ANALYSIS</code> subcommand. If <code>DESCRIP</code> is not followed by a <code>varlist</code> , Descriptives tables are displayed for all the variables in the <code>varlist</code> on the <code>ANALYSIS</code> subcommand.
<b>DISCRIM</b>	<i>Discrimination measures per variable and per dimension.</i>
<b>QUANT(varlist)</b>	<i>Category quantifications (centroid coordinates), mass, inertia of the categories, contribution of the categories to the inertia of the dimensions, and contribution of the dimensions to the inertia of the categories.</i> Any variable in the <code>ANALYSIS</code> subcommand may be specified in parentheses after <code>QUANT</code> . If <code>QUANT</code> is not followed by a <code>varlist</code> , Quantification tables are displayed for all variables in the <code>varlist</code> on the <code>ANALYSIS</code> subcommand.  <i>Note:</i> the coordinates and contributions (including the mass and inertia) are displayed in separate layers of the pivot table output, with the coordinates shown by default. To display the contributions, activate (double-click) on the table and select Contributions from the Layer dropdown list.
<b>HISTORY</b>	<i>History of iterations.</i> For each iteration, the variance accounted for, the loss, and the increase in variance accounted for are shown.
<b>CORR</b>	<i>Correlations of the transformed variables, and the eigenvalues of this correlation matrix.</i> Correlation tables are displayed for each set of quantifications, thus there are <i>ndim</i> (the number of dimensions in the analysis) correlation tables; the <i>i</i> th table contains the correlations of the quantifications of dimension <i>i</i> , <i>i</i> = 1, ..., <i>ndim</i> . For variables with missing values specified to be treated as <code>PASSIVE</code> on the <code>MISSING</code> subcommand, the missing values are imputed according to the specification on the <code>PASSIVE</code> keyword (if nothing is specified, mode imputation is used).
<b>OCORR</b>	<i>Correlations of the original variables, and the eigenvalues of this correlation matrix.</i> For variables with missing values specified to be treated as <code>PASSIVE</code> on the <code>MISSING</code> subcommand, the missing values are imputed with the variable mode.

<b>OBJECT((varname)varlist)</b>	<i>Object scores (component scores) and, in separate table, mass, inertia of the objects, contribution of the objects to the inertia of the dimensions, and contribution of the dimensions to the inertia of the objects.</i> Following the keyword, a <code>varlist</code> can be given in parentheses to display variables (category indicators) along with the object scores. If you want to use a variable to label the objects, this variable must occur in parenthesis as the first variable in the <code>varlist</code> . If no labeling variable is specified, the objects are labeled with case numbers. The variables to display along with the object scores and the variable to label the objects must be specified on the <code>VARIABLES</code> subcommand but need not appear on the <code>ANALYSIS</code> subcommand. If no <code>varlist</code> is given, only the object scores are displayed.
<b>NONE</b>	<i>No optional output is displayed.</i> The only output shown is the Model Summary and the <code>HISTORY</code> statistics for the last iteration.

The keyword `TO` in a variable list can only be used with variables that are in the `ANALYSIS` subcommand, and `TO` applies only to the order of the variables in the `ANALYSIS` subcommand. For variables that are in the `VARIABLES` subcommand but not in the `ANALYSIS` subcommand, the keyword `TO` cannot be used. For example, if `/VARIABLES = v1 TO v5` and the `ANALYSIS` subcommand has `/ANALYSIS v2 v1 v4`, then `/PLOT OBJECT(v1 TO v4)` will give two object plots, one labeled with `v1` and one labeled with `v4`. (`/PLOT OBJECT(v1 TO v4 v2 v3 v5)` will give objects plots labeled with `v1`, `v2`, `v3`, `v4`, and `v5`).

## ***PLOT Subcommand***

The `PLOT` subcommand controls the display of plots. The default keywords are `OBJECT` and `DISCRIM`. That is, the two keywords are in effect when the `PLOT` subcommand is omitted, or when the `PLOT` subcommand is given without any keyword. If a keyword is duplicated (for example, `/PLOT = RESID RESID`), then it will cause a syntax error and the procedure will terminate. If the keyword `NONE` is used together with other keywords (for example, `/PLOT = RESID NONE DISCRIM`), then only the keywords following `NONE` are effective. That is, when keywords contradict, the later one overwrites the earlier ones.

- All the variables to be plotted must be specified in the `ANALYSIS` subcommand.
- If the variable list following the keywords `CATEGORIES`, `TRANS`, and `RESID` is empty, then it will cause a syntax error and the procedure will terminate.
- The variables in the `varlist` for labeling the object points following `OBJECT` and `BIPLOT` must be specified on the `VARIABLES` subcommand but need not appear on the `ANALYSIS` subcommand. This means that variables not included in the analysis can still be used to label plots.
- The keyword `TO` in a variable list can only be used with variables that are in the `ANALYSIS` subcommand, and `TO` applies only to the order of the variables in the `ANALYSIS` subcommand. For variables that are in the `VARIABLES` subcommand but not in the `ANALYSIS` subcommand, the keyword `TO` cannot be used. For example, if `/VARIABLES = v1 TO v5` and `/ANALYSIS is v2 v1 v4`, then `/PLOT OBJECT(v1 TO v4)` will give two object plots, one

labeled with  $v1$  and one labeled with  $v4$ . (/PLOT OBJECT( $v1$  TO  $v4$   $v2$   $v3$   $v5$ )) will give objects plots labeled with  $v1$ ,  $v2$ ,  $v3$ ,  $v4$ , and  $v5$ ).

- For multidimensional plots, all of the dimensions in the solution are produced in a matrix scatterplot if the number of dimensions in the solution is greater than two and the `NDIM` keyword is not specified; if the specified number of dimensions is 2, a scatterplot is produced.

The following keywords can be specified:

<b>OBJECT (varlist)(n)</b>	<i>Plots of the object points.</i> Following the keyword, a list of variables in parentheses can be given to indicate that plots of object points labeled with the categories of the variables should be produced (one plot for each variable). If the variable list is omitted, a plot labeled with case numbers is produced.
<b>CATEGORY(varlist)(n)</b>	<i>Plots of the category points (centroid coordinates).</i> A list of variables must be given in parentheses following the keyword. Categories are in the centroids of the objects in the particular categories.
<b>DISCRIM(varlist)(n)</b>	<i>Plot of the discrimination measures.</i> <code>DISCRIM</code> can be followed by a <code>varlist</code> to select the variables to include in the plot. If the variable list is omitted, a plot including all variables is produced.
<b>TRANS(varlist)(n)</b>	<i>Transformation plots per variable (optimal category quantifications against category indicators).</i> Following the keyword, a list of variables in parentheses must be given. Each variable can be followed by a number of dimensions in parentheses to indicate you want to display $p$ residual plots, one for each of the first $p$ dimensions. If the number of dimensions is not specified, a plot for the first dimension is produced.
<b>RESID(varlist)(n)</b>	<i>Plot of residuals per variable (approximation against optimal category quantifications).</i> Following the keyword, a list of variables in parentheses must be given. Each variable can be followed by a number of dimensions in parentheses to indicate you want to display $p$ residual plots, one for each of the first $p$ dimensions. If the number of dimensions is not specified, a plot for the first dimension is produced.
<b>BI-PLOT((varlist))(varlist)(n)</b>	<i>Plot of objects and variables (centroids).</i> When <code>NORMALIZATION = INDEPENDENT</code> , this plot is incorrect and therefore not available. <code>BIPLOT</code> can be followed by a <code>varlist</code> in double parentheses to select the variables to include in the plot. If this variable list is omitted, a plot including all variables is produced. Following <code>BIPLOT</code> or <code>BIPLOT( (varlist) )</code> , a list of variables in single parentheses can be given to indicate that plots with objects labeled with the categories of the variables should be produced (one plot for each variable). If this variable list is omitted, a plot with objects labeled with case numbers is produced.
<b>JOINTCAT(varlist)(n)</b>	<i>Joint plot of the category points for the variables in the varlist.</i> If no <code>varlist</code> is given, the category points for all variables are displayed.
<b>NONE</b>	<i>No plots.</i>

- For all of the keywords except `TRANS` and `NONE`, the user can specify an optional parameter  $l$  in parentheses after the variable list in order to control the global upper boundary of variable name/label and value label lengths in the plot. Note that this boundary is applied uniformly to all variables in the list. The label length parameter  $l$  can take any non-negative integer less than or equal to the applicable maximum length (64 for variable names, 255 for variable labels, and 60 for value labels). If  $l = 0$ , names/values instead of variable/value labels are displayed to indicate variables/categories. If  $l$  is not specified, `MULTIPLE CORRESPONDENCE` assumes that each variable name/label and value label at its full length is displayed. If  $l$  is an integer larger than the applicable maximum, then we reset it to the applicable maximum but do

not issue a warning. If a positive value of  $l$  is given but if some or all of the variables/category values do not have labels, then for those variables/values the names/values themselves are used as the labels.

In addition to the plot keywords, the following can be specified:

**NDIM(value,value)**            *Dimension pairs to be plotted.* NDIM is followed by a pair of values in parentheses. If NDIM is not specified or if NDIM is specified without parameter values, a matrix scatterplot including all dimensions is produced.

- The first value (an integer that can range from 1 to the number of dimensions in the solution minus 1) indicates the dimension that is plotted against higher dimensions.
- The second value (an integer that can range from 2 to the number of dimensions in the solution) indicates the highest dimension to be used in plotting the dimension pairs.
- The NDIM specification applies to all requested multidimensional plots.

## SAVE Subcommand

The SAVE subcommand is used to add the transformed variables (category indicators replaced with optimal quantifications) and the object scores to the working data file.

- Excluded cases are represented by a dot (the `sysmis` symbol) on every saved variable.

**TRDATA**            *Transformed variables.* Missing values specified to be treated as passive are represented by a dot. Following TRDATA, a rootname, and the number of dimensions to be saved can be specified in parentheses (if the number of dimensions is not specified, all dimensions are saved).

**OBJECT**            *Object scores.*

MULTIPLE CORRESPONDENCE adds three numbers. The first number uniquely identifies the source variable names, the middle number corresponds to the dimension number, and the last number uniquely identifies the MULTIPLE CORRESPONDENCE procedures with the successfully executed SAVE subcommands. Only one rootname can be specified and it can contain up to three characters. If more than one rootname is specified, the first rootname is used; if a rootname contains more than three characters, the first three characters are used at most.

- If a rootname is not specified for TRDATA, rootname *TRA* is used to automatically generate unique variable names. The formula is *ROOTNAME<sub>k</sub>\_m\_n*, where  $k$  increments from 1 to identify the source variable names by using the source variables' position numbers in the ANALYSIS subcommand,  $m$  increments from 1 to identify the dimension number, and  $n$  increments from 1 to identify the MULTIPLE CORRESPONDENCE procedures with the successfully executed SAVE subcommands for a given data file in a continuous session. For example, with two variables specified on ANALYSIS and 2 dimensions to save, the first set of default names, if they do not exist in the data file, would be *TRAI\_1\_1*, *TRAI\_2\_1*, *TRA2\_1\_1*, *TRA2\_2\_1*. The next set of default names, if they do not exist in the data file, would be *TRAI\_1\_2*, *TRAI\_2\_2*, *TRA2\_1\_2*, *TRA2\_2\_2*. However, if, for example, *TRAI\_1\_2* already exists in the data file, then the default names should be attempted as *TRAI\_1\_3*, *TRAI\_2\_3*, *TRA2\_1\_3*, *TRA2\_2\_3*. That is, the last number increments to the next available integer.

- Following **OBJECT**, a rootname and the number of dimensions can be specified in parentheses (if the number of dimensions is not specified, all dimensions are saved), to which **MULTIPLE CORRESPONDENCE** adds two numbers separated by the underscore symbol (**\_**). The first number corresponds to the dimension number. The second number uniquely identifies the **MULTIPLE CORRESPONDENCE** procedures with the successfully executed **SAVE** subcommands. Only one rootname can be specified, and it can contain up to five characters. If more than one rootname is specified, the first rootname is used; if a rootname contains more than five characters, the first five characters are used at most.
- If a rootname is not specified for **OBJECT**, the rootname *OBSCO* is used to automatically generate unique variable names. The formula is *ROOTNAME<sub>m</sub>\_n*, where *m* increments from 1 to identify the dimension number and *n* increments from 1 to identify the **MULTIPLE CORRESPONDENCE** procedures with the successfully executed **SAVE** subcommands for a given data file in a continuous session. For example, if 2 dimensions are specified following **OBJECT**, the first set of default names, if they do not exist in the data file, would be *OBSCO1\_1*, *OBSCO2\_1*. The next set of default names, if they do not exist in the data file, would be *OBSCO1\_2*, *OBSCO2\_2*. However, if, for example, *OBSCO2\_2* already exists in the data file, then the default names should be attempted as *OBSCO1\_3*, *OBSCO2\_3*. That is, the second number increments to the next available integer.
- Variable labels are created automatically. They are shown in the Notes table and can also be displayed in the Data Editor window.
- If the number of dimensions is not specified, the **SAVE** subcommand saves all dimensions.

## ***OUTFILE Subcommand***

The **OUTFILE** subcommand is used to write the discretized data, transformed data (category indicators replaced with optimal quantifications), and the object scores to an external IBM® SPSS® Statistics data file or previously declared dataset. Excluded cases are represented by a dot (the **sysmis** symbol) on every saved variable.

<b>DISCRDATA</b> ( <b>'savfile'</b> 'dataset')	<i>Discretized data.</i>
<b>TRDATA</b> ( <b>'savfile'</b> 'dataset')	<i>Transformed variables.</i> Missing values specified to be treated as passive are represented by a dot.
<b>OBJECT</b> ( <b>'savfile'</b> 'dataset')	<i>Object scores.</i>

Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files. The names should be different for each of the keywords.

- The active dataset, in principle, should not be replaced by this subcommand, and the asterisk (\*) file specification is not supported. This strategy also helps prevent **OUTFILE** interference with the **SAVE** subcommand.



# MULTIPLE IMPUTATION

MULTIPLE IMPUTATION is available in the Missing Values option.

```
MULTIPLE IMPUTATION varlist
[/IMPUTE [METHOD={AUTO** }]
          {FCS
           }
          {MONOTONE}
          {NONE
           }
          [NIMPUTATIONS={5**}]
            {int}
          [SCALEMODEL = {LINEAR**}]
            {PMM
             }
          [INTERACTIONS = {NONE**}]
            {TWOWAY}
          [MAXPCTMISSING={NONE**}]
            {num
             }
          [MAXCASEDRAWS={50**}]
            {int
             }
          [MAXPARAMDRAWS={2**}]
            {int
             }
          [MAXITER={10**}]
            {int
             }
          [SINGULAR={1E-12**}]
            {num
             }
          [MAXMODELPARAM={100**}]
            {int
             }
[/CONSTRAINTS varlist ([MIN={NONE**}]
                       {num
                        }
                       [MAX={NONE**}]
                         {num
                          }
                       [RND={NO** }]
                         {num
                          }
                       [ROLE={BOTH**}])]
                       {DEP
                        }
                       {IND
                        }
[/MISSINGSUMMARIES [OVERALL**]
                   [VARIABLES [( [MAXVARS={25**}] [MINPCTMISSING={10**}])]
                               {int
                                }
                               {num
                                }
                   [PATTERNS]
                   [NONE]]
[/IMPUTATIONSUMMARIES [MODELS**] [DESCRIPTIVES[(varlist)]] [NONE]]
[/ANALYSISWEIGHT var]
[/OUTFILE [IMPUTATIONS = 'savfile' | 'dataset']]
          [FCSTITERATIONS = 'savfile' | 'dataset']
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 17.0

- Command introduced.

## Example

```
DATASET DECLARE imputedData.
MULTIPLE IMPUTATION x y z
/OUTFILE IMPUTATIONS = imputedData.
```

## Overview

The `MULTIPLE IMPUTATION` procedure performs multiple imputation of missing data values. Given a dataset containing missing values, it outputs one or more datasets in which missing values are replaced with plausible estimates. The procedure also summarizes missing values in the working dataset.

Datasets produced by the `MULTIPLE IMPUTATION` procedure can be analyzed using supported analysis procedures to obtain final (combined) parameter estimates that take into account the inherent uncertainty among the various sets of imputed values.

### Options

**Variables.** You can specify which variables to impute and specify constraints on the imputed values such as minimum and maximum values. You can also specify which variables are used as predictors when imputing missing values of other variables.

**Methods.** Three imputation methods are offered. The monotone method is an efficient method for data that have a monotone pattern of missingness. Fully conditional specification (FCS) is an iterative Markov Chain Monte Carlo (MCMC) method that is appropriate when the data have an arbitrary (monotone or nonmonotone) missing pattern. The default method (AUTO) scans the data to determine the best imputation method (monotone or FCS). For each method you can control the number of imputations.

**Output.** By default the procedure displays an overall summary of missingness in your data as well as an imputation summary and the imputation model for each variable whose values are imputed. You can obtain analysis of missing values by variable as well as tabulated patterns of missing values. If you request imputation you can obtain descriptive statistics for imputed values.

### Basic Specification

The basic specification is two or more variables and a file specification where the imputed data will be written.

- By default, the procedure imputes missing values using the `AUTO` method. Five imputations are performed.
- When imputing the default model type depends on the measurement level of the variable: for categorical variables, logistic regression is used; for scale variables linear regression is used.
- The procedure generates output that summarizes missingness in the data and summaries how values were imputed.

### Operations

- The output dataset contains the original (nonmissing) data and data for one or more imputations. Each imputation includes all of the observed data and imputed data values. The original and imputed data are stacked in the output dataset. A special variable, `Imputation_`, identifies whether a case represents original data (`Imputation_ = 0`) or imputed data (`Imputation_ = 1...m`).

- Multiple imputation datasets can be analyzed using supported analysis procedures to obtain final (combined) parameter estimates that take into account the inherent uncertainty in the various sets of imputed values. Variable *Imputation\_* must be defined as a split variable in order to obtain pooled parameter estimates.
- The procedure honors the random number generator and seed specified via the global `SET` command. Specify the same seed across invocations of the procedure if you want to be able to reproduce your imputed values.
- The procedure honors the `WEIGHT` variable. It is treated as a replication weight when summarizing missing values and estimating imputation models. Cases with negative or zero value of replication weight are ignored. Noninteger weights are rounded to the nearest integer. The procedure also accepts analysis weights (see `ANALYSISWEIGHT` subcommand).
- The procedure honors `SPLIT FILE`. A separate missing value analysis and set of imputations is produced for each combination of values of the split variables. An error occurs if imputation is requested and the input dataset has eight split variables.
- The procedure honors the `FILTER` command. Cases filtered out are ignored by the procedure.
- The procedure accepts string variables and treats them as categorical. Completely blank string values are treated as valid values, i.e., they are not replaced.
- The procedure treats both user- and system-missing values as invalid values. Both types of missing values are replaced when values are imputed and both are treated as invalid values of variables used as predictors in imputation models. User- and system-missing values are also treated as missing in analyses of missingness (counts of missing values, etc.).
- Cases that have a missing value for each analysis variable are included in analyses of missingness but are excluded from imputation. Specifically, values of such cases are not imputed and are excluded when building imputation models. The determination of which cases are completely missing is made after any variables are filtered out of the imputation model by the `MAXPCTMISSING` keyword.
- An error occurs if imputation is requested and the input dataset contains a variable named *Imputation\_*.
- An error occurs if imputation and iteration history are requested and the input dataset contains a variable named *Iteration\_*, or *SummaryStatistic\_*.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### **Syntax Rules**

- Two or more analysis variables are required.
- The `OUTFILE` subcommand is required unless imputation is turned off (`/IMPUTE METHOD=NONE`). All other subcommands are optional.
- Only a single instance of each subcommand is allowed except for the `CONSTRAINTS` command, which can be repeated.
- An error occurs if an attribute or keyword is specified more than once within a subcommand.
- Equals signs shown in the syntax chart are required.

- Subcommand names and keywords must be spelled in full.
- Empty subcommands are not allowed.

## Examples

### **Automatic Imputation**

```
DATASET DECLARE imputedData.
MULTIPLE IMPUTATION x y z
  /OUTFILE IMPUTATIONS = imputedData.
```

- The command requests multiple imputation of scale variables x, y, and z.
- By default, the procedure chooses an imputation method based on a scan of the data. Five imputations are produced by default.
- As all variables are scale, a linear regression model is used for each. When a particular variable is imputed, all other variables are treated as covariates in the imputation model if FCS is automatically chosen as the imputation method. If the monotone method is chosen, only preceding variables are used as predictors.
- The imputed data are written to a new dataset *imputedData*.
- By default, an overall summary of missingness is displayed, as well as information about how values were imputed.

### **Analysis of Missingness**

```
MULTIPLE IMPUTATION x y z
  /IMPUTE METHOD=NONE
  /MISSINGSUMMARIES VARIABLES.
```

- The command requests analysis of missing values for variables x, y, and z.
- No imputation is performed.

### **Custom Imputation Model**

```
DATASET DECLARE imputedData.
MULTIPLE IMPUTATION x y z
  /IMPUTE METHOD=FCS
  /CONSTRAINTS x y (RND=1 MIN=1)
  /CONSTRAINTS x (MAX=5)
  /CONSTRAINTS y (MAX=10)
  /OUTFILE IMPUTATIONS = imputedData.
```

- Scale variables x, y, and z are defined as analysis variables.
- Missing values are imputed using the FCS method. Variables are imputed in order of their appearance in the variable list. Variable x is imputed first, using x and z as predictors. Then y is imputed, treating x and z as predictors. Finally, z is imputed using x and y as predictors.
- CONSTRAINTS subcommands specify constraints on imputed values. Imputed values of x and y are rounded to the nearest integer and are constrained to be no less than 1. Maximum values of 5 and 10 are specified for x and y, respectively. No constraints are defined for variable z.

- Imputed data are written to a new dataset *imputedData*.
- By default, an overall summary of missingness is displayed, as well as information about how values were imputed.

## **Variable Lists**

Analysis variables are specified after the command name. Two or more variables must be specified. The `TO` and `ALL` keywords can be used to refer to multiple variables. If any variable is specified more than once, the last instance of that variable is honored.

The variable list specifies variables to impute and to include in analyses of missingness. By default, analysis variables are also used as predictors to in imputation models of other analysis variables. The lists of variables to impute and predictors can be restricted via the `CONSTRAINTS` subcommand.

### **Variable Order**

When imputing using the `FCS` and `MONOTONE` methods, variables are imputed sequentially in the order in which they are listed in the variables list. The order of the variables is ignored by the `AUTO` method except to break ties.

### **Predictors**

The set of predictors that is used when a particular variable is imputed depends on the imputation method:

- For the `FCS` method, when a variable is imputed all other analysis variables are used as predictors in the imputation model.
- When the `MONOTONE` method is used, only variables that precede the variable to be imputed in the variable list are used as predictors.
- When the `AUTO` method is used the set of predictors depends on the pattern of missingness in the data. If the data have a nonmonotone pattern of missingness, `FCS` is used and all other analysis variables are used as predictors. If the data have a monotone pattern, `MONOTONE` is used and all variables that precede the variable to be imputed are used predictors. Note that `AUTO` sorts the analysis variables to detect monotone pattern, so the actual order of variables may not correspond to their order in which they are specified in the variables list.

### **Measurement Level**

Measurement level recorded in the data dictionary is honored for each analysis variable. Measurement level determines the following:

- The default type of imputation model for variables whose values are imputed (linear regression or logistic regression).
- Whether a variable is treated as a factor (categorical) or covariate (scale) when used as a predictor in imputation models.
- Whether the variable is treated as scale or categorical in summaries of missing values.

The procedure treats ordinal and nominal variables equivalently as categorical variables.

## ***IMPUTE Subcommand***

The `IMPUTE` subcommand controls the imputation method and model. By default, the `AUTO` method is used to impute missing data values.

### ***METHOD Keyword***

The `METHOD` keyword specifies the imputation method. Specify one of the following options:

<b>AUTO</b>	<i>Automatic method.</i> This is the default. Chooses the best method based on a scan of the data. The monotone method is used when the data have a monotone pattern of missingness; otherwise, if the data have a nonmonotone pattern, <code>FCS</code> is used. The <code>AUTO</code> method internally sorts the analysis variables in order from the least missing to most missing to detect a monotone pattern of missingness if it exists; thus, order of variables in the variable list is ignored when <code>AUTO</code> is used.
<b>MONOTONE</b>	<i>Monotone method.</i> This is a noniterative method that can be used only when the data have a monotone pattern of missingness. A monotone pattern exists when you can order the variables such that, if a variable has a nonmissing value, all preceding variables also have nonmissing values. The monotone method imputes missing values sequentially in the order specified in the variable list. An error occurs if the input data do not have a monotone pattern for the order in which variables are specified.
<b>FCS</b>	<i>Fully conditional specification method.</i> This is an iterative Markov chain Monte Carlo (MCMC) method that can be used when the pattern of missing data is arbitrary (monotone or nonmonotone). In each iteration, the <code>FCS</code> method sequentially imputes missing values in the order specified in the variable list.
<b>NONE</b>	<i>None.</i> No imputation is performed. Choose this option if you want analyses of missingness only. A warning is issued if you turn off imputation and suppress all analyses of missing values.

### ***NIMPUTATIONS Keyword***

By default, five imputations are performed. To request a different number of imputations, specify a positive integer. As a rule of thumb, the higher the degree of missing information in the data, the more imputations that are needed to obtain reasonable efficiency relative to an infinite number of imputations.

- `NIMPUTATIONS` is ignored if imputation is turned off.

### ***SCALEMODEL Keyword***

The `AUTO`, `FCS`, and `MONOTONE` methods are multivariate methods that can be used when several variables have missing values. Each uses univariate models when imputing values of individual variables.

By default, the type of univariate model that is used depends on the measurement level of the variable whose missing values are to be imputed. Multinomial logistic regression is always used for categorical variables. For scale variables, linear regression (`LINEAR`) is used by default. Optionally you can use the predictive mean matching (`PMM`) method for scale variables. `PMM` is

a variant of linear regression that ensures that the imputed values are plausible. For `PMM` the imputed value always matches an observed value (specifically, the observed value that is closest to the value drawn by the imputation model).

- `SCALEMODEL` is honored only when the `FCS` or `MONOTONE` method is chosen explicitly. Linear regression is always used for scale variables when the `AUTO` method is requested. A warning is issued if `SCALEMODEL` is specified for the `AUTO` method.
- `SCALEMODEL` is ignored if imputation is turned off.

### ***INTERACTIONS Keyword***

By default, the imputation model for each variable includes a constant term and main effects for predictor variables. You can optionally include all possible two-way interactions among categorical predictor variables. Specify `INTERACTIONS=TWOWAY`. Interaction terms do not include scale predictors.

- `INTERACTIONS` is honored only when the `FCS` or `MONOTONE` method is chosen explicitly. Two-way interactions are not included when `AUTO` method is requested. A warning is issued if `INTERACTIONS` is specified for the `AUTO` method.
- `INTERACTIONS` is ignored if imputation is turned off or if there are fewer than two categorical predictors.

### ***MAXPCTMISSING Keyword***

By default analysis variables are imputed and used as predictors without regard to how many missing values they have provided they at least have sufficient data to estimate an imputation model. The optional `MAXPCTMISSING` keyword is used to exclude variables that have a high percentage of missing values. Specify the maximum allowable percentage of missing values as a positive number less than 100. For example, if you specify `MAXMISSINGPCT=50`, analysis variables that have more than 50% missing values are not imputed, nor are they used as predictors in imputation models.

- `MAXPCTMISSING` is ignored if imputation is turned off.

### ***Maximum Number of Draws***

If minimum or maximum values are specified for imputed values of scale variables (see `CONSTRAINTS` subcommand), the procedure attempts to draw values for a case until it finds a set of values that are within the specified ranges. By default, a maximum of 50 sets of values are drawn per case. To override the maximum, specify a positive integer for `MAXCASEDRAWS`. If a set of values is not obtained within the specified number of draws per case, the procedure draws another set of model parameters and the repeats the case-drawing process. By default, a maximum of 2 sets of model parameters are drawn. To override the maximum, specify a positive integer for `MAXPARAMDRAWS`. An error occurs if a set of values within the ranges is not obtained within the specified number of case and parameter draws.

Note that increasing these values can increase the processing time. If the procedure is taking a long time (or is unable) to find suitable draws, check the minimum and maximum values specified on the `CONSTRAINTS` subcommand to ensure they are appropriate.

- `MAXCASEDRAWS` and `MAXPARAMDRAWS` are ignored if imputation is turned off, when predictive mean matching is used, and when imputing categorical values.

### ***MAXITER Keyword***

Iteration stops when the maximum number of iterations is reached. To override the default number of iterations (10), specify a positive integer value for `MAXITER`.

- `MAXITER` is honored only if the `FCS` method is chosen explicitly. The keyword is ignored when the `AUTO` method is used.
- The `MAXITER` keyword is ignored if the monotone method is used or imputation is turned off.

### ***SINGULAR Keyword***

`SINGULAR` specifies the tolerance value used to test for singularity in univariate imputation models (linear regression, logistic regression, and predictive mean matching). The default value is  $10e-12$ . Specify a positive value.

- `SINGULAR` is honored only if the `FCS` or `MONOTONE` method is chosen explicitly. The keyword is ignored and the default singularity tolerance is used when `AUTO` method is requested.
- `SINGULAR` is ignored if imputation is turned off.

### ***MAXMODELPARAM Keyword***

`MAXMODELPARAM` specifies the maximum number of model parameters allowed when imputing any variable. If a model has more parameters than the specified limit, processing terminates with an error (no missing values are imputed). The default value is 100. Specify a positive integer value.

- `MAXMODELPARAM` is ignored if imputation is turned off.

## ***CONSTRAINTS Subcommand***

The `CONSTRAINTS` subcommand specifies constraints on imputed values. `CONSTRAINTS` specifications are ignored when imputation is turned off.

To define constraints, specify one or more analysis variables and one or more of following in parentheses:

**MIN = NONE | num**

*Minimum allowable imputed value for scale variables.* Specify a number. If an imputed value is less than the minimum, the procedure draws another value until it finds one that is greater than or equal to `MIN` or the `MAXCASEDRAWS` or `MAXPARAMDRAWS` threshold is reached (See `METHOD` subcommand). There is no default minimum. `MIN` is ignored when predictive mean matching is used or when applied to a categorical variable. An error occurs if `MIN` is greater than or equal to `MAX`. For date format variables, values must be enclosed in single or double quotes and expressed in the same date format as the defined date format for the variable.



**MAX = NONE | num**

*Maximum allowable imputed value for scale variables.* Specify a number. If an imputed value is greater than the maximum, the procedure draws another value until it finds one that is less than or equal to MAX or the MAXCASEDRAWS or MAXPARAMDRAWS threshold is reached (See METHOD subcommand). There is no default maximum. MAX is ignored when predictive mean matching is used or when applied to a categorical variable. For date format variables, values must be enclosed in single or double quotes and expressed in the same date format as the defined date format for the variable.

**RND = NO | num**

*Controls rounding of imputed values.* By default imputed values are not rounded. To round imputed values, specify a numeric multiplier value. Imputed values are rounded to the nearest integer multiple of the specified multiplier. For example, if you specify RND=1, imputed values are rounded to the nearest integer; if you specify RND=0.1, imputed values are rounded to the nearest tenth. Note that no fuzz factor is added before rounding (that is, rounding is 'fuzzless'). RND is ignored when applied to a categorical or date format variable.

**ROLE = BOTH | DEP | IND**

*Customize the set of variables to be imputed and/or treated as predictors.* By default, each analysis variable is considered as both a dependent variable and independent variable. ROLE can be used to turn off imputation for variables that you want to use as independent variables only (ROLE=IND) or to exclude variables from being used as independent variables (ROLE=DEP) and thereby make the prediction model more compact. Other CONSTRAINTS keywords are ignored for a variable if you define a variable as a predictor only.

When imputation is requested, an error is issued if ROLE=IND for all analysis variables.

*Note:* Missing values of variables that are defined as predictor-only (ROLE=IND) variables are imputed, but only for internal use (otherwise those values of the variable could not be used to impute other variables). Imputed values of such variables are not saved to the output dataset. When imputing predictor-only variables, a main-effects only model is always used; the INTERACTIONS keyword is ignored.

The variable list honors the TO and ALL keywords. Constraints are ignored when applied to a variable that is not an analysis variable.

The CONSTRAINTS subcommand can be specified more than once, and its settings are cumulative. That is, if more than one set of specifications is provided for a variable across CONSTRAINTS subcommands, the specifications are merged. When merging, if a particular keyword has different values for the same variable, the value specified in the last CONSTRAINTS subcommand for that variable is honored.

***Cumulative Effects of Multiple Constraints***

```
MULTIPLE IMPUTATION x y z
  /CONSTRAINTS x y z (RND=1 MIN=1)
  /CONSTRAINTS x      (MAX=5)
  /CONSTRAINTS y      (MIN=2 MAX=10)
  /CONSTRAINTS z      (ROLE=DEP)
  /OUTFILE IMPUTATIONS = imputedData.
```

- The first CONSTRAINTS subcommand specifies that imputed values of x, y, and z are rounded to the nearest integer and are constrained to be no less than 1.
- The second CONSTRAINTS subcommand specifies a maximum value of 5 for x.

- The third `CONSTRAINTS` subcommand specifies a maximum value of 10 for y, and overrules the previous minimum of 1 with a minimum of 2 for y.
- The fourth `CONSTRAINTS` subcommand specifies that z should be used as a dependent variable only in the imputation models.

The constraints can be summarized as follows:

Variable	Role	Minimum	Maximum	Rounding
x	Dependent and Predictor	1	5	1
y	Dependent and Predictor	2	10	1
z	Dependent only	1		1

## ***MISSINGSUMMARIES Subcommand***

The `MISSINGSUMMARIES` subcommand controls the display of missing values of analysis variables.

### ***OVERALL Keyword***

The `OVERALL` keyword displays a graph that summarizes missingness for cases, variables, and individual data (cell) values:

- Number and percent of analysis variables that have one or more missing values.
- Number and percent of cases that have one or more missing values on analysis variables.
- Number and percent of individual data values that are missing among all analysis variables and cases.

Overall summaries are displayed by default.

### ***VARIABLES Keyword***

The `VARIABLES` keyword displays a table of analysis variables sorted by percent of missing values. The table includes descriptive statistics (mean and standard deviation) for scale variables. The table is not shown by default.

You can optionally control the maximum number of variables to display (default=25) and minimum percentage missing (default=10) for a variable to be included in the display. Specify one or both in parentheses. For example, `VARIABLES (MAXVARS=50 MINPCTMISSING=25)` requests that the table display up to 50 variables that have at least 25% missing values. The set of variables that meet both criteria are displayed. For example, by default, if there are 25 analysis variables but only 5 have 10% or more missing values, the output includes only 5 variables. If there are more analysis variables than `MAXVARS` and all meet the `MINPCTMISSING` criteria, then variables with the most missing values are displayed.

- `MAXVARS` must be a positive integer.
- `MINPCTMISSING` must be a nonnegative number that is less than 100.

***PATTERNS Keyword***

The `PATTERNS` keyword displays tabulated patterns of missing values. Each pattern corresponds to a group of cases with the same pattern of incomplete and complete data on analysis variables. You can use `PATTERNS` output to determine whether monotone imputation method can be used for your data, or if not, how closely your data approximate a monotone pattern.

To display patterns, specify the `PATTERNS` keyword. The chart orders analysis variables and patterns to reveal monotonicity where it exists. Specifically, variables are ordered from left to right in increasing order of missingness. Rows (patterns) are the sorted first by the last variable (nonmissing values first, then missing values), then by the second to last variable, and so on, working from right to left.

You can use `PATTERNS` output to determine whether monotone imputation method can be used for your data, or, if not, how closely your data approximate a monotone pattern. If no nonmonotone pattern exists after reordering you can conclude that the data have a monotonic pattern when analysis variables are ordered as such. If data are not monotone, the chart shows how to achieve reach monotonicity.

***NONE Keyword***

`NONE` suppresses all `MISSINGSUMMARIES` output.

`NONE` generates an error when used with any other keyword.

***IMPUTATIONSUMMARIES Subcommand***

If imputation is requested an overall imputation summary is displayed. Contents include analysis specifications and results:

- Specifications include imputation method, number of imputations, model type for scale variables, whether two-way interactions included in models, `MAXPCTMISSING` setting, and analysis weight.
- Results include imputation method, iterations (for `FCS` method), dependent variables imputed, dependent variables excluded from imputation, and imputation sequence.

If specified, constants for analysis variables are also shown when imputation is requested. You can also obtain optional summaries of imputed data by specifying one or more of the following keywords. Imputation summary keywords are ignored if imputation is turned off.

***MODELS Keyword***

The `MODELS` keyword displays the imputation model for dependent variables and predictors. Imputation models are shown by default. Includes univariate model type, model effects, and number of values imputed.

**DESCRIPTIVES Keyword**

The `DESCRIPTIVES` keyword displays descriptive statistics for dependent variables for which values are imputed. Descriptive statistics are not shown by default.

- For scale variables the descriptive statistics include mean, count, standard deviation, min, and max for the original input data (prior to imputation), imputed values (by imputation), and complete data (original and imputed values together—by imputation).
- For categorical variables the descriptive statistics include count and percent by category for the original input data (prior to imputation), imputed values (by imputation), and complete data (original and imputed values together—by imputation).

If you specify `DESCRIPTIVES`, statistics are produced by default for each dependent variables for which values are imputed. You can optionally specify a list of dependent variables in parentheses. Descriptives are shown only for the requested variables. Non-analysis variables in the list are ignored.

**NONE Keyword**

The `NONE` keyword suppresses `IMPUTATIONSUMMARIES` output.

`NONE` generates an error if any other keyword is specified.

**ANALYSISWEIGHT Subcommand**

The `ANALYSISWEIGHT` command specifies a variable containing analysis (regression) weights. The procedure incorporates analysis weights in regression and classification models used to impute missing values. Analysis weights are also used in summaries of imputed values (e.g., mean, standard deviation, and standard error).

The analysis weight variable must be numeric. Cases with a negative or zero analysis weight are ignored.

**OUTFILE Subcommand**

The `OUTFILE` subcommand allows you to export IBM® SPSS® Statistics-format data containing imputed values or SPSS Statistics-format data containing an FCS iteration history. The `OUTFILE` subcommand is ignored if imputation is turned off.

**IMPUTATIONS Keyword**

The `IMPUTATIONS` keyword specifies an optional dataset containing imputed data. The output dataset consists of the original data and one complete dataset per imputation. All variables in the input dataset are included in the output dataset. Dictionary properties (names, labels, etc.) of existing variables are copied to the new dataset. The file also contains a new variable, *Imputation\_*, a numeric variable that indicates the imputation (0 for original data, 1..*n* for cases having imputed values).

The procedure automatically defines the *Imputation\_* variable as a split variable (see [SPLIT FILE on p. 1930](#)) when the output dataset is created. If `SPLIT FILE` is in effect when the procedure executes, the output dataset includes one set of imputations for each combination of values of split variables.

***FCSITERATIONS Keyword***

The `FCSITERATIONS` keyword specifies an optional dataset that contains iteration history data for FCS imputation.

The dataset contains means and standard deviations by iteration and imputation for each scale dependent variable for which values are imputed. You can plot the data to help assess model convergence.

- The `FCSITERATIONS` keyword is ignored if the monotone method is used. A warning is shown and the dataset is not generated if all imputed variables are categorical.

# MVA

MVA is available in the Missing Values Analysis option.

```
MVA VARIABLES= {varlist}
               {ALL      }

[/CATEGORICAL={varlist}]

[/MAXCAT={25**}]
           {n      }

[/ID={varname}]
```

## Description:

```
[/NOUNIVARIATE]

[/TTEST [PERCENT={5}] [{T  }] [{DF  }] [{PROB  }] [{COUNTS  }] [{MEANS  }]]
           {n} {NOT} {NODF} {NOPROB} {NOCOUNTS} {NOMEANS}

[/CROSSTAB [PERCENT={5}]]
           {n}

[/MISMATCH [PERCENT={5}] [NOSORT]]
           {n}

[/DPATTERN [SORT={varname}({ASCENDING })] [varname ... ]]
           {DESCENDING}
           [DESCRIBE={varlist}]

[/MPATTERN [NOSORT] [DESCRIBE={varlist}]]

[/TPATTERN [NOSORT] [DESCRIBE={varlist}] [PERCENT={1}]]
           {n}
```

## Estimation:

```
[/LISTWISE]

[/PAIRWISE]

[/EM [predicted_varlist] [WITH predictor_varlist]
    [( [TOLERANCE={0.001} ]
      {value}
      [CONVERGENCE={0.0001}]
      {value }
      [ITERATIONS={25} ]
      {n }
      [TDF=n ]
      [LAMBDA=a ]
      [PROPORTION=b ]
      [OUTFILE='file' ])]

[/REGRESSION [predicted_varlist] [WITH predictor_varlist]
    [( [TOLERANCE={0.001} ]
      {n }
      [FLIMIT={4.0} ]
      {N }
      [NPREDICTORS=number_of_predictor_variables]
      [ADDDTYPE={RESIDUAL*} ]
      {NORMAL }
      {T[({5}) ]
      {n}
      {NONE } ]
```

```
[OUTFILE='file'                                ]]]
```

\*If the number of complete cases is less than half the number of cases, the default ADDTYPE specification is NORMAL.

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### Examples

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /ID=country
  /MPATTERN DESCRIBE=region religion.
```

```
MVA VARIABLES=all
  /EM males msport WITH males msport gradrate facratio.
```

## Overview

MVA (Missing Value Analysis) describes the missing value patterns in a data file (data matrix). It can estimate the means, the covariance matrix, and the correlation matrix by using listwise, pairwise, regression, and EM estimation methods. Missing values themselves can be estimated (imputed), and you can then save the new data file.

### Options

**Categorical variables.** String variables are automatically defined as categorical. For a long string variable, only the first eight characters are used to define categories. Quantitative variables can be designated as categorical by using the CATEGORICAL subcommand.

MAXCAT specifies the maximum number of categories for any categorical variable. If any categorical variable has more than the specified number of distinct values, MVA is not executed.

**Analyzing Patterns.** For each quantitative variable, the TTEST subcommand produces a series of  $t$  tests. Values of the quantitative variable are divided into two groups, based on the presence or absence of other variables. These pairs of groups are compared using the  $t$  test.

**Crosstabulating Categorical Variables.** The CROSSTAB subcommand produces a table for each categorical variable, showing, for each category, how many nonmissing values are in the other variables and the percentages of each type of missing value.

**Displaying Patterns.** DPATTERN displays a case-by-case data pattern with codes for system-missing, user-missing, and extreme values. MPATTERN displays only the cases that have missing values and sorts by the pattern that is formed by missing values. TPATTERN tabulates the cases that have a common pattern of missing values. The pattern tables have sorting options. Also, descriptive variables can be specified.

**Labeling Cases.** For pattern tables, an ID variable can be specified to label cases.

**Suppression of Rows.** To shorten tables, the `PERCENT` keyword suppresses missing-value patterns that occur relatively infrequently.

**Statistics.** Displays of univariate, listwise, and pairwise statistics are available.

**Estimation.** `EM` and `REGRESSION` use different algorithms to supply estimates of missing values, which are used in calculating estimates of the mean vector, the covariance matrix, and the correlation matrix of dependent variables. The estimates can be saved as replacements for missing values in a new data file.

### ***Basic Specification***

The basic specification depends on whether you want to describe the missing data pattern or estimate statistics. Often, description is done first, and then, considering the results, an estimation is done. Alternatively, both description and estimation can be done by using the same `MVA` command.

**Descriptive Analysis.** A basic descriptive specification includes a list of variables and a statistics or pattern subcommand. For example, a list of variables and the subcommand `DPATTERN` would show missing value patterns for all cases with respect to the list of variables.

**Estimation.** A basic estimation specification includes a variable list and an estimation method. For example, if the `EM` method is specified, the following are estimated: the mean vector, the covariance matrix, and the correlation matrix of quantitative variables with missing values.

## ***Syntax Rules***

- A variables specification is required directly after the command name. The specification can be either a variable list or the keyword `ALL`.
- The `CATEGORICAL`, `MAXCAT`, and `ID` subcommands, if used, must be placed after the variables list and before any other subcommand. These three subcommands can be in any order.
- Any combination of description and estimation subcommands can be specified. For example, both the `EM` and `REGRESSION` subcommands can be specified in one `MVA` command.
- Univariate statistics are displayed unless the `NOUNIVARIATE` subcommand is specified. Thus, if only a list of variables is specified, with no description or estimation subcommands, univariate statistics are displayed.
- If a subcommand is specified more than once, only the last subcommand is honored.
- The following words are reserved as keywords or internal commands in the `MVA` procedure: `VARIABLES`, `SORT`, `NOSORT`, `DESCRIBE`, and `WITH`. They cannot be used as variable names in `MVA`.
- The tables *Summary of Estimated Means* and *Summary of Estimated Standard Deviations* are produced if you specify more than one way to estimate means and standard deviations. The methods include univariate (default), listwise, pairwise, `EM`, and regression. For example, these tables are produced when you specify both `LISTWISE` and `EM`.



## ***Symbols***

The symbols that are displayed in the `DPATTERN` and `MPATTERN` table cells are:

+	Extremely high value
-	Extremely low value
S	System-missing value
A	First type of user-missing value
B	Second type of user-missing value
C	Third type of user-missing value

- An extremely high value is more than 1.5 times the interquartile range above the 75th percentile, if  $(\text{number of variables}) \times n \log n \leq 150000$ , where  $n$  is the number of cases.
- An extremely low value is more than 1.5 times the interquartile range below the 25th percentile, if  $(\text{number of variables}) \times n \log n \leq 150000$ , where  $n$  is the number of cases.
- For larger files—that is,  $(\text{number of variables}) \times n \log n > 150000$ —extreme values are two standard deviations from the mean.

## ***Missing Indicator Variables***

For each variable in the variables list, a binary indicator variable is formed (internal to `MVA`), indicating whether a value is present or missing.

## ***VARIABLES Subcommand***

A list of variables or the keyword `ALL` is required.

- The order in which the variables are listed determines the default order in the output.
- If the keyword `ALL` is used, the default order is the order of variables in the active dataset.
- String variables that are specified in the variable list, whether short or long, are automatically defined as categorical. For a long string variable, only the first eight characters of the values are used to distinguish categories.
- The list of variables must precede all other subcommands.
- Multiple lists of variables are not allowed.

## ***CATEGORICAL Subcommand***

The `MVA` procedure automatically treats all string variables in the variables list as categorical. You can designate numeric variables as categorical by listing them on the `CATEGORICAL` subcommand. If a variable is designated categorical, it will be ignored if listed as a dependent or independent variable on the `REGRESSION` or `EM` subcommand.

## MAXCAT Subcommand

The `MAXCAT` subcommand sets the upper limit of the number of distinct values that each categorical variable in the analysis can have. The default is 25. This limit affects string variables in the variables list and also the categorical variables that are defined by the `CATEGORICAL` subcommand. A large number of categories can slow the analysis considerably. If any categorical variable violates this limit, `MVA` does not run.

### Example

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /MAXCAT=30
  /MPATTERN.
```

- The `CATEGORICAL` subcommand specifies that *region*, a numeric variable, is categorical. The variable *religion*, a string variable, is automatically categorical.
- The maximum number of categories in *region* or *religion* is 30. If either variable has more than 30 distinct values, `MVA` produces only a warning.
- Missing data patterns are shown for those cases that have at least one missing value in the specified variables.
- The summary table lists the number of missing and extreme values for each variable, including those with no missing values.

## ID Subcommand

The `ID` subcommand specifies a variable to label cases. These labels appear in the pattern tables. Without this subcommand, the case numbers are used.

### Example

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /MAXCAT=20
  /ID=country
  /MPATTERN.
```

- The values of the variable *country* are used as case labels.
- Missing data patterns are shown for those cases that have at least one missing value in the specified variables.

## NOUNIVARIATE Subcommand

By default, `MVA` computes univariate statistics for each variable—the number of cases with nonmissing values, the mean, the standard deviation, the number and percentage of missing values, and the counts of extreme low and high values. (Means, standard deviations, and extreme value counts are not reported for categorical variables.)

- To suppress the univariate statistics, specify `NOUNIVARIATE`.

**Examples**

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /CROSSTAB PERCENT=0.
```

- Univariate statistics (number of cases, means, and standard deviations) are displayed for *populatn*, *density*, *urban*, and *lifeexpf*. Also, the number of cases, counts and percentages of missing values, and counts of extreme high and low values are displayed.
- The total number of cases and counts and percentages of missing values are displayed for *region* and *religion* (a string variable).
- Separate crosstabulations are displayed for *region* and *religion*.

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region.
  /NOUNIVARIATE
  /CROSSTAB PERCENT=0.
```

- Only crosstabulations are displayed (no univariate statistics).

**TTEST Subcommand**

For each quantitative variable, a series of *t* tests are computed to test the difference of means between two groups defined by a missing indicator variable for each of the other variables. (For more information, see the topic [Missing Indicator Variables on p. 1299](#).) For example, a *t* test is performed on *populatn* between two groups defined by whether their values are present or missing for *calories*. Another *t* test is performed on *populatn* for the two groups defined by whether their values for *density* are present or missing, and the tests continue for the remainder of the variable list.

**PERCENT=n**            *Omit indicator variables with less than the specified percentage of missing values. You can specify a percentage from 0 to 100. The default is 5, indicating the omission of any variable with less than 5% missing values. If you specify 0, all rows are displayed.*

**Display of Statistics**

The following statistics can be displayed for a *t* test:

- The *t* **statistic**, for comparing the means of two groups defined by whether the indicator variable is coded as missing or nonmissing. (For more information, see the topic [Missing Indicator Variables on p. 1299](#).)

**T**                      *Display the t statistics. This setting is the default.*

**NOT**                    *Suppress the t statistics.*

- The **degrees of freedom** associated with the *t* statistic.

**DF**                     *Display the degrees of freedom. This setting is the default.*

**NODF**                  *Suppress the degrees of freedom.*

- The **probability** (two-tailed) associated with the  $t$  test, calculated for the variable that is tested without reference to other variables. Care should be taken when interpreting this probability.

**PROB**                    *Display probabilities.*

**NOPROB**                *Suppress probabilities.* This setting is the default.

- The **number of values in each group**, where groups are defined by values that are coded as missing and present in the indicator variable.

**COUNTS**                *Display counts.* This setting is the default.

**NOCOUNTS**             *Suppress counts.*

- The **means** of the groups, where groups are defined by values that are coded as missing and present in the indicator variable.

**MEANS**                 *Display means.* This setting is the default.

**NOMEANS**              *Suppress means.*

### **Example**

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /ID=country
  /TTEST.
```

- The `TTEST` subcommand produces a table of  $t$  tests. For each quantitative variable named in the variables list, a  $t$  test is performed, comparing the mean of the values for which the other variable is present against the mean of the values for which the other variable is missing.
- The table displays default statistics, including values of  $t$ , degrees of freedom, counts, and means.

## **CROSSTAB Subcommand**

`CROSSTAB` produces a table for each categorical variable, showing the frequency and percentage of values that are present (nonmissing) and the percentage of missing values for each category as related to the other variables.

- No tables are produced if there are no categorical variables.
- Each categorical variable yields a table, whether it is a string variable that is assumed to be categorical or a numeric variable that is declared on the `CATEGORICAL` subcommand.
- The categories of the categorical variable define the columns of the table.

- Each of the remaining variables defines several rows—one each for the number of values present, the percentage of values present, and the percentage of system-missing values; and one each for the percentage of values defined as each discrete type of user-missing (if they are defined).

**PERCENT=*n***      *Omit rows for variables with less than the specified percentage of missing values. You can specify a percentage from 0 to 100. The default is 5, indicating the omission of any variable with less than 5% missing values. If you specify 0, all rows are displayed.*

### Example

```
MVA VARIABLES=age income91 childs jazz folk
/CATEGORICAL=jazz folk
/CROSSTAB PERCENT=0.
```

- A table of univariate statistics is displayed by default.
- In the output are two crosstabulations (one crosstabulation for *jazz* and one crosstabulation for *folk*). The table for *jazz* displays, for each category of *jazz*, the number and percentage of present values for *age*, *income91*, *childs*, and *folk*. It also displays, for each category of *jazz*, the percentage of each type of missing value (system-missing and user-missing) in the other variables. The second crosstabulation shows similar counts and percentages for each category of *folk*.
- No rows are omitted, because PERCENT=0.

## MISMATCH Subcommand

MISMATCH produces a matrix showing percentages of cases for a pair of variables in which one variable has a missing value and the other variable has a nonmissing value (a mismatch). The diagonal elements are percentages of missing values for a single variable, while the off-diagonal elements are the percentage of mismatch of the indicator variables. [For more information, see the topic Missing Indicator Variables on p. 1299.](#) Rows and columns are sorted on missing patterns.

**PERCENT=*n***      *Omit patterns involving less than the specified percentage of cases. You can specify a percentage from 0 to 100. The default is 5, indicating the omission of any pattern that is found in less than 5% of the cases.*

**NOSORT**      *Suppress sorting of the rows and columns. The order of the variables in the variables list is used. If ALL was used in the variables list, the order is that of the data file.*

## DPATTERN Subcommand

DPATTERN lists the missing values and extreme values for each case symbolically. For a list of the symbols that are used, see [Symbols](#).

By default, the cases are listed in the order in which they appear in the file. The following keywords are available:

- SORT=varname [(order)]**      *Sort the cases according to the values of the named variables. You can specify more than one variable for sorting. Each sort variable can be in ASCENDING or DESCENDING order. The default order is ASCENDING.*
- DESCRIBE=varlist**              *List values of each specified variable for each case.*

### Example

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /ID=country
  /DPATTERN DESCRIBE=region religion SORT=region.
```

- In the data pattern table, the variables form the columns, and each case, identified by its country, defines a row.
- Missing and extreme values are indicated in the table, and, for each row, the number missing and percentage of variables that have missing values are listed.
- The values of *region* and *religion* are listed at the end of the row for each case.
- The cases are sorted by *region* in ascending order.
- Univariate statistics are displayed.

## MPATTERN Subcommand

The MPATTERN subcommand symbolically displays patterns of missing values for cases that have missing values. The variables form the columns. Each case that has any missing values in the specified variables forms a row. The rows are sorted by missing-value patterns. For use of symbols, see [Symbols](#).

- The rows are sorted to minimize the differences between missing patterns of consecutive cases.
- The columns are also sorted according to missing patterns of the variables.

The following keywords are available:

- NOSORT**                              *Suppress the sorting of variables. The order of the variables in the variables list is used. If ALL was used in the variables list, the order is that of the data file.*
- DESCRIBE=varlist**                  *List values of each specified variable for each case.*

### Example

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /ID=country
  /MPATTERN DESCRIBE=region religion.
```

- A table of missing data patterns is produced.
- The *region* and the *religion* are named for each listed case.

## TPATTERN Subcommand

The TPATTERN subcommand displays a tabulated patterns table, which lists the frequency of each missing value pattern. The variables in the variables list form the columns. Each pattern of missing values forms a row, and the frequency of the pattern is displayed.

- An *X* is used to indicate a missing value.
- The rows are sorted to minimize the differences between missing patterns of consecutive cases.
- The columns are sorted according to missing patterns of the variables.

The following keywords are available:

<b>NOSORT</b>	<i>Suppress the sorting of the columns.</i> The order of the variables in the variables list is used. If ALL was used in the variables list, the order is that of the data file.
<b>DESCRIBE=varlist</b>	<i>Display values of variables for each pattern.</i> Categories for each named categorical variable form columns in which the number of each pattern of missing values is tabulated. For quantitative variables, the mean value is listed for the cases having the pattern.
<b>PERCENT=n</b>	<i>Omit patterns that describe less than 1% of the cases.</i> You can specify a percentage from 0 to 100. The default is 1, indicating the omission of any pattern representing less than 1% of the total cases. If you specify 0, all patterns are displayed.

### Example

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /TPATTERN NOSORT DESCRIBE=populatn region.
```

- Missing value patterns are tabulated. Each row displays a missing value pattern and the number of cases having that pattern.
- DESCRIBE causes the mean value of *populatn* to be listed for each pattern. For the categories in *region*, the frequency distribution is given for the cases having the pattern in each row.

## LISTWISE Subcommand

For each quantitative variable in the variables list, the LISTWISE subcommand computes the mean, the covariance between the variables, and the correlation between the variables. The cases that are used in the computations are listwise nonmissing; that is, they have no missing value in any variable that is listed in the VARIABLES subcommand.

### Example

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /LISTWISE.
```

- Means, covariances, and correlations are displayed for *populatn*, *density*, *urban*, and *lifeexpf*. Only cases that have values for all of these variables are used.

## ***PAIRWISE Subcommand***

For each pair of quantitative variables, the `PAIRWISE` subcommand computes the number of pairwise nonmissing values, the pairwise means, the pairwise standard deviations, the pairwise covariances, and the pairwise correlation matrices. These results are organized as matrices. The cases that are used are all cases having nonmissing values for the pair of variables for which each computation is done.

### ***Example***

```
MVA VARIABLES=populatn density urban religion lifeexpf region
  /CATEGORICAL=region
  /PAIRWISE.
```

- Frequencies, means, standard deviations, covariances, and the correlations are displayed for *populatn*, *density*, *urban*, and *lifeexpf*. Each calculation uses all cases that have values for both variables under consideration.

## ***EM Subcommand***

The `EM` subcommand uses an EM (expectation-maximization) algorithm to estimate the means, the covariances, and the Pearson correlations of quantitative variables. This process is an iterative process, which uses two steps for each iteration. The E step computes expected values conditional on the observed data and the current estimates of the parameters. The M step calculates maximum-likelihood estimates of the parameters based on values that are computed in the E step.

- If no variables are listed in the `EM` subcommand, estimates are performed for all quantitative variables in the variables list.
- If you want to limit the estimation to a subset of the variables in the list, specify a subset of quantitative variables to be estimated after the subcommand name `EM`. You can also list, after the keyword `WITH`, the quantitative variables to be used in estimating.
- The output includes tables of means, correlations, and covariances.
- The estimation, by default, assumes that the data are normally distributed. However, you can specify a multivariate  $t$  distribution with a specified number of degrees of freedom or a mixed normal distribution with any mixture proportion (`PROPORTION`) and any standard deviation ratio (`LAMBDA`).
- You can save a data file with the missing values filled in. You must specify a filename and its complete path in single or double quotation marks.
- Criteria keywords and `OUTFILE` specifications must be enclosed in a single pair of parentheses.



The criteria for the `EM` subcommand are as follows:

<b>TOLERANCE=value</b>	<i>Numerical accuracy control.</i> Helps eliminate predictor variables that are highly correlated with other predictor variables and would reduce the accuracy of the matrix inversions that are involved in the calculations. The smaller the tolerance, the more inaccuracy is tolerated. The default value is 0.001.
<b>CONVERGENCE=value</b>	<i>Convergence criterion.</i> Determines when iteration ceases. If the relative change in the likelihood function is less than this value, convergence is assumed. The value of this ratio must be between 0 and 1. The default value is 0.0001.
<b>ITERATIONS=n</b>	<i>Maximum number of iterations.</i> Limits the number of iterations in the EM algorithm. Iteration stops after this many iterations even if the convergence criterion is not satisfied. The default value is 25.

Possible distribution assumptions are as follows:

<b>TDF=n</b>	<i>Student's t distribution with n degrees of freedom.</i> The degrees of freedom must be specified if you use this keyword. The degrees of freedom must be an integer that is greater than or equal to 2.
<b>LAMBDA=a</b>	<i>Ratio of standard deviations of a mixed normal distribution.</i> Any positive real number can be specified.
<b>PROPORTION=b</b>	<i>Mixture proportion of two normal distributions.</i> Any real number between 0 and 1 can specify the mixture proportion of two normal distributions.

The following keyword produces a new data file:

<b>OUTFILE='file'</b>	<i>Specify a filename or previously declared dataset name.</i> Filenames should be enclosed in quotation marks and are stored in the working directory unless a path is included as part of the file specification. Datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files. Missing values for predicted variables in the file are filled in by using the EM algorithm. (Note that the data that are completed with EM-based imputations will not in general reproduce the EM estimates from MVA.)
-----------------------	--

### Examples

```
MVA VARIABLES=males to tuition
  /EM (OUTFILE='/colleges/emdata.sav').
```

- All variables on the variables list are included in the estimations.
- The output includes the means of the listed variables, a correlation matrix, and a covariance matrix.
- A new data file named *emdata.sav* with imputed values is saved in the */colleges* directory.

```
MVA VARIABLES=all
  /EM males msport WITH males msport gradrate facratio.
```

- For *males* and *msport*, the output includes a vector of means, a correlation matrix, and a covariance matrix.

- The values in the tables are calculated by using imputed values for *males* and *msport*. Existing observations for *males*, *msport*, *gradrate*, and *facratio* are used to impute the values that are used to estimate the means, correlations, and covariances.

```
MVA VARIABLES=males to tuition
/EM verbal math WITH males msport gradrate facratio
(TDF=3 OUTFILE '/colleges/emdata.sav').
```

- The analysis uses a *t* distribution with three degrees of freedom.
- A new data file named *emdata.sav* with imputed values is saved in the */colleges* directory.

## REGRESSION Subcommand

The `REGRESSION` subcommand estimates missing values by using multiple linear regression. It can add a random component to the regression estimate. Output includes estimates of means, a covariance matrix, and a correlation matrix of the variables that are specified as predicted.

- By default, all of the variables that are specified as predictors (after `WITH`) are used in the estimation, but you can limit the number of predictors (independent variables) by using `NPREDICTORS`.
- Predicted and predictor variables, if specified, must be quantitative.
- By default, `REGRESSION` adds the observed residuals of a randomly selected complete case to the regression estimates. However, you can specify that the program add random normal, *t*, or no variates instead. The normal and *t* distributions are properly scaled, and the degrees of freedom can be specified for the *t* distribution.
- If the number of complete cases is less than half the total number of cases, the default `ADDTYPE` is `NORMAL` instead of `RESIDUAL`.
- You can save a data file with the missing values filled in. You must specify a filename and its complete path in single or double quotation marks.
- The criteria and `OUTFILE` specifications for the `REGRESSION` subcommand must be enclosed in a single pair of parentheses.

The criteria for the `REGRESSION` subcommand are as follows:

<b>TOLERANCE=value</b>	<i>Numerical accuracy control.</i> Helps eliminate predictor variables that are highly correlated with other predictor variables and would reduce the accuracy of the matrix inversions that are involved in the calculations. If a variable passes the tolerance criterion, it is eligible for inclusion. The smaller the tolerance, the more inaccuracy is tolerated. The default value is 0.001.
<b>FLIMIT=n</b>	<i>F-to-enter limit.</i> The minimum value of the <i>F</i> statistic that a variable must achieve in order to enter the regression estimation. You may want to change this limit, depending on the number of variables and the correlation structure of the data. The default value is 4.
<b>NPREDICTORS=n</b>	<i>Maximum number of predictor variables.</i> Limits the total number of predictors in the analysis. The analysis uses the stepwise selected <i>n</i> best predictors, entered in accordance with the tolerance. If <i>n</i> =0, it is equivalent to replacing each variable with its mean.

**ADDDTYPE**

*Type of distribution from which the error term is randomly drawn.* Random errors can be added to the regression estimates before the means, correlations, and covariances are calculated. You can specify one of the following types:

**RESIDUAL.** Error terms are chosen randomly from the observed residuals of complete cases to be added to the regression estimates.

**NORMAL.** Error terms are randomly drawn from a distribution with the expected value 0 and the standard deviation equal to the square root of the mean squared error term (sometimes called the **root mean squared error**, or RMSE) of the regression.

**T(n).** Error terms are randomly drawn from the t(n) distribution and scaled by the RMSE. The degrees of freedom can be specified in parentheses. If T is specified without a value, the default degrees of freedom is 5.

**NONE.** Estimates are made from the regression model with no error term added.

The following keyword produces a new data file:

**OUTFILE**

*Specify a filename or previously declared dataset name.* Filenames should be enclosed in quotation marks and are stored in the working directory unless a path is included as part of the file specification. Datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files. Missing values for the dependent variables in the file are imputed (filled in) by using the regression algorithm.

**Examples**

```
MVA VARIABLES=males to tuition
  /REGRESSION (OUTFILE=' /colleges/regdata.sav').
```

- All variables in the variables list are included in the estimations.
- The output includes the means of the listed variables, a correlation matrix, and a covariance matrix.
- A new data file named *regdata.sav* with imputed values is saved in the */colleges* directory.

```
MVA VARIABLES=males to tuition
  /REGRESSION males verbal math WITH males verbal math faculty
  (ADDDTYPE = T(7)).
```

- The output includes the means of the listed variables, a correlation matrix, and a covariance matrix.
- A *t* distribution with 7 degrees of freedom is used to produce the randomly assigned additions to the estimates.

# N OF CASES

```
N OF CASES n
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
N OF CASES 100.
```

## **Overview**

N OF CASES (alias N) limits the number of cases in the active dataset to the first *n* cases.

### **Basic Specification**

The basic specification is N OF CASES followed by at least one space and a positive integer. Cases in the active dataset are limited to the specified number.

### **Syntax Rules**

- To limit the number of cases for the next procedure only, use the TEMPORARY command before N OF CASES (see TEMPORARY).
- In some versions of the program, N OF CASES can be specified only after a active dataset is defined.

### **Operations**

- N OF CASES takes effect as soon as it is encountered in the command sequence. Thus, special attention should be paid to the position of N OF CASES among commands. [For more information, see the topic Command Order on p. 41.](#)
- N OF CASES limits the number of cases that are analyzed by all subsequent procedures in the session. The active dataset will have no more than *n* cases after the first data pass following the N OF CASES command. Any subsequent N OF CASES command that specifies a greater number of cases will be ignored.
- If N OF CASES specifies more cases than can actually be built, the program builds as many cases as possible.
- If N OF CASES is used with SAMPLE or SELECT IF, the program reads as many records as required to build the specified *n* cases. It makes no difference whether N OF CASES precedes or follows SAMPLE or SELECT IF.

## **Example**

```
GET FILE='/data/city.sav'.  
N 100.
```

- N OF CASES limits the number of cases on the active dataset to the first 100 cases. Cases are limited for all subsequent analyses.

### Example

```
DATA LIST FILE='/data/prsnnl.txt'  
  / NAME 1-20 (A) AGE 22-23 SALARY 25-30.  
N 25.  
SELECT IF (SALARY GT 20000).  
LIST.
```

- DATA LIST defines variables from file *prsnnl.txt*.
- N OF CASES limits the active dataset to 25 cases after cases have been selected by SELECT IF.
- SELECT IF selects only cases in which *SALARY* is greater than \$20,000.
- LIST produces a listing of the cases in the active dataset. If the original active dataset has fewer than 25 cases in which *SALARY* is greater than 20,000, fewer than 25 cases will be listed.

### Example

```
DATA LIST FILE='/data/prsnnl.txt'  
  / NAME 1-20 (A) AGE 22-23 SALARY 25-30 DEPT 32.  
LIST.  
TEMPORARY.  
N 25.  
FREQUENCIES VAR=SALARY.  
N 50.  
FREQUENCIES VAR=AGE.  
REPORT FORMAT=AUTO /VARS=NAME AGE SALARY /BREAK=DEPT  
  /SUMMARY=MEAN.
```

- The first N OF CASES command is temporary. Only 25 cases are used in the first FREQUENCIES procedure.
- The second N OF CASES command is permanent. The second frequency table and the report are based on 50 cases from file *prsnnl.txt*. The active dataset now contains 50 cases (assuming that the original active dataset had at least that many cases).

# NAIVEBAYES

NAIVEBAYES is available in IBM® SPSS® Statistics Server.

NAIVEBAYES dependent variable BY factor list WITH covariate list

```
[/EXCEPT VARIABLES=varlist]
[/FORCE [FACTORS=varlist] [COVARIATES=varlist]]
[/TRAININGSAMPLE {PERCENT=number }
                 {VARIABLE=varname}
[/SUBSET {MAXSIZE={AUTO** } [(BESTSUBSET={PSEUDOBIC } )]}]
          {integer}           {TESTDATA }
          {EXACTSIZE=integer }
          {NOSELECTION        }
[/CRITERIA [BINS={10** }
            {integer}
            [MEMALLOCATE {1024** }
                          {number }
            [TIMER={5** }
                   {number}
[/MISSING USERMISSING={EXCLUDE**}]
           {INCLUDE }
[/PRINT [CPS**] [EXCLUDED**] [SUMMARY**]
        [SELECTED**] [CLASSIFICATION**]
        [NONE]]
[/SAVE [PREDVAL[(varname)]]
       [PREDPROB[(rootname[:{25 }])]]]
       {integer}
[/OUTFILE MODEL=file]
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- Command introduced.

## **Example**

NAIVEBAYES default.

## Overview

The NAIVEBAYES procedure can be used in three ways:

1. **Predictor selection followed by model building.** The procedure submits a set of predictor variables and selects a smaller subset. Based on the Naïve Bayes model for the selected predictors, the procedure then classifies cases.
2. **Predictor selection only.** The procedure selects a subset of predictors for use in subsequent predictive modeling procedures but does not report classification results.
3. **Model building only.** The procedure fits the Naïve Bayes classification model by using all input predictors.

NAIVEBAYES is available for categorical dependent variables only and is not intended for use with a very large number of predictors.

### Options

**Methods.** The NAIVEBAYES procedure performs predictor selection followed by model building, or the procedure performs predictor selection only, or the procedure performs model building only.

**Training and test data.** NAIVEBAYES optionally divides the dataset into training and test samples. Predictor selection uses the training data to compute the underlying model, and either the training or the test data can be used to determine the “best” subset of predictors. If the dataset is partitioned, classification results are given for both the training and test samples. Otherwise, results are given for the training sample only.

**Binning.** The procedure automatically distributes scale predictors into 10 bins, but the number of bins can be changed.

**Memory allocation.** The NAIVEBAYES procedure automatically allocates 128MB of memory for storing training records when computing average log-likelihoods. The amount of memory that is allocated for this task can be modified.

**Timer.** The procedure automatically limits processing time to 5 minutes, but a different time limit can be specified.

**Maximum or exact subset size.** Either a maximum or an exact size can be specified for the subset of selected predictors. If a maximum size is used, the procedure creates a sequence of subsets, from an initial (smaller) subset to the maximum-size subset. The procedure then selects the “best” subset from this sequence.

**Missing values.** Cases with missing values for the dependent variable or for all predictors are excluded. The NAIVEBAYES procedure has an option for treating user-missing values of categorical variables as valid. User-missing values of scale variables are always treated as invalid.

**Output.** NAIVEBAYES displays pivot table output by default but offers an option for suppressing most such output. The procedure displays the lists of selected categorical and scale predictors in a text block. These lists can be copied for use in subsequent modeling procedures. The NAIVEBAYES procedure also optionally saves predicted values and probabilities based on the Naïve Bayes model.

**Basic Specification**

The basic specification is the `NAIVEBAYES` command followed by a dependent variable.

By default, `NAIVEBAYES` treats all variables — except the dependent variable and the weight variable if it is defined — as predictors, with the dictionary setting of each predictor determining its measurement level. `NAIVEBAYES` selects the “best” subset of predictors (based on the Naïve Bayes model) and then classifies cases by using the selected predictors. User-missing values are excluded and pivot table output is displayed by default.

**Syntax Rules**

- All subcommands are optional.
- Subcommands may be specified in any order.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equal signs, and slashes that are shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not honored.

**Operations**

The `NAIVEBAYES` procedure automatically excludes cases and predictors with any of the following properties:

- Cases with a missing value for the dependent variable.
- Cases with missing values for all predictors.
- Predictors with missing values for all cases.
- Predictors with the same value for all cases.

The `NAIVEBAYES` procedure requires predictors to be categorical. Any scale predictors that are input to the procedure are temporarily binned into categorical variables for the procedure.

If predictor selection is used, the `NAIVEBAYES` procedure selects a subset of predictors that “best” predict the dependent variable, based on the training data. The procedure first creates a sequence of subsets, with an increasing number of predictors in each subset. The predictor that is added to each subsequent subset is the predictor that increases the average log-likelihood the most. The procedure uses simulated data to compute the average log-likelihood when the training dataset cannot fit into memory.

The final subset is obtained by using one of two approaches:

- By default, a maximum subset size is used. This approach creates a sequence of subsets from the initial subset to the maximum-size subset. The “best” subset is chosen by using a BIC-like criterion or a test data criterion.
- A particular subset size may be used to select the subset with the specified size.



If model building is requested, the NAIVEBAYES procedure classifies cases based on the Naïve Bayes model for the input or selected predictors, depending on whether predictor selection is requested. For a given case, the classification—or predicted category—is the dependent variable category with the highest posterior probability.

The NAIVEBAYES procedure uses the IBM® SPSS® Statistics random number generator in the following two scenarios: (1) if a percentage of cases in the active dataset is randomly assigned to the test dataset, and (2) if the procedure creates simulated data to compute the average log-likelihood when the training records cannot fit into memory. To ensure that the same results are obtained regardless of which scenario is in effect when NAIVEBAYES is invoked repeatedly, specify a seed on the SET command. If a seed is not specified, a default random seed is used, and results may differ across runs of the NAIVEBAYES procedure.

### **Frequency Weight**

If a WEIGHT variable is in effect, its values are used as frequency weights by the NAIVEBAYES procedure.

- Cases with missing weights or weights that are less than 0.5 are not used in the analyses.
- The weight values are rounded to the nearest whole numbers before use. For example, 0.5 is rounded to 1, and 2.4 is rounded to 2.

### **Limitations**

SPLIT FILE settings are ignored by the NAIVEBAYES procedure.

## **Examples**

### **Predictor selection followed by model building**

```
NAIVEBAYES default
  /EXCEPT VARIABLES=preddef1 preddef2 preddef3 training
  /TRAININGSAMPLE VARIABLE=training
  /SAVE PREDVAL PREDPROB.
```

- This analysis specifies *default* as the response variable.
- All other variables are to be considered as possible predictors, with the exception of *preddef1*, *preddef2*, *preddef3*, and *training*.
- Cases with a value of 1 on the variable *training* are assigned to the training sample and used to create the series of predictor subsets, while all other cases are assigned to the test sample and used to select the “best” subset.
- Model-predicted values of *default* are saved to the variable *PredictedValue*.
- Model-estimated probabilities for the values of *default* are saved to the variables *PredictedProbability\_1* and *PredictedProbability\_2*.

### **Predictor selection only**

```
NAIVEBAYES default
```

```

/EXCEPT VARIABLES=preddef1 preddef2 preddef3 validate
/TRAININGSAMPLE VARIABLE=validate
/SUBSET EXACTSIZE=5
/PRINT CLASSIFICATION=NO.

```

- The NAIVEBAYES procedure treats *default* as the dependent variable and selects a subset of five predictors from all other variables, with the exception of *preddef1*, *preddef2*, *preddef3*, and *validate*.

### **Model building only**

```

NAIVEBAYES response_01
  BY addresscat callcard callid callwait card card2 churn
    commutecarpool confer ebill edcat equip forward internet
    multline owngame ownipod ownpc spousedcat tollfree voice
  WITH cardmon ed equipmon equipten lncardmon lntollmon
    pets_saltnfish spoused tollmon tollten
/SUBSET NOSELECTION
/SAVE PREDPROB.

```

- This analysis specifies *response\_01* as the response variable.
- Variables following the BY keyword are treated as categorical predictors, while those following the WITH keyword are treated as scale.
- The SUBSET subcommand specifies that the procedure should not perform predictor selection. All specified predictors are to be used in creating the classification.
- Model-estimated probabilities for the values of *response\_01* are saved to the variables *PredictedProbability\_1* and *PredictedProbability\_2*.

## **Variable Lists**

The variable lists specify the dependent variable, any categorical predictors (also known as factors), and any scale predictors (also known as covariates).

- The dependent variable must be the first specification on the NAIVEBAYES command.
- The dependent variable may not be the weight variable.
- The dependent variable is the only required specification on the NAIVEBAYES command.
- The dependent variable must have a dictionary setting of ordinal or nominal. In either case, NAIVEBAYES treats the dependent variable as categorical.
- The names of the factors, if any, must be preceded by the keyword BY.
- If keyword BY is specified with no factors, a warning is issued and the keyword is ignored.
- The names of covariates must be preceded by the keyword WITH.
- If keyword WITH is specified with no covariates, a warning is issued and the keyword is ignored.
- If the dependent variable or the weight variable is specified within a factor list or a covariate list, the variable is ignored in the list.
- All variables that are specified within a factor or covariate list must be unique. If duplicate variables are specified within a list, the duplicates are ignored.
- If duplicate variables are specified across the factor and covariate lists, an error is issued.

- The universal keywords `TO` and `ALL` may be specified in the factor and covariate lists.
- If the `BY` and `WITH` keywords are not specified, all variables in the active dataset—except the dependent variable, the weight variable, and any variables that are specified on the `EXCEPT` subcommand—are treated as predictors. If the dictionary setting of a predictor is nominal or ordinal, the predictor is treated as a factor. If the dictionary setting is scale, the predictor is treated as a covariate. (Note that any variables on the `FORCE` subcommand are still forced into each subset of selected predictors.)
- The dependent variable and factor variables can be numeric or string.
- The covariates must be numeric.

## ***EXCEPT Subcommand***

The `EXCEPT` subcommand lists any variables that the `NAIVEBAYES` procedure should exclude from the factor or covariate lists on the command line. This subcommand is useful if the factor or covariate lists contain a large number of variables—specified by using the `TO` or `ALL` keyword, for example—but a few variables (e.g., Case ID or a weight variable) should be excluded.

- The `EXCEPT` subcommand ignores the following types of variables if they are specified: Duplicate variables; the dependent variable; variables that are not specified on the command line's factor or covariate lists; and variables that are specified on the `FORCE` subcommand.
- There is no default variable list on the `EXCEPT` subcommand.

## ***FORCE Subcommand***

The `FORCE` subcommand specifies any predictors that will be in the initial predictor subset and all subsequent predictor subsets. The specified predictors are considered important and will be in the final subset irrespective of any other chosen predictors.

- Variables that are specified on the `FORCE` subcommand do not need to be specified in the variable lists on the command line.
- The `FORCE` subcommand overrides variable lists on the command line and overrides the `EXCEPT` subcommand. If a variable specified on the `FORCE` subcommand is also specified on the command line or the `EXCEPT` subcommand, the variable is forced into all subsets.
- There is no default list of forced variables; the default initial subset is the empty set.

### ***FACTORS Keyword***

The `FACTORS` keyword specifies any factors that should be forced into each subset.

- If duplicate variables are specified, the duplicates are ignored.
- The specified variables may not include the dependent variable, the weight variable, or any variable that is specified on the `COVARIATES` keyword.
- Specified variables may be numeric or string.

***COVARIATES Keyword***

The `COVARIATES` keyword specifies any covariates that should be forced into each subset.

- If duplicate variables are specified, the duplicates are ignored.
- The specified variables may not include the dependent variable, the weight variable, or any variable that is specified on the `FACTORS` keyword.
- Specified variables must be numeric.

***TRAININGSAMPLE Subcommand***

The `TRAININGSAMPLE` subcommand indicates the method of partitioning the active dataset into training and test samples. You can specify either a percentage of cases to assign to the training sample, or you can specify a variable that indicates whether a case is assigned to the training sample.

- If `TRAININGSAMPLE` is not specified, all cases in the active dataset are treated as training data records.

***PERCENT Keyword***

The `PERCENT` keyword specifies the percentage of cases in the active dataset to randomly assign to the training sample. All other cases are assigned to the test sample. The percentage must be a number that is greater than 0 and less than 100. There is no default percentage.

If a weight variable is defined, the `PERCENT` keyword may not be used.

***VARIABLE Keyword***

The `VARIABLE` keyword specifies a variable that indicates which cases in the active dataset are assigned to the training sample. Cases with a value of 1 on the variable are assigned to the training sample. All other cases are assigned to the test sample.

- The specified variable may not be the dependent variable, the weight variable, any variable that is specified in the factor or covariate lists of the command line, or any variable that is specified in the factor or covariate lists of the `FORCE` subcommand.
- The variable must be numeric.

***SUBSET Subcommand***

The `SUBSET` subcommand gives settings for the subset of selected predictors.

- There are three mutually exclusive settings: (1) specify a maximum subset size and a method of selecting the best subset, (2) specify an exact subset size, or (3) do not specify a selection.
- Only one of the keywords `MAXSIZE`, `EXACTSIZE`, or `NOSELECTION` may be specified. The `BESTSUBSET` option is available only if `MAXSIZE` is specified.

**MAXSIZE Keyword**

The `MAXSIZE` keyword specifies the maximum subset size to use when creating the sequence of predictor subsets. The `MAXSIZE` value is the size of the largest subset beyond any predictors that were forced via the `FORCE` subcommand. If no predictors are forced, the `MAXSIZE` value is simply the size of the largest subset.

- Value `AUTO` indicates that the number should be computed automatically. Alternatively, a positive integer may be specified. The integer must be less than or equal to the number of unique predictors on the `NAIVEBAYES` command.
- By default, `MAXSIZE` is used and `AUTO` is the default value.

**BESTSUBSET Keyword**

The `BESTSUBSET` keyword indicates the criterion for finding the best subset when a maximum subset size is used.

- This keyword is honored only if the `MAXSIZE` keyword is in effect and must be given in parentheses immediately following the `MAXSIZE` specification.

**PSEUDOBI** *Use the pseudo-BIC criterion.* The pseudo-BIC criterion is based on the training sample. If the active dataset is not partitioned into training and test samples, `PSEUDOBI` is the default. If the active dataset is partitioned, `PSEUDOBI` is available but is not the default.

**TESTDATA** *Use the test data criterion.* The test data criterion is based on the test sample. If the active dataset is partitioned into training and test samples, `TESTDATA` is the default. If the active dataset is not partitioned, `TESTDATA` may not be specified.

**EXACTSIZE Keyword**

The `EXACTSIZE` keyword specifies a particular subset size to use. The `EXACTSIZE` value is the size of the subset beyond any predictors forced via the `FORCE` subcommand. If no predictors are forced, then the `EXACTSIZE` value is simply the size of the subset.

- A positive integer may be specified. The integer must be less than the number of unique predictors on the `NAIVEBAYES` command.
- There is no default value.

**NOSELECTION Keyword**

The `NOSELECTION` keyword indicates that all predictors that are specified on the `NAIVEBAYES` command—excluding any predictors that are also specified on the `EXCEPT` subcommand—are included in the final subset. This specification is useful if the `NAIVEBAYES` procedure is used for model building but not predictor selection.

**CRITERIA Subcommand**

The `CRITERIA` subcommand specifies computational and resource settings for the `NAIVEBAYES` procedure.

**BINS Keyword**

The `BINS` keyword specifies the number of bins to use when dividing the domain of a scale predictor into equal-width bins. A positive integer greater than 1 may be specified. The default is 10.

**MEMALLOCATE Keyword**

The `MEMALLOCATE` keyword specifies the maximum amount of memory in megabytes (MB) that the `NAIVEBAYES` procedure uses to store training data records when computing the average log-likelihood. If the amount of memory that is required to store records is larger, simulated data are used instead.

- Any number that is greater than or equal to 4 may be specified. Consult your system administrator for the largest value that can be specified on your system. The default is 1024.

**TIMER Keyword**

The `TIMER` keyword specifies the maximum number of minutes during which the `NAIVEBAYES` procedure can run. If the time limit is exceeded, the procedure is terminated and no results are given. Any number that is greater than or equal to 0 may be specified. Specifying 0 turns the timer off completely. The default is 5.

**MISSING Subcommand**

The `MISSING` subcommand controls whether user-missing values for categorical variables are treated as valid values. By default, user-missing values for categorical variables are treated as invalid.

- User-missing values for scale variables are always treated as invalid.
- System-missing values for any variables are always treated as invalid.

**USERMISSING=EXCLUDE**     *User-missing values for categorical variables are treated as invalid values. This setting is the default.*

**USERMISSING=INCLUDE**     *User-missing values for categorical variables are treated as valid values.*

**PRINT Subcommand**

The `PRINT` subcommand indicates the statistical output to display.

**CPS**     *Case processing summary.* The table summarizes the number of cases that are included and excluded in the analysis. This table is shown by default.

**EXCLUDED**     *Predictors excluded due to missing or constant values for all cases.* The table lists excluded predictors by type (categorical or scale) and the reasons for being excluded.

<b>SUMMARY</b>	<i>Statistical summary of the sequence of predictor subsets.</i> This table is shown by default. The <code>SUMMARY</code> keyword is ignored if <code>NOSELECTION</code> is specified on the <code>SUBSET</code> subcommand.
<b>SELECTED</b>	<i>Selected predictors by type (categorical or scale).</i> This table is shown by default. The <code>SELECTED</code> keyword is ignored if <code>NOSELECTION</code> is specified on the <code>SUBSET</code> subcommand.
<b>CLASSIFICATION</b>	<i>Classification table.</i> The table gives the number of cases that are classified correctly and incorrectly for each dependent variable category. If test data are defined, classification results are given for the training and the test samples. If test data are not defined, classification results are given only for the training sample. This table is shown by default.
<b>NONE</b>	<i>Suppress all displayed output except the Notes table and any warnings.</i> This keyword may not be specified with any other keywords.

## ***SAVE Subcommand***

The `SAVE` subcommand writes optional temporary variables to the active dataset.

<b>PREDVAL(varname)</b>	<i>Predicted value.</i> The predicted value is the dependent variable category with the highest posterior probability as estimated by the Naïve Bayes model. A valid variable name must be specified. The default variable name is <i>PredictedValue</i> .
<b>PREDPROB(rootname:n)</b>	<i>Predicted probability.</i> The predicted probabilities of the first <i>n</i> categories of the dependent variable are saved. Suffixes are added to the root name to get a group of variable names that correspond to the dependent variable categories. If a root name is specified, it must be a valid variable name. The root name can be followed by a colon and a positive integer that indicates the number of probabilities to save. The default root name is <i>PredictedProbability</i> . The default <i>n</i> is 25. To specify <i>n</i> without a root name, enter a colon before the number.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand writes the Naïve Bayes model to an XML file. The Naïve Bayes model is based on the training sample even if the active dataset is partitioned into training and test samples.

A valid file name must be specified on the `MODEL` keyword.

# ***NEW FILE***

NEW FILE

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Overview***

The NEW FILE command clears the active dataset. NEW FILE is used when you want to build a new active dataset by generating data within an input program (see INPUT PROGRAM–END INPUT PROGRAM).

### ***Basic Specification***

NEW FILE is always specified by itself. No other keyword is allowed.

### ***Operations***

- NEW FILE creates a new, blank active dataset. The command takes effect as soon as it is encountered.
- When you build an active dataset with GET, DATA LIST, or other file-definition commands (such as ADD FILES or MATCH FILES), the active dataset is automatically replaced. It is not necessary to specify NEW FILE.



# NLR

NLR and CNLR are available in the Regression option.

```
MODEL PROGRAM parameter=value [parameter=value ...]
transformation commands

[DERIVATIVES
transformation commands]

[CLEAR MODEL PROGRAMS]
```

## *Procedure CNLR (Constrained Nonlinear Regression):*

```
[CONSTRAINED FUNCTIONS
transformation commands]

CNLR dependent var

[/FILE=file]    [/OUTFILE=file]

[/PRED=varname]

[/SAVE [PRED] [RESID[(varname)]] [DERIVATIVES] [LOSS]]

[/CRITERIA=[ITER n] [MITER n] [CKDER {0.5**}]
           {n }
           [ISTEP {1E+20**}] [FPR n] [LFTOL n]
           {n }
           [LSTOL n] [STEPLIMIT {2**}] [NFTOL n]
           {n }
           [FTOL n] [OPTOL n] [CRSHTOL {.01**}]]
           {n }

[/BOUNDS=expression, expression, ...]

[/LOSS=varname]

[/BOOTSTRAP [=n]]
```

## *Procedure NLR (Nonlinear Regression):*

```
NLR dependent var

[/FILE=file]    [/OUTFILE=file]

[/PRED=varname]

[/SAVE [PRED] [RESID [(varname)]] [DERIVATIVES]]

[/CRITERIA=[ITER {100**}] [CKDER {0.5**}]
           {n }           {n }
           [SSCON {1E-8**}] [PCON {1E-8**}] [RCON {1E-8**}]]
           {n }           {n }           {n }
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

**Example**

```
MODEL PROGRAM A=.6.  
COMPUTE PRED=EXP(A*X).  
  
NLR Y.
```

**Overview**

Nonlinear regression is used to estimate parameter values and regression statistics for models that are not linear in their parameters. There are two procedures for estimating nonlinear equations. CNLR (constrained nonlinear regression), which uses a sequential quadratic programming algorithm, is applicable for both constrained and unconstrained problems. NLR (nonlinear regression), which uses a Levenberg-Marquardt algorithm, is applicable only for unconstrained problems.

CNLR is more general. It allows linear and nonlinear constraints on any combination of parameters. It will estimate parameters by minimizing any smooth loss function (objective function) and can optionally compute bootstrap estimates of parameter standard errors and correlations. The individual bootstrap parameter estimates can optionally be saved in a separate IBM® SPSS® Statistics data file.

Both programs estimate the values of the parameters for the model and, optionally, compute and save predicted values, residuals, and derivatives. Final parameter estimates can be saved in SPSS Statistics data files and used in subsequent analyses.

CNLR and NLR use much of the same syntax. Some of the following sections discuss features that are common to both procedures. In these sections, the notation [C]NLR means that either the CNLR or NLR procedure can be specified. Sections that apply only to CNLR or only to NLR are clearly identified.

**Options**

**The Model.** You can use any number of transformation commands under `MODEL PROGRAM` to define complex models.

**Derivatives.** You can use any number of transformation commands under `DERIVATIVES` to supply derivatives.

**Adding Variables to Active Dataset.** You can add predicted values, residuals, and derivatives to the active dataset with the `SAVE` subcommand.

**Writing Parameter Estimates to a New Data File.** You can save final parameter estimates as an external SPSS Statistics data file by using the `OUTFILE` subcommand; you can retrieve them in subsequent analyses by using the `FILE` subcommand.

**Controlling Model-Building Criteria.** You can control the iteration process that is used in the regression with the `CRITERIA` subcommand.

**Additional CNLR Controls.** For CNLR, you can impose linear and nonlinear constraints on the parameters with the `BOUNDS` subcommand. Using the `LOSS` subcommand, you can specify a loss function for CNLR to minimize and, using the `BOOTSTRAP` subcommand, you can provide bootstrap estimates of the parameter standard errors, confidence intervals, and correlations.

### **Basic Specification**

The basic specification requires three commands: `MODEL PROGRAM`, `COMPUTE` (or any other computational transformation command), and `[C]NLR`.

- The `MODEL PROGRAM` command assigns initial values to the parameters and signifies the beginning of the model program.
- The computational transformation command generates a new variable to define the model. The variable can take any legitimate name, but if the name is not `PRED`, the `PRED` subcommand will be required.
- The `[C]NLR` command provides the regression specifications. The minimum specification is the dependent variable.
- By default, the residual sum of squares and estimated values of the model parameters are displayed for each iteration. Statistics that are generated include regression and residual sums of squares and mean squares, corrected and uncorrected total sums of squares,  $R^2$ , parameter estimates with their asymptotic standard errors and 95% confidence intervals, and an asymptotic correlation matrix of the parameter estimates.

### **Command Order**

- The model program, beginning with the `MODEL PROGRAM` command, must precede the `[C]NLR` command.
- The derivatives program (when used), beginning with the `DERIVATIVES` command, must follow the model program but precede the `[C]NLR` command.
- The constrained functions program (when used), beginning with the `CONSTRAINED FUNCTIONS` command, must immediately precede the `CNLR` command. The constrained functions program cannot be used with the `NLR` command.
- The `CNLR` command must follow the block of transformations for the model program and the derivatives program when specified; the `CNLR` command must also follow the constrained functions program when specified.
- Subcommands on `[C]NLR` can be named in any order.

### **Syntax Rules**

- The `FILE`, `OUTFILE`, `PRED`, and `SAVE` subcommands work the same way for both `CNLR` and `NLR`.
- The `CRITERIA` subcommand is used by both `CNLR` and `NLR`, but iteration criteria are different. Therefore, the `CRITERIA` subcommand is documented separately for `CNLR` and `NLR`.
- The `BOUNDS`, `LOSS`, and `BOOTSTRAP` subcommands can be used only with `CNLR`. They cannot be used with `NLR`.

## **Operations**

- By default, the predicted values, residuals, and derivatives are created as temporary variables. To save these variables, use the `SAVE` subcommand.

## Weighting Cases

- If case weighting is in effect, [C]NLR uses case weights when calculating the residual sum of squares and derivatives. However, the degrees of freedom in the ANOVA table are always based on unweighted cases.
- When the model program is first invoked for each case, the weight variable's value is set equal to its value in the active dataset. The model program may recalculate that value. For example, to effect a robust estimation, the model program may recalculate the weight variable's value as an inverse function of the residual magnitude. [C]NLR uses the weight variable's value after the model program is executed.

## Missing Values

Cases with missing values for any of the dependent or independent variables that are named on the [C]NLR command are excluded.

- Predicted values, but not residuals, can be calculated for cases with missing values on the dependent variable.
- [C]NLR ignores cases that have missing, negative, or zero weights. The procedure displays a warning message if it encounters any negative or zero weights at any time during its execution.
- If a variable that is used in the model program or the derivatives program is omitted from the independent variable list on the [C]NLR command, the predicted value and some or all of the derivatives may be missing for every case. If this situation happens, an error message is generated.

## Examples

```
MODEL PROGRAM A=.5 B=1.6.
COMPUTE PRED=A*SPEED**B.

DERIVATIVES.
COMPUTE D.A=SPEED**B.
COMPUTE D.B=A*LN(SPEED)*SPEED**B.

NLR STOP.
```

- MODEL PROGRAM assigns values to the model parameters *A* and *B*.
- COMPUTE generates the variable *PRED* to define the nonlinear model using parameters *A* and *B* and the variable *SPEED* from the active dataset. Because this variable is named *PRED*, the PRED subcommand is not required on NLR.
- DERIVATIVES indicates that calculations for derivatives are being supplied.
- The two COMPUTE statements on the DERIVATIVES transformations list calculate the derivatives for the parameters *A* and *B*. If either parameter had been omitted, NLR would have calculated it numerically.
- NLR specifies *STOP* as the dependent variable. It is not necessary to specify *SPEED* as the independent variable because it has been used in the model and derivatives programs.

### **Asymptotic Regression**

```
* NonLinear Regression.
MODEL PROGRAM b1=13 b2=-6 b3=-1.33 .
COMPUTE PRED_ = b1 + b2 * exp(b3*advert).
CNLR sales
  /OUTFILE='C:\TEMP\SPSSFNLR.TMP'
  /PRED PRED_
  /BOUNDS b1 >= 0; b2 <= 0; b3 <= 0
  /SAVE PRED RESID
  /CRITERIA STEPLIMIT 2 ISTEP 1E+20 .
```

- MODEL PROGRAM assigns initial values to the model parameters *b1*, *b2*, and *b3*.
- COMPUTE generates the variable *PRED\_* to define the nonlinear model using the parameters from MODEL PROGRAM and the variable *advert* from the active dataset.
- CNLR specifies *sales* as the dependent variable. It is not necessary to specify *advert* as the independent variable because it has been used in the computation of *PRED\_*.
- The OUTFILE subcommand saves the final parameter estimates to an external file.
- The PRED subcommand specifies that *PRED\_* is the nonlinear model.
- The BOUNDS subcommand specifies that *b1* must be nonnegative and *b2* and *b3* must be nonpositive.
- The SAVE subcommand requests the predicted values and residuals be saved as new variables to the active dataset.
- All other options are set to their default values.

### **Using Conditional Logic to Fit a Segmented Model**

```
* NonLinear Regression.
MODEL PROGRAM b1=.65 b2=5 b3=0.03 a1=0.0 a2=.5 a3=-0.025.
COMPUTE PRED_ = (time<20)*b1/(1 + b2*exp(-b3*time)) + (time>=20)*(a1 + a2*exp(a3*(time-19))
CNLR infected
  /OUTFILE='C:\TEMP\SPSSFNLR.TMP'
  /PRED PRED_
  /BOUNDS b1 >= 0; b2 >= 0; b3 >= 0; a1 >= 0; a2 >= 0; a3 <= 0
  /CRITERIA STEPLIMIT 2 ISTEP 1E+20
  /SAVE PRED RESID.
```

- MODEL PROGRAM assigns initial values to the model parameters *b1*, *b2*, *b3*, *a1*, *a2*, and *a3*.
- COMPUTE generates the variable *PRED\_* to define the nonlinear model using the parameters from MODEL PROGRAM and the variable *time* from the active dataset.
- CNLR specifies *infected* as the dependent variable. It is not necessary to specify *time* as the independent variable because it has been used in the computation of *PRED\_*.
- The OUTFILE subcommand saves the final parameter estimates to an external file.
- The PRED subcommand specifies that *PRED\_* is the nonlinear model.
- The BOUNDS subcommand specifies that *b1*, *b2*, *b3*, *a1*, and *a2* must be nonnegative and *a3* must be nonpositive.
- The SAVE subcommand requests the predicted values and residuals be saved as new variables to the active dataset.
- All other options are set to their default values.

## **MODEL PROGRAM Command**

The `MODEL PROGRAM` command assigns initial values to the parameters and signifies the beginning of the model program. The model program specifies the nonlinear equation that is chosen to model the data. There is no default model.

- The model program is required and must precede the `[C]NLR` command.
- The `MODEL PROGRAM` command must specify all parameters in the model program. Each parameter must be individually named. Keyword `TO` is not allowed.
- Parameters can be assigned any acceptable variable name. However, if you intend to write the final parameter estimates to a file with the `OUTFILE` subcommand, do not use the name `SSE` or `NCASES` (see [OUTFILE Subcommand on p. 1331](#)).
- Each parameter in the model program must have an assigned value. The value can be specified on `MODEL PROGRAM` or read from an existing parameter data file named on the `FILE` subcommand.
- Zero should be avoided as an initial value because it provides no information about the scale of the parameters. This situation is especially true for `CNLR`.
- The model program must include at least one command that uses the parameters and the independent variables (or preceding transformations of these) to calculate the predicted value of the dependent variable. This predicted value defines the nonlinear model. There is no default model.
- By default, the program assumes that `PRED` is the name assigned to the variable for the predicted values. If you use a different variable name in the model program, you must supply the name on the `PRED` subcommand (see [PRED Subcommand on p. 1332](#)).
- In the model program, you can assign a label to the variable holding predicted values and also change its print and write formats, but you should not specify missing values for this variable.
- You can use any computational commands (such as `COMPUTE`, `IF`, `DO IF`, `LOOP`, `END LOOP`, `END IF`, `RECODE`, or `COUNT`) or output commands (`WRITE`, `PRINT`, or `XSAVE`) in the model program, but you cannot use input commands (such as `DATA LIST`, `GET`, `MATCH FILES`, or `ADD FILES`).
- Transformations in the model program are used only by `[C]NLR`, and they do not affect the active dataset. The parameters that are created by the model program do not become a part of the active dataset. Permanent transformations should be specified before the model program.

### **Caution: Initial Values**

The selection of good initial values for the parameters in the model program is very important to the operation of `[C]NLR`. The selection of poor initial values can result in no solution, a local solution rather than a general solution, or a physically impossible solution.

#### **Example**

```
MODEL PROGRAM A=10 B=1 C=5 D=1 .  
COMPUTE PRED= A*exp(B*X) + C*exp(D*X) .
```

- The `MODEL PROGRAM` command assigns starting values to the four parameters *A*, *B*, *C*, and *D*.
- `COMPUTE` defines the model to be fit as the sum of two exponentials.

## ***DERIVATIVES Command***

The optional `DERIVATIVES` command signifies the beginning of the derivatives program. The derivatives program contains transformation statements for computing some or all of the derivatives of the model. The derivatives program must follow the model program but precede the `[C]NLR` command.

If the derivatives program is not used, `[C]NLR` numerically estimates derivatives for all the parameters. Providing derivatives reduces computation time and, in some situations, may result in a better solution.

- The `DERIVATIVES` command has no further specifications but must be followed by the set of transformation statements that calculate the derivatives.
- You can use any computational commands (such as `COMPUTE`, `IF`, `DO IF`, `LOOP`, `END LOOP`, `END IF`, `RECODE`, or `COUNT`) or output commands (`WRITE`, `PRINT`, or `XSAVE`) in the derivatives program, but you cannot use input commands (such as `DATA LIST`, `GET`, `MATCH FILES`, or `ADD FILES`).
- To name the derivatives, specify the prefix *D*. before each parameter name. For example, the derivative name for the parameter *PARMI* must be *D.PARMI*.
- When a derivative has been calculated by a transformation, the variable for that derivative can be used in subsequent transformations.
- You do not need to supply all of the derivatives. Those derivatives that are not supplied will be estimated by the program. During the first iteration of the nonlinear estimation procedure, derivatives that are calculated in the derivatives program are compared with numerically calculated derivatives. This process serves as a check on the supplied values (see [CRITERIA Subcommand on p. 1334](#)).
- Transformations in the derivatives program are used by `[C]NLR` only and do not affect the active dataset.
- For `NLR`, the derivative of each parameter must be computed with respect to the predicted function (see [LOSS Subcommand on p. 1338](#)).

### ***Example***

```
MODEL PROGRAM A=1, B=0, C=1, D=0.
COMPUTE PRED = A * exp(B * X) + C * exp(D * X) .
DERIVATIVES .
COMPUTE D.A = exp (B * X) .
COMPUTE D.B = A * exp (B * X) * X .
COMPUTE D.C = exp (D * X) .
COMPUTE D.D = C * exp (D * X) * X .
```

- The derivatives program specifies derivatives of the `PRED` function for the sum of the two exponentials in the model described by the following equation:

$$Y = Ae^{Bx} + Ce^{Dx}$$

**Example**

```
DERIVATIVES.
COMPUTE D.A = exp (B * X) .
COMPUTE D.B = A * X * D.A.
COMPUTE D.C = exp (D * X) .
COMPUTE D.D = C * X * D.C.
```

- This example is an alternative way to express the same derivatives program that was specified in the previous example.

**CONSTRAINED FUNCTIONS Command**

The optional `CONSTRAINED FUNCTIONS` command signifies the beginning of the constrained functions program, which specifies nonlinear constraints. The constrained functions program is specified after the model program and the derivatives program (when used). It can only be used with, and must precede, the `CNLR` command. [For more information, see the topic BOUNDS Subcommand on p. 1337.](#)

**Example**

```
MODEL PROGRAM A=.5 B=1.6.
COMPUTE PRED=A*SPEED**B.

CONSTRAINED FUNCTIONS.
COMPUTE CF=A-EXP(B) .

CNLR STOP
/BOUNDS CF LE 0.
```

**CLEAR MODEL PROGRAMS Command**

`CLEAR MODEL PROGRAMS` deletes all transformations that are associated with the previously submitted model program, derivative program, and/or constrained functions program. It is primarily used in interactive mode to remove temporary variables that were created by these programs without affecting the active dataset or variables that were created by other transformation programs or temporary programs. It allows you to specify new models, derivatives, or constrained functions without having to run `[C]NLR`.

It is not necessary to use this command if you have already executed the `[C]NLR` procedure. Temporary variables that are associated with the procedure are automatically deleted.

**CNLR and NLR Commands**

Either the `CNLR` or the `NLR` command is required to specify the dependent and independent variables for the nonlinear regression.

- For either `CNLR` or `NLR`, the minimum specification is a dependent variable.
- Only one dependent variable can be specified. It must be a numeric variable in the active dataset and cannot be a variable that is generated by the model or the derivatives program.



## **OUTFILE Subcommand**

OUTFILE stores final parameter estimates for use on a subsequent [C]NLR command. The only specification on OUTFILE is the target file. Some or all of the values from this file can be read into a subsequent [C]NLR procedure with the FILE subcommand. The parameter data file that is created by OUTFILE stores the following variables:

- All of the split-file variables. OUTFILE writes one case of values for each split-file group in the active dataset.
- All of the parameters named on the MODEL PROGRAM command.
- The labels, formats, and missing values of the split-file variables and parameters defined for them previous to their use in the [C]NLR procedure.
- The sum of squared residuals (named *SSE*). *SSE* has no labels or missing values. The print and write format for *SSE* is F10.8.
- The number of cases on which the analysis was based (named *NCASES*). *NCASES* has no labels or missing values. The print and write format for *NCASES* is F8.0.

When OUTFILE is used, the model program cannot create variables named *SSE* or *NCASES*.

### **Example**

```
MODEL PROGRAM A=.5 B=1.6.  
COMPUTE PRED=A*SPEED*B.  
NLR STOP /OUTFILE=PARAM.
```

- OUTFILE generates a parameter data file containing one case for four variables: *A*, *B*, *SSE*, and *NCASES*.

## **FILE Subcommand**

FILE reads starting values for the parameters from a parameter data file that was created by an OUTFILE subcommand from a previous [C]NLR procedure. When starting values are read from a file, they do not have to be specified on the MODEL PROGRAM command. Rather, the MODEL PROGRAM command simply names the parameters that correspond to the parameters in the data file.

- The only specification on FILE is the file that contains the starting values.
- Some new parameters may be specified for the model on the MODEL PROGRAM command while other parameters are read from the file that is specified on the FILE subcommand.
- You do not have to name the parameters on MODEL PROGRAM in the order in which they occur in the parameter data file. In addition, you can name a partial list of the variables that are contained in the file.
- If the starting value for a parameter is specified on MODEL PROGRAM, the specification overrides the value that is read from the parameter data file.
- If split-file processing is in effect, the starting values for the first subfile are taken from the first case of the parameter data file. Subfiles are matched with cases in order until the starting-value file runs out of cases. All subsequent subfiles use the starting values for the last case.

- To read starting values from a parameter data file and then replace those values with the final results from [C]NLR, specify the same file on the `FILE` and `OUTFILE` subcommands. The input file is read completely before anything is written in the output file.

### Example

```
MODEL PROGRAM A B C=1 D=3.
COMPUTE PRED=A*SPEED**B + C*SPEED**D.
NLR STOP /FILE=PARAM /OUTFILE=PARAM.
```

- `MODEL PROGRAM` names four of the parameters that are used to calculate *PRED* but assigns values to only *C* and *D*. The values of *A* and *B* are read from the existing data file *PARAM*.
- After `NLR` computes the final estimates of the four parameters, `OUTFILE` writes over the old input file. If, in addition to these new final estimates, the former starting values of *A* and *B* are still desired, specify a different file on the `OUTFILE` subcommand.

## PRED Subcommand

`PRED` identifies the variable holding the predicted values.

- The only specification is a variable name, which must be identical to the variable name that is used to calculate predicted values in the model program.
- If the model program names the variable *PRED*, the `PRED` subcommand can be omitted. Otherwise, the `PRED` subcommand is required.
- The variable for predicted values is not saved in the active dataset unless the `SAVE` subcommand is used.

### Example

```
MODEL PROGRAM A=.5 B=1.6.
COMPUTE PSTOP=A*SPEED**B.
NLR STOP /PRED=PSTOP.
```

- `COMPUTE` in the model program creates a variable named *PSTOP* to temporarily store the predicted values for the dependent variable *STOP*.
- `PRED` identifies *PSTOP* as the variable that is used to define the model for the `NLR` procedure.

## SAVE Subcommand

`SAVE` is used to save the temporary variables for the predicted values, residuals, and derivatives that are created by the model and the derivatives programs.

- The minimum specification is a single keyword.
- The variables to be saved must have unique names on the active dataset. If a naming conflict exists, the variables are not saved.
- Temporary variables—for example, variables that are created after a `TEMPORARY` command and parameters that are specified by the model program—are not saved in the active dataset. They will not cause naming conflicts.

The following keywords are available and can be used in any combination and in any order. The new variables are always appended to the active dataset in the order in which these keywords are presented here:

<b>PRED</b>	<i>Save the predicted values.</i> The variable's name, label, and formats are those specified for it (or assigned by default) in the model program.
<b>RESID [(varname)]</b>	<i>Save the residuals variable.</i> You can specify a variable name in parentheses following the keyword. If no variable name is specified, the name of this variable is the same as the specification that you use for this keyword. For example, if you use the three-character abbreviation RES, the default variable name will be RES. The variable has the same print and write format as the predicted values variable that is created by the model program. It has no variable label and no user-defined missing values. It is system-missing for any case in which either the dependent variable is missing or the predicted value cannot be computed.
<b>DERIVATIVES</b>	<i>Save the derivative variables.</i> The derivative variables are named with the prefix <i>D.</i> to the first six characters of the parameter names. Derivative variables use the print and write formats of the predicted values variable and have no value labels or user-missing values. Derivative variables are saved in the same order as the parameters named on MODEL PROGRAM. Derivatives are saved for all parameters, whether or not the derivative was supplied in the derivatives program.
<b>LOSS</b>	<i>Save the user-specified loss function variable.</i> This specification is available only with CNLR and only if the LOSS subcommand has been specified.

Asymptotic standard errors of predicted values and residuals, and special residuals used for outlier detection and influential case analysis are not provided by the [C]NLR procedure. However, for a squared loss function, the asymptotically correct values for all these statistics can be calculated by using the SAVE subcommand with [C]NLR and then using the REGRESSION procedure. In REGRESSION, the dependent variable is still the same, and derivatives of the model parameters are used as independent variables. Casewise plots, standard errors of prediction, partial regression plots, and other diagnostics of the regression are valid for the nonlinear model.

### Example

```
MODEL PROGRAM A=.5 B=1.6.
COMPUTE PSTOP=A*SPEED**B.
NLR STOP /PRED=PSTOP
      /SAVE=RESID(RSTOP) DERIVATIVES PRED.
REGRESSION VARIABLES=STOP D.A D.B /ORIGIN
      /DEPENDENT=STOP /ENTER D.A D.B /RESIDUALS.
```

- The SAVE subcommand creates the residuals variable *RSTOP* and the derivative variables *D.A* and *D.B*.
- Because the PRED subcommand identifies *PSTOP* as the variable for predicted values in the nonlinear model, keyword PRED on SAVE adds the variable *PSTOP* to the active dataset.
- The new variables are added to the active dataset in the following order: *PSTOP*, *RSTOP*, *D.A*, and *D.B*.
- The subcommand RESIDUALS for REGRESSION produces the default analysis of residuals.

## **CRITERIA Subcommand**

CRITERIA controls the values of the cutoff points that are used to stop the iterative calculations in [C]NLR.

- The minimum specification is any of the criteria keywords and an appropriate value. The value can be specified in parentheses after an equals sign, a space, or a comma. Multiple keywords can be specified in any order. Defaults are in effect for keywords that are not specified.
- Keywords available for CRITERIA differ between CNLR and NLR and are discussed separately. However, with both CNLR and NLR, you can specify the critical value for derivative checking.

### **Checking Derivatives for CNLR and NLR**

Upon entering the first iteration, [C]NLR always checks any derivatives that are calculated on the derivatives program by comparing them with numerically calculated derivatives. For each comparison, it computes an agreement score. A score of 1 indicates agreement to machine precision; a score of 0 indicates definite disagreement. If a score is less than 1, either an incorrect derivative was supplied or there were numerical problems in estimating the derivative. The lower the score, the more likely it is that the supplied derivatives are incorrect. Highly correlated parameters may cause disagreement even when a correct derivative is supplied. Be sure to check the derivatives if the agreement score is not 1.

During the first iteration, [C]NLR checks each derivative score. If any score is below 1, it begins displaying a table to show the worst (lowest) score for each derivative. If any score is below the critical value, the program stops.

To specify the critical value, use the following keyword on CRITERIA:

**CKDER n**                      *Critical value for derivative checking.* Specify a number between 0 and 1 for *n*. The default is 0.5. Specify 0 to disable this criterion.

### **Iteration Criteria for CNLR**

The CNLR procedure uses NPSOL (Version 4.0) Fortran Package for Nonlinear Programming (Gill, Murray, Saunders, and Wright, 1986). The CRITERIA subcommand of CNLR gives the control features of NPSOL. The following section summarizes the NPSOL documentation.

CNLR uses a sequential quadratic programming algorithm, with a quadratic programming subproblem to determine the search direction. If constraints or bounds are specified, the first step is to find a point that is feasible with respect to those constraints. Each major iteration sets up a quadratic program to find the search direction,  $p$ . Minor iterations are used to solve this subproblem. Then, the major iteration determines a steplength  $\alpha$  by a line search, and the function is evaluated at the new point. An optimal solution is found when the optimality tolerance criterion is met.

The CRITERIA subcommand has the following keywords when used with CNLR:

<b>ITER n</b>	<i>Maximum number of major iterations.</i> Specify any positive integer for $n$ . The default is $\max(50, 3(p+m_L)+10m_N)$ , where $p$ is the number of parameters, $m_L$ is the number of linear constraints, and $m_N$ is the number of nonlinear constraints. If the search for a solution stops because this limit is exceeded, CNLR issues a warning message.
<b>MINORITERATION n</b>	<i>Maximum number of minor iterations.</i> Specify any positive integer. This value is the number of minor iterations allowed within each major iteration. The default is $\max(50, 3(n+m_L+m_N))$ .
<b>CRSHTOL n</b>	<i>Crash tolerance.</i> CRSHTOL is used to determine whether initial values are within their specified bounds. Specify any value between 0 and 1. The default value is 0.01. A constraint of the form $a'X \geq l$ is considered a valid part of the working set if $ a'X-l  < \text{CRSHTOL}(1+ l )$ .
<b>STEPLIMIT n</b>	<i>Step limit.</i> The CNLR algorithm does not allow changes in the length of the parameter vector to exceed a factor of $n$ . The limit prevents very early steps from going too far from good initial estimates. Specify any positive value. The default value is 2.
<b>FTOLERANCE n</b>	<i>Feasibility tolerance.</i> This value is the maximum absolute difference allowed for both linear and nonlinear constraints for a solution to be considered feasible. Specify any value greater than 0. The default value is the square root of your machine's epsilon.
<b>LFTOLERANCE n</b>	<i>Linear feasibility tolerance.</i> If specified, this value overrides FTOLERANCE for linear constraints and bounds. Specify any value greater than 0. The default value is the square root of your machine's epsilon.
<b>NFTOLERANCE n</b>	<i>Nonlinear feasibility tolerance.</i> If specified, this value overrides FTOLERANCE for nonlinear constraints. Specify any value greater than 0. The default value is the square root of your machine's epsilon.
<b>LSTOLERANCE n</b>	<i>Line search tolerance.</i> This value must be between 0 and 1 (but not including 1). It controls the accuracy required of the line search that forms the innermost search loop. The default value, 0.9, specifies an inaccurate search. This setting is appropriate for many problems, particularly if nonlinear constraints are involved. A smaller positive value, corresponding to a more accurate line search, may give better performance if there are no nonlinear constraints, all (or most) derivatives are supplied in the derivatives program, and the data fit in memory.
<b>OPTOLERANCE n</b>	<i>Optimality tolerance.</i> If an iteration point is a feasible point, and the next step will not produce a relative change in either the parameter vector or the objective function of more than the square root of OPTOLERANCE, an optimal solution has been found. OPTOLERANCE can also be thought of as the number of significant digits in the objective function at the solution. For example, if $\text{OPTOLERANCE}=10^{-6}$ , the objective function should have approximately six significant digits of accuracy. Specify any number between the FPRECISION value and 1. The default value for OPTOLERANCE is $\text{epsilon}^{**}0.8$ .

<b>FPRECISION n</b>	<i>Function precision.</i> This measure is a measure of the accuracy with which the objective function can be checked. It acts as a relative precision when the function is large and an absolute precision when the function is small. For example, if the objective function is larger than 1, and six significant digits are desired, <code>FPRECISION</code> should be 1E-6. If, however, the objective function is of the order 0.001, <code>FPRECISION</code> should be 1E-9 to get six digits of accuracy. Specify any number between 0 and 1. The choice of <code>FPRECISION</code> can be very complicated for a badly scaled problem. Chapter 8 of Gill et al. (1981) gives some scaling suggestions. The default value is <code>epsilon**0.9</code> .
<b>ISTEP n</b>	<i>Infinite step size.</i> This value is the magnitude of the change in parameters that is defined as infinite. That is, if the change in the parameters at a step is greater than <code>ISTEP</code> , the problem is considered unbounded, and estimation stops. Specify any positive number. The default value is 1E+20.

### Iteration Criteria for NLR

The NLR procedure uses an adaptation of subroutine LMSTR from the MINPACK package by Garbow et al. Because the NLR algorithm differs substantially from CNLR, the `CRITERIA` subcommand for NLR has a different set of keywords.

NLR computes parameter estimates by using the Levenberg-Marquardt method. At each iteration, NLR evaluates the estimates against a set of control criteria. The iterative calculations continue until one of five cutoff points is met, at which point the iterations stop and the reason for stopping is displayed.

The `CRITERIA` subcommand has the following keywords when used with NLR:

<b>ITER n</b>	<i>Maximum number of major and minor iterations allowed.</i> Specify any positive integer for <i>n</i> . The default is 100 iterations per parameter. If the search for a solution stops because this limit is exceeded, NLR issues a warning message.
<b>SSCON n</b>	<i>Convergence criterion for the sum of squares.</i> Specify any non-negative number for <i>n</i> . The default is 1E-8. If successive iterations fail to reduce the sum of squares by this proportion, the procedure stops. Specify 0 to disable this criterion.
<b>PCON n</b>	<i>Convergence criterion for the parameter values.</i> Specify any non-negative number for <i>n</i> . The default is 1E-8. If successive iterations fail to change any of the parameter values by this proportion, the procedure stops. Specify 0 to disable this criterion.
<b>RCON n</b>	<i>Convergence criterion for the correlation between the residuals and the derivatives.</i> Specify any non-negative number for <i>n</i> . The default is 1E-8. If the largest value for the correlation between the residuals and the derivatives equals this value, the procedure stops because it lacks the information that it needs to estimate a direction for its next move. This criterion is often referred to as a gradient convergence criterion. Specify 0 to disable this criterion.

#### Example

```
MODEL PROGRAM A=.5 B=1.6.
COMPUTE PRED=A*SPEED**B.
NLR STOP /CRITERIA=ITER(80) SSCON=.000001.
```

- `CRITERIA` changes two of the five cutoff values affecting iteration, `ITER` and `SSCON`, and leaves the remaining three, `PCON`, `RCON`, and `CKDER`, at their default values.

## **BOUNDS Subcommand**

The BOUNDS subcommand can be used to specify both linear and nonlinear constraints. It can be used only with CNLR; it cannot be used with NLR.

### **Simple Bounds and Linear Constraints**

BOUNDS can be used to impose bounds on parameter values. These bounds can involve either single parameters or a linear combination of parameters and can be either equalities or inequalities.

- All bounds are specified on the same BOUNDS subcommand and are separated by semicolons.
- The only variables that are allowed on BOUNDS are parameter variables (those variables that are named on MODEL PROGRAM).
- Only \* (multiplication), + (addition), - (subtraction), = or EQ, >= or GE, and <= or LE can be used. When two relational operators are used (as in the third bound in the example below), they must both be in the same direction.

#### **Example**

```
/BOUNDS 5 >= A;
        B >= 9;
        .01 <= 2*A + C <= 1;
        D + 2*E = 10
```

- BOUNDS imposes bounds on the parameters *A*, *B*, *C*, and *D*. Specifications for each parameter are separated by a semicolon.

### **Nonlinear Constraints**

Nonlinear constraints on the parameters can also be specified with the BOUNDS subcommand. The constrained function must be calculated and stored in a variable by a constrained functions program directly preceding the CNLR command. The constraint is then specified on the BOUNDS subcommand.

In general, nonlinear bounds will not be obeyed until an optimal solution has been found. This process is different from simple and linear bounds, which are satisfied at each iteration. The constrained functions must be smooth near the solution.

#### **Example**

```
MODEL PROGRAM A=.5 B=1.6.
COMPUTE PRED=A*SPEED**B.

CONSTRAINED FUNCTIONS.
COMPUTE DIFF=A-10**B.

CNLR STOP /BOUNDS DIFF LE 0.
```

- The constrained function is calculated by a constrained functions program and stored in variable *DIFF*. The constrained functions program immediately precedes CNLR.

- BOUNDS imposes bounds on the function (less than or equal to 0).
- CONSTRAINED FUNCTIONS variables and parameters that are named on MODEL PROGRAM cannot be combined in the same BOUNDS expression. For example, you *cannot* specify  $(DIFF + A) \geq 0$  on the BOUNDS subcommand.

## LOSS Subcommand

LOSS specifies a loss function for CNLR to minimize. By default, CNLR minimizes the sum of squared residuals. LOSS can be used only with CNLR; it cannot be used with NLR.

- The loss function must first be computed in the model program. LOSS is then used to specify the name of the computed variable.
- The minimizing algorithm may fail if it is given a loss function that is not smooth, such as the absolute value of residuals.
- If derivatives are supplied, the derivative of each parameter must be computed with respect to the loss function, rather than the predicted value. The easiest way to do this is in two steps: First compute derivatives of the model, and then compute derivatives of the loss function with respect to the model and multiply by the model derivatives.
- When LOSS is used, the usual summary statistics are not computed. Standard errors, confidence intervals, and correlations of the parameters are available only if the BOOTSTRAP subcommand is specified.

### Example

```
MODEL PROGRAM A=1 B=1.
COMPUTE PRED=EXP(A+B*T) / (1+EXP(A+B*T)) .
COMPUTE LOSS=-W*(Y*LN(PRED) + (1-Y)*LN(1-PRED)) .

DERIVATIVES.
COMPUTE D.A=PRED / (1+EXP(A+B*T)) .
COMPUTE D.B=T*PRED / (1+EXP(A+B*T)) .
COMPUTE D.A=(-W*(Y/PRED - (1-Y)/(1-PRED))) * D.A .
COMPUTE D.B=(-W*(Y/PRED - (1-Y)/(1-PRED))) * D.B .

CNLR Y /LOSS=LOSS.
```

- The second COMPUTE command in the model program computes the loss function and stores its values in the variable *LOSS*, which is then specified on the LOSS subcommand.
- Because derivatives are supplied in the derivatives program, the derivatives of all parameters are computed with respect to the loss function, rather than the predicted value.

## BOOTSTRAP Subcommand

BOOTSTRAP provides bootstrap estimates of the parameter standard errors, confidence intervals, and correlations. BOOTSTRAP can be used only with CNLR; it cannot be used with NLR.

**Bootstrapping** is a way of estimating the standard error of a statistic, using repeated samples from the original data set. This process is done by sampling with replacement to get samples of the same size as the original data set.



- The minimum specification is the subcommand keyword. Optionally, specify the number of samples to use for generating bootstrap results.
- By default, `BOOTSTRAP` generates bootstrap results based on  $10 * p * (p + 1) / 2$  samples, where  $p$  is the number of parameters. That is, 10 samples are drawn for each statistic (standard error or correlation) to be calculated.
- When `BOOTSTRAP` is used, the nonlinear equation is estimated for each sample. The standard error of each parameter estimate is then calculated as the standard deviation of the bootstrapped estimates. Parameter values from the original data are used as starting values for each bootstrap sample. Even so, bootstrapping is computationally expensive.
- If the `OUTFILE` subcommand is specified, a case is written to the output file for each bootstrap sample. The first case in the file will be the actual parameter estimates, followed by the bootstrap samples. After the first case is eliminated (using `SELECT IF`), other procedures (such as `FREQUENCIES`) can be used to examine the bootstrap distribution.

### Example

```
MODEL PROGRAM A=.5 B=1.6.
COMPUTE PSTOP=A*SPEED**B.
CNLR STOP /BOOTSTRAP /OUTFILE=PARAM.
GET FILE=PARAM.
LIST.
COMPUTE ID=$CASENUM.
SELECT IF (ID > 1).
FREQUENCIES A B /FORMAT=NOTABLE /HISTOGRAM.
```

- `CNLR` generates the bootstrap standard errors, confidence intervals, and parameter correlation matrix. `OUTFILE` saves the bootstrap estimates in the file *PARAM*.
- `GET` retrieves the system file *PARAM*.
- `LIST` lists the different sample estimates, along with the original estimate. *NCASES* in the listing (see [OUTFILE Subcommand on p. 1331](#)) refers to the number of distinct cases in the sample because cases are duplicated in each bootstrap sample.
- `FREQUENCIES` generates histograms of the bootstrapped parameter estimates.

## References

Gill, P. E., W. M. Murray, M. A. Saunders, and M. H. Wright. 1986. *User's guide for NPSOL (version 4.0): A FORTRAN package for nonlinear programming. Technical Report SOL 86-2*. Stanford University: Department of Operations Research.

# NOMREG

NOMREG is available in the Regression option.

```
NOMREG dependent varname [(BASE = {FIRST } ORDER = {ASCENDING**})] [BY factor list]
                               {LAST**}           {DATA      }
                               {value  }           {DESCENDING }
      [WITH covariate list]

[/CRITERIA = [CIN({95**})] [DELTA({0**})] [MXITER({100**})] [MXSTEP({5**})]
              {n }           {n }           {n }           {n }
              [LCONVERGE({0**})] [PCONVERGE({1.0E-6**})] [SINGULAR({1E-8**})]
              {n }           {n }           {n }           {n }
              [BIAS({0**})] [CHKSEP({20**})] ]
              {n }           {n }

[/FULLFACTORIAL]

[/INTERCEPT = {EXCLUDE  }
                {INCLUDE** }

[/MISSING = {EXCLUDE**}
            {INCLUDE  }

[/MODEL = {[effect effect ...]} [| {BACKWARD} = { effect effect ...}]
           {FORWARD }
           {BSTEP  }
           {FSTEP  }

[/STEPWISE = [RULE({SINGLE**  })] [MINEFFECT({0**  })] [MAXEFFECT(n)]
             {SFACOR      } {value}
             {CONTAINMENT}
             {NONE       }

             [PIN({0.05**})] [POUT({0.10**})]
             {value } {value }

             [ENTRYMETHOD({LR** })] [REMOVALMETHOD({LR**})]
             {SCORE} {WALD}

[/OUTFILE = [{MODEL  } (filename)]
            {PARAMETER}

[/PRINT = [CELLPROB] [CLASSTABLE] [CORB] [HISTORY({1**})] [IC] ]
          {n }
          [SUMMARY ] [PARAMETER ] [COVB] [FIT] [LRT] [KERNEL]
          [ASSOCIATION] [CPS**] [STEP**] [MFI**] [NONE]

[/SAVE = [ACPROB[(newname)]] [ESTPROB[(rootname[:{25**})]] ]
        {n }
        [PCPROB[(newname)]] [PREDCAT[(newname)]]

[/SCALE = {1**  }
          {n }
          {DEVIANCE}
          {PEARSON }

[/SUBPOP = varlist]

[/TEST[(valuelist)] = {'label' effect valuelist effect valuelist...;}
                    {'label' ALL list; }
                    {'label' ALL list }
                    }
```

\*\* Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- ENTRYMETHOD keyword introduced on STEPWISE subcommand.
- REMOVALMETHOD keyword introduced on STEPWISE subcommand.
- IC keyword introduced on PRINT subcommand.

Release 15.0

- ASSOCIATION keyword introduced on PRINT subcommand.

Release 17.0

- Added support for SET THREADS and SET MCACHE.

### **Example**

NOMREG response.

## **Overview**

NOMREG is a procedure for fitting a multinomial logit model to a polytomous nominal dependent variable.

### **Options**

**Tuning the algorithm.** You can control the values of algorithm-tuning parameters with the CRITERIA subcommand.

**Optional output.** You can request additional output through the PRINT subcommand.

**Exporting the model.** You can export the model to an external file. The model information will be written using the Extensible Markup Language (XML).

### **Basic Specification**

The basic specification is one dependent variable.

### **Operations**

- This procedure uses the multithreaded options specified by SET THREADS and SET MCACHE.

### **Syntax Rules**

- Minimum syntax—at least one dependent variable must be specified.
- The variable specification must come first.
- Subcommands can be specified in any order.

- Empty subcommands except the `MODEL` subcommand are ignored.
- The `MODEL` and the `FULLFACTORIAL` subcommands are mutually exclusive. Only one of them can be specified at any time.
- The `MODEL` subcommand stepwise options and the `TEST` subcommand are mutually exclusive. Only one of them can be specified at any time.
- When repeated subcommands except the `TEST` subcommand are specified, all specifications except the last valid one are discarded.
- The following words are reserved as keywords or internal commands in the `NOMREG` procedure: `BY`, `WITH`, and `WITHIN`.
- The set of factors and covariates used in the `MODEL` subcommand (or implied on the `FULLFACTORIAL` subcommand) must be a subset of the variable list specified or implied on the `SUBPOP` subcommand.

## Examples

```
NOMREG
  bfast BY agecat gender active
  /CRITERIA = CIN(95) DELTA(0) MXITER(100) MXSTEP(5) CHKSEP(20) LCONVERGE(0)
  PCONVERGE(1.0E-6) SINGULAR(1.0E-8)
  /MODEL = agecat active
  /INTERCEPT = INCLUDE
  /PRINT = CELLPROB CLASSTABLE FIT PARAMETER SUMMARY LRT MFI CPS.
```

- The procedure fits a model for *bfast* using *agecat*, *gender*, and *active*.
- The `MODEL` subcommand specifies a main effects model using *agecat* and *active*. Though *gender* is not in the model, it is still used to define subpopulations for fit statistics.
- The `PRINT` subcommand requests cell probabilities, the classification table, fit statistics, parameter estimates, the model summary, likelihood ratio tests, model fitting information, and the case processing summary.
- All other options are set to their default values.

### 1-1 Matched Case Control

```
NOMREG
  claim WITH difhgt difwgt difnic difcvd difhaz
  /CRITERIA CIN(95) DELTA(0) MXITER(100) MXSTEP(5) CHKSEP(20) LCONVERGE(0)
  PCONVERGE(0.000001) SINGULAR(0.00000001)
  /MODEL
  /INTERCEPT EXCLUDE
  /PRINT PARAMETER SUMMARY LRT .
```

- The procedure fits a model for *claim* using variables *difhgt*, *difwgt*, *difnic*, *difcvd*, and *difhaz* as covariates. The values of these variables represent the difference in values between cases and controls.
- The empty `MODEL` subcommand indicates that a main-effects model using all the covariates will be created. The `INTERCEPT` subcommand excludes the intercept from the model.
- The `PRINT` subcommand requests the parameter estimates, model summary, and likelihood ratio tests.
- All other options are set to their default values.

## Variable List

The variable list specifies the dependent variable and the factors in the model.

- The dependent variable must be the first specification on NOMREG. It can be of any type (numeric or string). Values of the dependent variable are sorted according to the ORDER specification.

**ORDER = ASCENDING**      *Response categories are sorted in ascending order.* The lowest value defines the first category, and the highest value defines the last category.

**ORDER = DATA**      *Response categories are not sorted.* The first value encountered in the data defines the first category. The last distinct value defines the last category.

**ORDER = DESCENDING**      *Response categories are sorted in descending order.* The highest value defines the first category, and the lowest value defines the last category.

- By default, the last response category is used as the base (or reference) category. No model parameters are assigned to the base category. Use the BASE attribute to specify a custom base category.

**BASE = FIRST**      *The first category is the base category.*

**BASE = LAST**      *The last category is the base category.*

**BASE = value**      *The category with the specified value is the base category.* Put the value inside a pair of quotes if either the value is formatted (such as date or currency) or if the dependent variable is the string type.

- Factor variables can be of any type (numeric or string). The factors follow the dependent variable separated by the keyword BY.
- Covariate variables must be numeric. The covariates follow the factors, separated by the keyword WITH.
- Listwise deletion is used. If any variables in a case contain missing values, that case will be excluded.
- If the WEIGHT command was specified, the actual weight values are used for the respective category combination. No rounding or truncation will be done. However, cases with negative and zero weight values are excluded from the analyses.

### Example

```
NOMREG response (ORDER = DESCENDING BASE='No') BY factor1.
```

- Values of the variable *response* are sorted in descending order, and the category whose value is *No* is the base category.

### Example

```
NOMREG
movie BY gender date
/CRITERIA = CIN(95) DELTA(0) MXITER(100) MXSTEP(5) LCONVERGE(0)
           PCONVERGE(0)
/INTERCEPT = EXCLUDE
/PRINT = CLASSTABLE FIT PARAMETER SUMMARY LRT .
```

- The dependent variable is *movie*, and *gender* and *date* are factors.
- CRITERIA specifies that the confidence level to use is 95, no delta value should be added to cells with observed zero frequency, and neither the log-likelihood nor parameter estimates convergence criteria should be used. This means that the procedure will stop when either 100 iterations or five step-halving operations have been performed.
- INTERCEPT specifies that the intercept should be excluded from the model.
- PRINT specifies that the classification table, goodness-of-fit statistics, parameter statistics, model summary, and likelihood-ratio tests should be displayed.

## **CRITERIA Subcommand**

The CRITERIA subcommand offers controls on the iterative algorithm used for estimation and specifies numerical tolerance for checking singularity.

<b>BIAS(n)</b>	<i>Bias value added to observed cell frequency.</i> Specify a non-negative value less than 1. The default value is 0.
<b>CHKSEP(n)</b>	<i>Starting iteration for checking for complete separation.</i> Specify a non-negative integer. The default value is 20.
<b>CIN(n)</b>	<i>Confidence interval level.</i> Specify a value greater than or equal to 0 and less than 100. The default value is 95.
<b>DELTA(n)</b>	<i>Delta value added to zero cell frequency.</i> Specify a non-negative value less than 1. The default value is 0.
<b>LCONVERGE(n)</b>	<i>Log-likelihood function convergence criterion.</i> Convergence is assumed if the absolute change in the log-likelihood function is less than this value. The criterion is not used if the value is 0. Specify a non-negative value. The default value is 0.
<b>MXITER(n)</b>	<i>Maximum number of iterations.</i> Specify a positive integer. The default value is 100.
<b>MXSTEP(n)</b>	<i>Maximum step-halving allowed.</i> Specify a positive integer. The default value is 5.
<b>PCONVERGE(a)</b>	<i>Parameter estimates convergence criterion.</i> Convergence is assumed if the absolute change in the parameter estimates is less than this value. The criterion is not used if the value is 0. Specify a non-negative value. The default value is $10^{-6}$ .
<b>SINGULAR(a)</b>	<i>Value used as tolerance in checking singularity.</i> Specify a positive value. The default value is $10^{-8}$ .

## **FULLFACTORIAL Subcommand**

The FULLFACTORIAL subcommand generates a specific model: first, the intercept (if included); second, all of the covariates (if specified), in the order in which they are specified; next, all of the main factorial effects; next, all of the two-way factorial interaction effects, all of the three-way factorial interaction effects, and so on, up to the highest possible interaction effect.

- The FULLFACTORIAL and the MODEL subcommands are mutually exclusive. Only one of them can be specified at any time.
- The FULLFACTORIAL subcommand does not take any keywords.

## **INTERCEPT Subcommand**

The `INTERCEPT` subcommand controls whether intercept terms are included in the model. The number of intercept terms is the number of response categories less one.

<b>INCLUDE</b>	<i>Includes the intercept terms. This is the default.</i>
<b>EXCLUDE</b>	<i>Excludes the intercept terms.</i>

## **MISSING Subcommand**

By default, cases with missing values for any of the variables on the `NOMREG` variable list are excluded from the analysis. The `MISSING` subcommand allows you to include cases with user-missing values.

- Note that missing values are deleted at the subpopulation level.

<b>EXCLUDE</b>	<i>Excludes both user-missing and system-missing values. This is the default.</i>
<b>INCLUDE</b>	<i>User-missing values are treated as valid. System-missing values cannot be included in the analysis.</i>

## **MODEL Subcommand**

The `MODEL` subcommand specifies the effects in the model.

- The `MODEL` and the `FULLFACTORIAL` subcommands are mutually exclusive. Only one of them can be specified at any time.
- If more than one `MODEL` subcommand is specified, only the last one is in effect.
- Specify a list of terms to be included in the model, separated by commas or spaces. If the `MODEL` subcommand is omitted or empty, the default model is generated. The default model contains: first, the intercept (if included); second, all of the covariates (if specified), in the order in which they are specified; and next, all of the main factorial effects, in the order in which they are specified.
- If a `SUBPOP` subcommand is specified, then effects specified in the `MODEL` subcommand can only be composed using the variables listed on the `SUBPOP` subcommand.
- To include a main-effect term, enter the name of the factor on the `MODEL` subcommand.
- To include an interaction-effect term among factors, use the keyword `BY` or the asterisk (\*) to join factors involved in the interaction. For example, `A*B*C` means a three-way interaction effect of A, B, and C, where A, B, and C are factors. The expression `A BY B BY C` is equivalent to `A*B*C`. Factors inside an interaction effect must be distinct. Expressions such as `A*C*A` and `A*A` are invalid.
- To include a nested-effect term, use the keyword `WITHIN` or a pair of parentheses on the `MODEL` subcommand. For example, `A(B)` means that A is nested within B, where A and B are factors. The expression `A WITHIN B` is equivalent to `A(B)`. Factors inside a nested effect must be distinct. Expressions such as `A(A)` and `A(B*A)` are invalid.

- Multiple-level nesting is supported. For example,  $A(B(C))$  means that  $B$  is nested within  $C$ , and  $A$  is nested within  $B(C)$ . When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus,  $A(B)(C)$  is not valid.
- Nesting within an interaction effect is valid. For example,  $A(B*C)$  means that  $A$  is nested within  $B*C$ .
- Interactions among nested effects are allowed. The correct syntax is the interaction followed by the common nested effect inside the parentheses. For example, interaction between  $A$  and  $B$  within levels of  $C$  should be specified as  $A*B(C)$  instead of  $A(C)*B(C)$ .
- To include a covariate term in the model, enter the name of the covariate on the `MODEL` subcommand.
- Covariates can be connected, but not nested, using the keyword `BY` or the asterisk (`*`) operator. For example,  $X*X$  is the product of  $X$  and itself. This is equivalent to a covariate whose values are the square of those of  $X$ . However,  $X(Y)$  is invalid.
- Factor and covariate effects can be connected in many ways. No effects can be nested within a covariate effect. Suppose  $A$  and  $B$  are factors, and  $X$  and  $Y$  are covariates. Examples of valid combination of factor and covariate effects are  $A*X$ ,  $A*B*X$ ,  $X(A)$ ,  $X(A*B)$ ,  $X*A(B)$ ,  $X*Y(A*B)$ , and  $A*B*X*Y$ .
- A stepwise method can be specified by following the model effects with a vertical bar (`|`), a stepwise method keyword, an equals sign (`=`), and a list of variables (or interactions or nested effects) for which the method is to be used.
- If a stepwise method is specified, then the `TEST` subcommand is ignored.
- If a stepwise method is specified, then it begins with the results of the model defined on the left side of the `MODEL` subcommand.
- If a stepwise method is specified but no effects are specified on the left side of the `MODEL` subcommand, then the initial model contains the intercept only (if `INTERCEPT = INCLUDE`) or the initial model is the null model (if `INTERCEPT = EXCLUDE`).
- The intercept cannot be specified as an effect in the stepwise method option.
- For all stepwise methods, if two effects have tied significance levels, then the removal or entry is performed on the effect specified first. For example, if the right side of the `MODEL` subcommand specifies `FORWARD A*B A(B)`, where  $A*B$  and  $A(B)$  have the same significance level less than `PIN`, then  $A*B$  is entered because it is specified first.

The available stepwise method keywords are:

- BACKWARD** *Backward elimination.* As a first step, the variables (or interaction effects or nested effects) specified on `BACKWARD` are entered into the model together and are tested for removal one by one. The variable with the largest significance level of the likelihood-ratio statistic, provided that the value is larger than `POUT`, is removed, and the model is reestimated. This process continues until no more variables meet the removal criterion or when the current model is the same as a previous model.
- FORWARD** *Forward entry.* The variables (or interaction effects or nested effects) specified on `FORWARD` are tested for entry into the model one by one, based on the significance level of the likelihood-ratio statistic. The variable with the smallest significance level less than `PIN` is entered into the model, and the model is reestimated. Model building stops when no more variables meet the entry criteria.



- BSTEP** *Backward stepwise.* As a first step, the variables (or interaction effects or nested effects) specified on **BSTEP** are entered into the model together and are tested for removal one by one. The variable with the largest significance level of the likelihood-ratio statistic, provided that the value is larger than **POUT**, is removed, and the model is reestimated. This process continues until no more variables meet the removal criterion. Next, variables not in the model are tested for possible entry, based on the significance level of the likelihood-ratio statistic. The variable with the smallest significance level less than **PIN** is entered, and the model is reestimated. This process repeats, with variables in the model again evaluated for removal. Model building stops when no more variables meet the removal or entry criteria or when the current model is the same as a previous model.
- FSTEP** *Forward stepwise.* The variables (or interaction effects or nested effects) specified on **FSTEP** are tested for entry into the model one by one, based on the significance level of the likelihood-ratio statistic. The variable with the smallest significance level less than **PIN** is entered into the model, and the model is reestimated. Next, variables that are already in the model are tested for removal, based on the significance level of the likelihood-ratio statistic. The variable with the largest probability greater than the specified **POUT** value is removed, and the model is reestimated. Variables in the model are then evaluated again for removal. Once no more variables satisfy the removal criterion, variables not in the model are evaluated again for entry. Model building stops when no more variables meet the entry or removal criteria or when the current model is the same as a previous one.

### Examples

```
NOMREG y BY a b c
/INTERCEPT = INCLUDE
/MODEL = a b c | BACKWARD = a*b a*c b*c a*b*c.
```

- The initial model contains the intercept and main effects a, b, and c. Backward elimination is used to select among the two- and three-way interaction effects.

```
NOMREG y BY a b c
/MODEL = INTERCEPT | FORWARD = a b c.
```

- The initial model contains the intercept. Forward entry is used to select among the main effects a, b, and c.

```
NOMREG y BY a b c
/INTERCEPT = INCLUDE
/MODEL = | FORWARD = a b c.
```

- The initial model contains the intercept. Forward entry is used to select among the main effects a, b, and c.

```
NOMREG y BY a b c
/INTERCEPT = EXCLUDE
/MODEL = | BSTEP = a b c.
```

- The initial model is the null model. Backward stepwise is used to select among the main effects a, b, and c.

```
NOMREG y BY a b c
/MODEL = | FSTEP =.
```

- This **MODEL** specification yields a syntax error.

## **STEPWISE Subcommand**

The `STEPWISE` subcommand gives you control of the statistical criteria when stepwise methods are used to build a model. This subcommand is ignored if a stepwise method is not specified on the `MODEL` subcommand.

<b>RULE(keyword)</b>	<i>Rule for entering or removing terms in stepwise methods.</i> The default <code>SINGLE</code> indicates that only one effect can be entered or removed at a time, provided that the hierarchy requirement is satisfied for all effects in the model. <code>SFACTOR</code> indicates that only one effect can be entered or removed at a time, provided that the hierarchy requirement is satisfied for all factor-only effects in the model. <code>CONTAINMENT</code> indicates that only one effect can be entered or removed at a time, provided that the containment requirement is satisfied for all effects in the model. <code>NONE</code> indicates that only one effect can be entered or removed at a time, where neither the hierarchy nor the containment requirement need be satisfied for any effects in the model.
<b>MINEFFECT(n)</b>	<i>Minimum number of effects in final model.</i> The default is 0. The intercept, if any, is not counted among the effects. This criterion is ignored unless one of the stepwise methods <code>BACKWARD</code> or <code>BSTEP</code> is specified.
<b>MAXEFFECT(n)</b>	<i>Maximum number of effects in final model.</i> The default value is the total number of effects specified or implied on the <code>NOMREG</code> command. The intercept, if any, is not counted among the effects. This criterion is ignored unless one of the stepwise methods <code>FORWARD</code> or <code>FSTEP</code> is specified.
<b>ENTRYMETHOD(keyword)</b>	<i>Method for entering terms in stepwise methods.</i> The default <code>LR</code> indicates that the likelihood ratio test is used to determine whether a term is entered into the model. <code>SCORE</code> indicates that the score test is used. This criterion is ignored unless one of the stepwise methods <code>FORWARD</code> , <code>BSTEP</code> , or <code>FSTEP</code> is specified.
<b>REMOVALMETHOD(keyword)</b>	<i>Method for removing terms in stepwise methods.</i> The default <code>LR</code> indicates that the likelihood ratio test is used to determine whether a term is entered into the model. <code>WALD</code> indicates that the Wald test is used. This criterion is ignored unless one of the stepwise methods <code>BACKWARD</code> , <code>BSTEP</code> , or <code>FSTEP</code> is specified.
<b>PIN(a)</b>	<i>Probability of the likelihood-ratio statistic for variable entry.</i> The default is 0.05. The larger the specified probability, the easier it is for a variable to enter the model. This criterion is ignored unless one of the stepwise methods <code>FORWARD</code> , <code>BSTEP</code> , or <code>FSTEP</code> is specified.
<b>POUT(a)</b>	<i>Probability of the likelihood-ratio statistic for variable removal.</i> The default is 0.1. The larger the specified probability, the easier it is for a variable to remain in the model. This criterion is ignored unless one of the stepwise methods <code>BACKWARD</code> , <code>BSTEP</code> , or <code>FSTEP</code> is specified.

The hierarchy requirement stipulates that among the effects specified or implied on the `MODEL` subcommand, for any effect to be in a model, all lower-order effects that are part of the former effect must also be in the model. For example, if `A`, `X`, and `A*X` are specified, then for `A*X` to be in a model, the effects `A` and `X` must also be in the model.

The containment requirement stipulates that among the effects specified or implied on the `MODEL` subcommand, for any effect to be in the model, all effects contained in the former effect must also be in the model. For any two effects  $F$  and  $F'$ ,  $F$  is contained in  $F'$  if:

- Both effects  $F$  and  $F'$  involve the same covariate effect, if any. (Note that effects  $A*X$  and  $A*X*X$  are not considered to involve the same covariate effect because the first involves covariate effect  $X$  and the second involves covariate effect  $X**2$ .)
- $F'$  consists of more factors than  $F$ .
- All factors in  $F$  also appear in  $F'$ .

The following table illustrates how the hierarchy and containment requirements relate to the `RULE` options. Each row of the table gives a different set of effects specified on the `MODEL` subcommand. The columns correspond to the `RULE` options `SINGLE`, `SFACTOR`, and `CONTAINMENT`. The cells contain the order in which effects must occur in the model. For example, unless otherwise noted, all effects numbered 1 must be in the model for any effects numbered 2 to be in the model.

Table 153-1  
Hierarchy and containment requirements

Effects	SINGLE	SFACTOR	CONTAINMENT
A, B, A*B	1. A, B 2. A*B	1. A, B 2. A*B	1. A, B 2. A*B
X, X**2, X**3	1. X 2. X**2 3. X**3	Effects can occur in the model in any order.	Effects can occur in the model in any order.
A, X, X(A)	1. A, X 2. X(A)	Effects can occur in the model in any order.	1. X 2. X(A)  Effect A can occur in the model in any order.
A, X, X**2(A)	1. A, X 2. X**2(A)	Effects can occur in the model in any order.	Effects can occur in the model in any order.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand allows you to specify files to which output is written.

- Only one `OUTFILE` subcommand is allowed. If you specify more than one, only the last one is executed.

- You must specify at least one keyword and a valid filename in parentheses. There is no default.
- Neither MODEL nor PARAMETER is honored if split file processing is on (SPLIT FILE command) or if more than one dependent (DEPENDENT subcommand) variable is specified.

<b>MODEL(filename)</b>	<i>Write parameter estimates and their covariances to an XML (PMML) file. Specify the filename in full. NOMREG does not supply an extension. You can use this model file to apply the model information to other data files for scoring purposes. For more information, see the topic Scoring expressions on p. 116.</i>
<b>PARAMETER(filename)</b>	<i>Write parameter estimates only to an XML (PMML) file. Specify the filename in full. NOMREG does not supply an extension. You can use this model file to apply the model information to other data files for scoring purposes. For more information, see the topic Scoring expressions on p. 116.</i>

## PRINT Subcommand

The PRINT subcommand displays optional output. If no PRINT subcommand is specified, the default output includes a factor information table.

<b>ASSOCIATION</b>	<i>Measures of Monotone Association.</i> Displays a table with information on the number of concordant pairs, discordant pairs, and tied pairs. The Somers' D, Goodman and Kruskal's Gamma, Kendall's tau-a, and Concordance Index C are also displayed in this table.
<b>CELLPROB</b>	<i>Observed proportion, expected probability, and the residual for each covariate pattern and each response category.</i>
<b>CLASSTABLE</b>	<i>Classification table.</i> The square table of frequencies of observed response categories versus the predicted response categories. Each case is classified into the category with the highest predicted probability.
<b>CORB</b>	<i>Asymptotic correlation matrix of the parameter estimates.</i>
<b>COVB</b>	<i>Asymptotic covariance matrix of the parameter estimates.</i>
<b>FIT</b>	<i>Goodness-of-fit statistics.</i> The change in chi-square statistics with respect to a model with intercept terms only (or to a null model when INTERCEPT=EXCLUDE). The table contains the Pearson chi-square and the likelihood-ratio chi-square statistics. The statistics are computed based on the subpopulation classification specified on the SUBPOP subcommand or the default classification.
<b>HISTORY(n)</b>	<i>Iteration history.</i> The table contains log-likelihood function value and parameter estimates at every <i>n</i> th iteration beginning with the 0th iteration (the initial estimates). The default is to print every iteration ( <i>n</i> = 1). The last iteration is always printed if HISTORY is specified, regardless of the value of <i>n</i> .
<b>IC</b>	<i>Information criteria.</i> The Akaike Information Criterion (AIC) and the Schwarz Bayesian Information Criterion (BIC) are displayed.
<b>KERNEL</b>	<i>Kernel of the log-likelihood.</i> Displays the value of the kernel of the $-2$ log-likelihood. The default is to display the full $-2$ log-likelihood. Note that this keyword has no effect unless the MFI or LRT keywords are specified.
<b>LRT</b>	<i>Likelihood-ratio tests.</i> The table contains the likelihood-ratio test statistics for the model and model partial effects. If LRT is not specified, just the model test statistic is printed.
<b>PARAMETER</b>	<i>Parameter estimates.</i>
<b>SUMMARY</b>	<i>Model summary.</i> Cox and Snell's, Nagelkerke's, and McFadden's $R^2$ statistics.
<b>CPS</b>	<i>Case processing summary.</i> This table contains information about the specified categorical variables. Displayed by default.

<b>STEP</b>	<i>Step summary.</i> This table summarizes the effects entered or removed at each step in a stepwise method. Displayed by default if a stepwise method is specified. This keyword is ignored if no stepwise method is specified.
<b>MFI</b>	<i>Model fitting information.</i> This table compares the fitted and intercept-only or null models. Displayed by default.
<b>NON</b>	<i>No statistics are displayed.</i> This option overrides all other specifications on the PRINT subcommand.

## **SAVE Subcommand**

The SAVE subcommand puts casewise post-estimation statistics back into the active file.

- The new names must be valid variable names and not currently used in the active dataset.
- The rootname must be a valid variable name.
- The rootname can be followed by the number of predicted probabilities saved. The number is a positive integer. For example, if the integer is 5, then the first five predicted probabilities across all split files (if applicable) are saved. The default is 25.
- The new variables are saved into the active file in the order in which the keywords are specified on the subcommand.

<b>ACPROB(newname)</b>	<i>Estimated probability of classifying a factor/covariate pattern into the actual category.</i>
<b>ESTPROB(rootname:n)</b>	<i>Estimated probabilities of classifying a factor/covariate pattern into the response categories.</i> There are as many number of probabilities as the number of response categories. The predicted probabilities of the first <i>n</i> response categories will be saved. The default value for <i>n</i> is 25. To specify <i>n</i> without a rootname, enter a colon before the number.
<b>PCPROB(newname)</b>	<i>Estimated probability of classifying a factor/covariate pattern into the predicted category.</i> This probability is also the maximum of the estimated probabilities of the factor/covariate pattern.
<b>PREDCAT(newname)</b>	<i>The response category that has the maximum expected probability for a factor/covariate pattern.</i>

## **SCALE Subcommand**

The SCALE subcommand specifies the dispersion scaling value. Model estimation is not affected by this scaling value. Only the asymptotic covariance matrix of the parameter estimates is affected.

<b>N</b>	<i>A positive number corresponding to the amount of overdispersion or underdispersion.</i> The default scaling value is 1, which corresponds to no overdispersion or underdispersion.
<b>DEVIANCE</b>	<i>Estimates the scaling value by using the deviance function statistic.</i>
<b>PEARSON</b>	<i>Estimates the scaling value by using the Pearson chi-square statistic.</i>

## **SUBPOP Subcommand**

The SUBPOP subcommand allows you to define the subpopulation classification used in computing the goodness-of-fit statistics.

- A variable list is expected if the `SUBPOP` subcommand is specified. The variables in the list must be a subset of the combined list of factors and covariates specified on the command line.
- Variables specified or implied on the `MODEL` subcommand must be a subset of the variables specified or implied on the `SUBPOP` subcommand.
- If the `SUBPOP` subcommand is omitted, the default classification is based on all of the factors and the covariates specified.
- Missing values are deleted listwise on the subpopulation level.

### Example

```
NOMREG
movie BY gender date WITH age
/CRITERIA = CIN(95) DELTA(0) MXITER(100) MXSTEP(5) LCONVERGE(0)
          PCONVERGE(1.0E-6) SINGULAR(1.0E-8)
/MODEL = gender
/SUBPOP = gender date
/INTERCEPT = EXCLUDE .
```

- Although the model consists only of *gender*, the `SUBPOP` subcommand specifies that goodness-of-fit statistics should be computed based on both *gender* and *date*.

## TEST Subcommand

The `TEST` subcommand allows you to customize your hypothesis tests by directly specifying null hypotheses as linear combinations of parameters.

- `TEST` is offered only through syntax.
- Multiple `TEST` subcommands are allowed. Each is handled independently.
- The basic format for the `TEST` subcommand is an optional list of values enclosed in parentheses, an optional label in quotes, an effect name or the keyword `ALL`, and a list of values.
- The value list preceding the first effect or the keyword `ALL` are the constants to which the linear combinations are equated under the null hypotheses. If this value list is omitted, the constants are assumed to be all zeros.
- The label is a string with a maximum length of 255 characters (or 127 double-byte characters). Only one label per linear combination can be specified.
- When `ALL` is specified, only a list of values can follow. The number of values must equal the number of parameters (including the redundant ones) in the model.
- When effects are specified, only valid effects appearing or implied on the `MODEL` subcommand can be named. The number of values following an effect name must equal the number of parameters (including the redundant ones) corresponding to that effect. For example, if the effect `A*B` takes up six parameters, then exactly six values must follow `A*B`. To specify the coefficient for the intercept, use the keyword `INTERCEPT`. Only one value is expected to follow `INTERCEPT`.
- When multiple linear combinations are specified within the same `TEST` subcommand, use semicolons to separate each hypothesis.

- The linear combinations are first tested separately for each logit and then simultaneously tested for all of the logits.
- A number can be specified as a fraction with a positive denominator. For example,  $1/3$  or  $-1/3$  are valid, but  $1/-3$  is invalid.
- Effects appearing or implied on the MODEL subcommand but not specified on the TEST are assumed to take the value 0 for all of their parameters.

### **Example**

```
NOMREG
  movie BY gender date
  /CRITERIA = CIN(95) DELTA(0) MXITER(100) MXSTEP(5) LCONVERGE(0)
              PCONVERGE(1.0E-6) SINGULAR(1.0E-8)
  /INTERCEPT = EXCLUDE
  /PRINT = CELLPROB CLASSTABLE FIT CORB COVB HISTORY(1) PARAMETER
              SUMMARY LRT
  /TEST (0 0) = ALL 1 0 0 0;
              ALL 0 1 0 0 .
```

- TEST specifies two separate tests: one in which the coefficient corresponding to the first category for gender is tested for equality with zero, and one in which the coefficient corresponding to the second category for gender is tested for equality with zero.

# NONPAR CORR

NONPAR CORR is available in the Statistics Base option.

```
NONPAR CORR VARIABLES= varlist [WITH varlist] [/varlist...]

[/PRINT={TWOTAIL**} {SIG**} {SPEARMAN**}]
        {ONETAIL } {NOSIG} {KENDALL }
        {BOTH }

[/SAMPLE]

[/MISSING={PAIRWISE**} [INCLUDE]]
        {LISTWISE }

[/MATRIX=OUT({* })]
        {'savfile'|'dataset'}
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
NONPAR CORR
  VARIABLES=PRESTIGE SPPRES PAPRES16 DEGREE PADEG MADEG.
```

## Overview

NONPAR CORR computes two rank-order correlation coefficients, Spearman's rho and Kendall's tau-*b*, with their significance levels. You can obtain one or both coefficients. NONPAR CORR automatically computes the ranks and stores the cases in memory. Therefore, memory requirements are directly proportional to the number of cases that are being analyzed.

### Options

**Coefficients and Significance Levels.** By default, NONPAR CORR computes Spearman coefficients and displays the two-tailed significance level. You can request a one-tailed test, and you can display the significance level for each coefficient as an annotation by using the PRINT subcommand.

**Random Sampling.** You can request a random sample of cases by using the SAMPLE subcommand when there is not enough space to store all cases.

**Matrix Output.** You can write matrix materials to a matrix data file by using the MATRIX subcommand. The matrix materials include the number of cases that are used to compute each coefficient and the Spearman or Kendall coefficients for each variable. These materials can be read by other procedures.



### **Basic Specification**

The basic specification is `VARIABLES` and a list of numeric variables. By default, Spearman correlation coefficients are calculated.

### **Subcommand Order**

- `VARIABLES` must be specified first.
- The remaining subcommands can be used in any order.

### **Operations**

- `NONPAR CORR` produces one or more matrices of correlation coefficients. For each coefficient, `NONPAR CORR` displays the number of used cases and the significance level.
- The number of valid cases is always displayed. Depending on the specification on the `MISSING` subcommand, the number of valid cases can be displayed for each pair or in a single annotation.
- If all cases have a missing value for a given pair of variables, or if all cases have the same value for a variable, the coefficient cannot be computed. If a correlation cannot be computed, `NONPAR CORR` displays a decimal point.
- If both Spearman and Kendall coefficients are requested, and `MATRIX` is used to write matrix materials to a matrix data file, only Spearman's coefficient will be written with the matrix materials.

### **Limitations**

- A maximum of 25 variable lists is allowed.
- A maximum of 100 variables total per `NONPAR CORR` command is allowed.

## **Examples**

```
NONPAR CORR VARIABLES=PRESTIGE SPPRES PAPRES16 DEGREE PADEG MADEG.
```

- By default, Spearman correlation coefficients are calculated. The number of cases upon which the correlations are based and the two-tailed significance level are displayed for each correlation.

## **VARIABLES Subcommand**

`VARIABLES` specifies the variable list.

- All variables must be numeric.
- If keyword `WITH` is not used, `NONPAR CORR` displays the correlations of each variable with every other variable in the list.
- To obtain a rectangular matrix, specify two variable lists that are separated by keyword `WITH`. `NONPAR CORR` writes a rectangular matrix of variables in the first list correlated with variables in the second list.
- Keyword `WITH` cannot be used when the `MATRIX` subcommand is used.

- You can request more than one analysis. Use a slash to separate the specifications for each analysis.

### Example

```
NONPAR CORR
  VARIABLES = PRESTIGE SPPRES PAPRES16 WITH DEGREE PADEG MADEG.
```

- The three variables that are listed before `WITH` define the rows; the three variables that are listed after `WITH` define the columns of the correlation matrix.
- Spearman's rho is displayed by default.

### Example

```
NONPAR CORR VARIABLES=SPPRES PAPRES16 PRESTIGE
  /SATCITY WITH SATHOBBY SATFAM.
```

- `NONPAR CORR` produces two Correlations tables.
- By default, Spearman's rho is displayed.

## PRINT Subcommand

By default, `NONPAR CORR` displays Spearman correlation coefficients. The significance levels are displayed below the coefficients. The significance level is based on a two-tailed test. Use `PRINT` to change these defaults.

- The Spearman and Kendall coefficients are both based on ranks.

<b>SPEARMAN</b>	<i>Spearman's rho.</i> Only Spearman coefficients are displayed. This specification is the default.
<b>KENDALL</b>	<i>Kendall's tau-b.</i> Only Kendall coefficients are displayed.
<b>BOTH</b>	<i>Kendall and Spearman coefficients.</i> Both coefficients are displayed. If <code>MATRIX</code> is used to write the correlation matrix to a matrix data file, only Spearman coefficients are written with the matrix materials.
<b>SIG</b>	<i>Do not flag significant values.</i> <code>SIG</code> is the default.
<b>NOSIG</b>	<i>Flag significant values.</i> Values significant at the 0.05 level are flagged with a single asterisk; those that are significant at the 0.01 level are flagged with two asterisks.
<b>TWOTAIL</b>	<i>Two-tailed test of significance.</i> This test is appropriate when the direction of the relationship cannot be determined in advance, as is often the case in exploratory data analysis. This specification is the default.
<b>ONETAIL</b>	<i>One-tailed test of significance.</i> This test is appropriate when the direction of the relationship between a pair of variables can be specified in advance of the analysis.

## SAMPLE Subcommand

`NONPAR CORR` must store cases in memory to build matrices. `SAMPLE` selects a random sample of cases when computer resources are insufficient to store all cases. `SAMPLE` has no additional specifications.

## MISSING Subcommand

MISSING controls the treatment of missing values.

- PAIRWISE and LISTWISE are alternatives. You can specify INCLUDE with either PAIRWISE or LISTWISE.

<b>PAIRWISE</b>	<i>Exclude missing values pairwise.</i> Cases with a missing value for one or both variables for a specific correlation coefficient are excluded from the computation of that coefficient. This process allows the maximum available information to be used in every calculation. This process also results in a set of coefficients based on a varying number of cases. The number is displayed for each pair. This specification is the default.
<b>LISTWISE</b>	<i>Exclude missing values listwise.</i> Cases with a missing value for any variable that is named in a list are excluded from the computation of all coefficients in the Correlations table. The number of used cases is displayed in a single annotation. Each variable list on a command is evaluated separately. Thus, a case that is missing for one matrix might be used in another matrix. This option decreases the amount of required memory and significantly decreases computational time.
<b>INCLUDE</b>	<i>Include user-missing values.</i> User-missing values are treated as valid values.

## MATRIX Subcommand

MATRIX writes matrix materials to a matrix data file. The matrix materials always include the number of cases that are used to compute each coefficient, and the materials include either the Spearman or the Kendall correlation coefficient for each variable, whichever is requested. [For more information, see the topic Format of the Matrix Data File on p. 1358.](#)

- You cannot write both Spearman's and Kendall's coefficients to the same matrix data file. To obtain both Spearman's and Kendall's coefficients in matrix format, specify separate NONPAR CORR commands for each coefficient and define different matrix data files for each command.
- If PRINT=BOTH is in effect, NONPAR CORR displays a matrix in the listing file for both coefficients but writes only the Spearman coefficients to the matrix data file.
- NONPAR CORR cannot write matrix materials for rectangular matrices (variable lists containing keyword WITH). If more than one variable list is specified, only the last variable list that does not use keyword WITH is written to the matrix data file.
- The specification on MATRIX is keyword OUT and a quoted file specification or previously declared dataset name (DATASET DECLARE command), enclosed in parentheses.
- If you want to use a correlation matrix that is written by NONPAR CORR in another procedure, change the ROWTYPE\_ value RHO or TAUB to CORR by using the RECODE command.
- Any documents that are contained in the active dataset are not transferred to the matrix file.

**OUT ('savfile']dataset')** *Write a matrix data file or dataset.* Specify either a filename, a previously declared dataset name, or an asterisk, enclosed in parentheses. Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. If you specify an asterisk (\*), the matrix data file replaces the active dataset.

**Multiple nonparametric correlation tables**

```
NONPAR CORR
  VARIABLES=PRESTIGE SPPRES PAPRES16 DEGREE PADEG MADEG
  /PRESTIGE TO DEGREE /PRESTIGE WITH DEGREE
  /MATRIX OUT('/data/npmat.sav').
```

- Only the matrix for *PRESTIGE* to *DEGREE* is written to the matrix data file because it is the last variable list that does not use keyword *WITH*.

**Format of the Matrix Data File**

- The matrix data file has two special variables that are created by the program: *ROWTYPE\_* and *VARNAME\_*.
- *ROWTYPE\_* is a short string variable with values *N* and *RHO* for Spearman's correlation coefficient. If you specify Kendall's coefficient, the values are *N* and *TAUB*.
- *VARNAME\_* is a short string variable whose values are the names of the variables that are used to form the correlation matrix. When *ROWTYPE\_* is *RHO* (or *TAUB*), *VARNAME\_* gives the variable that is associated with that row of the correlation matrix.
- The remaining variables in the file are the variables that are used to form the correlation matrix.

**Split Files**

- When split-file processing is in effect, the first variables in the matrix data file are the split variables, followed by *ROWTYPE\_*, *VARNAME\_*, and the variables that are used to form the correlation matrix.
- A full set of matrix materials is written for each split-file group that is defined by the split variables.
- A split variable cannot have the same name as any other variable that is written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split file must be in effect when that matrix is read by a procedure.

**Missing Values**

- With *PAIRWISE* treatment of missing values (the default), the matrix of *NS* that is used to compute each coefficient is included with the matrix materials.
- With *LISTWISE* or *INCLUDE* treatments, a single *N* that is used to calculate all coefficients is included with the matrix materials.

**Examples****Writing results to a matrix data file**

```
GET FILE='/data/GSS80.sav'
  /KEEP PRESTIGE SPPRES PAPRES16 DEGREE PADEG MADEG.
NONPAR CORR VARIABLES=PRESTIGE TO MADEG
  /MATRIX OUT('/data/npmat.sav').
```

- NONPAR CORR reads data from file *GSS80.sav* and writes one set of correlation matrix materials to the file *npmat.sav*.
- The active dataset is still *GSS80.sav*. Subsequent commands are executed on file *GSS80.sav*.

***Replacing the active dataset with matrix results***

```
GET FILE='/data/GSS80.sav'  
  /KEEP PRESTIGE SPPRES PAPRES16 DEGREE PADEG MADEG.  
NONPAR CORR VARIABLES=PRESTIGE TO MADEG  
  /MATRIX OUT(*).  
LIST.  
DISPLAY DICTIONARY.
```

- NONPAR CORR writes the same matrix as in the example above. However, the matrix data file replaces the active dataset. The LIST and DISPLAY commands are executed on the matrix file (not on the original active dataset *GSS80.sav*).

# NPTESTS

NPTESTS is available in the Statistics Base option.

```
NPTESTS
[/MISSING [SCOPE={ANALYSIS**}] [USERMISSING={EXCLUDE**}]]
      {LISTWISE } {INCLUDE }
[/CRITERIA ALPHA={0.05**}] CILEVEL={95** }
      {value } {value}

[/ONESAMPLE TEST(fieldlist) ]
  [CHISQUARE(
    [EXPECTED={EQUAL** }]]
    {CUSTOM(FREQUENCIES=valuelist
            CATEGORIES=valuelist) }
  )]
  [BINOMIAL ( [TESTVALUE={0.5**}]
              {value}
    [SUCCESSCATEGORICAL={FIRST** }]]
    {LIST(valuelist) }
    [SUCCESSCONTINUOUS=CUTPOINT({MIDPOINT**})]
    {value }
    [CLOPPERPEARSON] [JEFFREYS] [LIKELIHOOD]
  )]
  [KOLMOGOROV_SMIRNOV(
    [NORMAL={SAMPLE** }]]
    {CUSTOM(MEAN=value
            SD=value )}
    [UNIFORM={SAMPLE** }]]
    {CUSTOM(MIN=value
            MAX=value )}
    [EXPONENTIAL={SAMPLE** }]]
    {CUSTOM(MEAN=value )}
    [POISSON={SAMPLE** }]]
    {CUSTOM(MEAN=value )}
  )]
  [RUNS ( [GROUPCATEGORICAL={SAMPLE** }]]
    {LIST(valuelist) }
    [GROUPCONTINUOUS=CUTPOINT({SAMPLEMEDIAN**})]
    {SAMPLEMEAN }
    {value }
  )]
  [WILCOXON(TESTVALUE=value)]

[/INDEPENDENT TEST(fieldlist) ]
  GROUP(fieldname)

  [MANN_WHITNEY] [WALD_WOLFOWITZ] [KOLMOGOROV_SMIRNOV]
  [HODGES_LEHMAN]
  [MOSES ( [TRIMOUTLIERS={SAMPLE**}])]
    {integer }

  [KRUSKAL_WALLIS ( [COMPARE={NONE }]]]
    {PAIRWISE**}
    {STEPWISE }

  [MEDIAN ( [TESTVALUE={SAMPLE**}] [COMPARE={NONE }]]]
    {value } {PAIRWISE**}
    {STEPWISE }

  [JONCKHEERE_TERPSTRA (
    [ORDER={ASCENDING** }]] [COMPARE={NONE }]]
    {DESCENDING } {PAIRWISE**}
    {STEPWISE }
  )]
```

```

[/RELATED TEST(fieldlist) ]
  [MCNEMAR( [SUCCESS={FIRST** }])]
    {LIST(valuelist)}

  [WILCOXON] [SIGN] [MARGINAL_HOMOGENEITY]
  [HODGES_LEHMAN]

  [COCHRAN(
    [SUCCESS={FIRST } ]
    {LIST(valuelist)}
    [COMPARE={NONE } ]
    {PAIRWISE**}
    {STEPWISE }
  )]

  [FRIEDMAN( [COMPARE={NONE }])]
    {PAIRWISE**}
    {STEPWISE }

  [KENDALL( [COMPARE={NONE }])]
    {PAIRWISE**}
    {STEPWISE }

```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 18

- Command introduced.

### **Example**

```

NPTESTS
  /ONESAMPLE TEST(field1 TO field10).

```

## **Overview**

NPTESTS is a collection of nonparametric tests. These tests make minimal assumptions about the underlying distribution of the data.

The tests that are available in NPTESTS can be grouped into three broad categories based on how the data are organized.

- A one-sample test analyzes one field.
- A test for related samples compares two or more fields for the same set of cases.
- An independent-samples test analyzes one field that is grouped by categories of another field.

### **Options**

**Automatic test selection.** If a test subcommand (ONESAMPLE, INDEPENDENT, or RELATED) is specified without any test specifications, then NPTESTS automatically chooses the “best” tests to perform for each field specified on the subcommand. See the individual subcommands for details.

**Multiple comparisons.** In addition to the “omnibus” tests for  $k$  independent or related samples, the samples can also be tested pairwise, either by all pairwise comparisons or through a stepwise stepdown procedure.

**Missing value handling.** Records with missing values can be handled listwise or analysis by analysis. User-missing values on categorical fields can be treated as valid or excluded from analysis.

### ***Basic Specification***

The basic specification is a single test subcommand (ONESAMPLE, INDEPENDENT, or RELATED) and a TEST keyword with a list of fields to be tested.

### ***Syntax Rules***

- At least one test subcommand must be specified; all other subcommands are optional.
- Subcommands may be specified in any order.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not allowed.
- Any split field defined on the SPLIT FILE command cannot be used on this command.
- Any field specified on the WEIGHT command cannot be used on this command.

### ***Operations***

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

## ***MISSING Subcommand***

The MISSING subcommand controls the treatment of records with missing values.

```
[/MISSING [SCOPE={ANALYSIS**}] [USERMISSING={EXCLUDE**}]]
           {LISTWISE }           {INCLUDE }
```



**SCOPE Keyword**

The SCOPE keyword specifies how to determine the case basis for tests.

- ANALYSIS**      *Exclude records with missing values on a test-by-test basis.* Records with missing values for a field that is used for a specific test are omitted from that test. On subcommands that specify several tests, each test is evaluated separately. This setting is the default.
- LISTWISE**      *Exclude records with missing values listwise.* Records with missing values for any field that is named on any subcommand are excluded from all analyses.

**USERMISSING Keyword**

The USERMISSING keyword is used to control whether user-missing values for categorical fields are treated as valid values. By default, user-missing values for categorical fields are treated as invalid.

- User-missing values for continuous fields are always treated as invalid.
- System-missing values for any fields are always treated as invalid.

- EXCLUDE**      *User-missing values for categorical fields are treated as invalid.* This is the default.
- INCLUDE**      *User-missing values for categorical fields are treated as valid values.*

**CRITERIA Subcommand**

The CRITERIA subcommand controls the significance level used for tests and the confidence level used for intervals.

```
[ /CRITERIA ALPHA={0.05**}] CILEVEL={95** }
                        {value }           {value}
```

**ALPHA Keyword**

The ALPHA keyword specifies the significance level for all tests specified on the NPTESTS command.

- Specify a numeric value between 0 and 1. 0.05 is the default.

**CILEVEL Keyword**

The CILEVEL keyword specifies the confidence level for all confidence intervals produced by the NPTESTS command.

- Specify a numeric value between 0 and 100. 95 is the default.

## **ONESAMPLE Subcommand**

The `ONESAMPLE` subcommand produces one-sample nonparametric tests. The `TEST` keyword is required, all other keywords are optional. If `ONESAMPLE` is specified with none of the optional keywords, the following tests are performed automatically:

- Categorical fields with two values are tested using a binomial test.
- Categorical fields with more than two values are tested using a chi-square test with equal frequencies on category values found in the sample.
- Continuous fields are tested using a Kolmogorov-Smirnov test against a normal distribution with the sample mean and standard deviation.

### **TEST Keyword**

The `TEST` keyword lists the fields that you want to test.

- Specify one or more fields. Note that certain tests are not applicable to fields of a particular measurement level; for example, the chi-square test is only performed for categorical fields. `NPTESTS` automatically determines which tests are applicable to which fields. See the individual keyword descriptions for details.

### **CHISQUARE Keyword**

```
[CHISQUARE ( [EXPECTED={EQUAL**
                {CUSTOM (FREQUENCIES=valuelist
                        CATEGORIES=valuelist) } ] ] ]
```

The `CHISQUARE` keyword produces a one-sample test that computes a chi-square statistic based on the differences between the observed and expected frequencies of categories of a field.

- A separate chi-square test is performed for each and every categorical field specified on the `TEST` keyword.
- The test specifications given on the `CHISQUARE` keyword apply to all chi-square tests performed.
- If `CHISQUARE` is specified without any keywords, equal frequencies are expected in each category.

#### **EXPECTED = EQUAL|CUSTOM(FREQUENCIES=valuelist CATEGORIES=valuelist)**

*Expected frequencies.* The `EXPECTED` keyword defines how expected frequencies are derived. The default is `EQUAL`.

`EQUAL` produces equal frequencies among all categories in the sample. This is the default when `CHISQUARE` is specified without any other keywords..

`CUSTOM` allows you to specify unequal frequencies for a specified list of categories.

On the `CATEGORIES` keyword, specify a list of string or numeric values. The values in the list do not need to be present in the sample.

On the `FREQUENCIES` keyword, specify a value greater than 0 for each category, and in the same order as the categories, on the `CATEGORIES` keyword. Custom frequencies are treated as ratios so that, for example, `FREQUENCIES=1 2 3` is equivalent to `FREQUENCIES=10 20 30`, and both specify that 1/6 of the records are expected to fall into the first category on the `CATEGORIES` keyword, 1/3 into the second, and 1/2 into the third.

When `CUSTOM` is specified, the number of expected frequencies must match the number of category values; otherwise the test is not performed for that field.

### ***BINOMIAL Keyword***

```
[BINOMIAL ( [TESTVALUE={0.5**} ]
            {value}
            [SUCCESSCATEGORICAL={FIRST**           } ]
            {LIST(valuelist)                       } ]
            [SUCCESSCONTINUOUS=CUTPOINT({MIDPOINT**}) ]
            {value }
            [CLOPPERPEARSON] [JEFFREYS] [LIKELIHOOD]
        ) ]
```

The `BINOMIAL` keyword produces a one-sample test of whether the observed distribution of a dichotomous field is the same as what is expected from a specified binomial distribution. In addition, you can request confidence intervals.

- A separate binomial test is performed for each and every field specified on the `TEST` keyword.
- The test specifications given on the `BINOMIAL` keyword apply to all binomial tests performed.
- If `BINOMIAL` is specified without any keywords, each categorical field is assumed to have only two values and each continuous field is dichotomized using the average of the minimum and maximum as a cut point. The distribution of each named field is compared to a binomial distribution with  $p$  (the proportion of cases expected in the first category) equal to 0.5.

#### **TESTVALUE**

*Hypothesized proportion.* The `TESTVALUE` keyword specifies the expected proportion of cases in the first category. Specify a value greater than 0 and less than 1. The default is 0.5.

#### **SUCCESSCATEGORICAL=FIRST|LIST(valuelist)**

The `SUCCESSCATEGORICAL` keyword specifies how “success”, the data value(s) tested against the test value, is defined for categorical fields.

`FIRST` performs the binomial test using the first value found in the sample to define “success”. This option is only applicable to nominal or ordinal fields with only two values; all other categorical fields specified on a `ONESAMPLE` subcommand where `FIRST` is used will not be tested. This is the default.

`LIST` performs the binomial test using the specified list of values to define “success”. Specify a list of string or numeric values. The values in the list do not need to be present in the sample.

#### **SUCCESSCONTINUOUS=CUTPOINT(MIDPOINT|value)**

*Define success for continuous fields.* The `SUCCESSCONTINUOUS` keyword specifies how “success”, the data value(s) tested against the test value, is defined for continuous fields.

`CUTPOINT` defines values that are equal to or less than the cut point as “success”.

`MIDPOINT` sets the cut point at the average of the minimum and maximum values. Alternatively, specify a value for the cut point. The default is `MIDPOINT`.

#### **CLOPPERPEARSON**

*Exact interval based on the cumulative binomial distribution.*

### **JEFFREYS**

*Bayesian interval based on the posterior distribution of  $p$  using the Jeffreys prior.*

### **LIKELIHOOD**

*Interval based on the likelihood function for  $p$ .*

### **KOLMOGOROV\_SMIRNOV Keyword**

```
[KOLMOGOROV_SMIRNOV (
  [NORMAL={SAMPLE**          }]
    {CUSTOM(MEAN=value
             SD=value   )}
  [UNIFORM={SAMPLE
            {CUSTOM(MIN=value
                    MAX=value )}
  [EXPONENTIAL={SAMPLE          }]
    {CUSTOM(MEAN=value )}
  [POISSON={SAMPLE
           {CUSTOM(MEAN=value )}
  )]
```

The `KOLMOGOROV_SMIRNOV` keyword produces a one-sample test of whether the sample cumulative distribution function for a field is homogenous with a uniform, normal, Poisson, or exponential distribution.

- A separate Kolmogorov-Smirnov test is performed for each and every continuous field specified on the `TEST` keyword.
- The test specifications given on the `KOLMOGOROV_SMIRNOV` keyword apply to all Kolmogorov-Smirnov tests performed.
- If `KOLMOGOROV_SMIRNOV` is specified without any keywords, each field is tested against a normal distribution using its sample mean and sample standard deviation.

#### **NORMAL(SAMPLE|CUSTOM(MEAN=value SD=value))**

*Normal distribution.* `SAMPLE` uses the observed mean and standard deviation, `CUSTOM` allows you to specify values.

#### **UNIFORM(SAMPLE|CUSTOM(MIN=value MAX=value))**

*Uniform distribution.* `SAMPLE` uses the observed minimum and maximum, `CUSTOM` allows you to specify values.

#### **POISSON(SAMPLE|CUSTOM(MEAN=value))**

*Poisson distribution.* `SAMPLE` uses the observed mean, `CUSTOM` allows you to specify a value.

#### **EXPONENTIAL(SAMPLE|CUSTOM(MEAN=value))**

*Exponential distribution.* `SAMPLE` uses the observed mean, `CUSTOM` allows you to specify a value.

### **RUNS Keyword**

```
[RUNS ( [GROUPCATEGORICAL={SAMPLE**          }]
        {LIST(valuelist)          }
  [GROUPCONTINUOUS=CUTPOINT ( {SAMPLEMEDIAN** } )]
        {SAMPLEMEAN              }
        {value                    }
  )]
```

The `RUNS` keyword produces a one-sample test of whether the sequence of values of a dichotomized field is random.

- A separate runs test is performed for each and every field specified on the `TEST` keyword.
- The test specifications given on the `RUNS` keyword apply to all runs tests performed.
- If `RUNS` is specified without any keywords, each categorical field is assumed to have only two values and each continuous field is dichotomized using the sample median as a cut point.

#### **GROUPCATEGORICAL=SAMPLE|LIST(value list)**

*Determine groups for categorical fields.* `SAMPLE` is the default.

`SAMPLE` performs the runs test using the values found in the sample to define the groups. This option is only applicable to nominal or ordinal fields with only two values; all other categorical fields specified on a `ONESAMPLE` subcommand where `SAMPLE` is used will not be tested.

`LIST` performs the runs test using the specified list of values to define one of the groups. All other values in the sample define the other group. The values in the list do not all need to be present in the sample, but at least one record must be in each group.

#### **GROUPCONTINUOUS=CUTPOINT(SAMPLEMEDIAN | SAMPLEMEAN | value)**

*Determine groups for continuous fields.* `CUTPOINT` defines values that are equal to or less than the cut point as the first group; all other values define the other group. `SAMPLEMEDIAN` sets the cut point at the sample median. `SAMPLEMEAN` sets the cut point at the sample mean. Alternatively, specify a value for the cut point. The default is `SAMPLEMEDIAN`.

### ***WILCOXON Keyword***

[`WILCOXON (TESTVALUE=value)` ]

The `WILCOXON` keyword produces a one sample test of median value of a field.

- A separate Wilcoxon test is performed for each and every continuous field specified on the `TEST` keyword.
- The test specifications given on the `WILCOXON` keyword apply to all Wilcoxon tests performed.
- The `TESTVALUE` keyword is required.

#### **TESTVALUE=value**

*Hypothesized median.* The Wilcoxon test is performed using the specified value. The `TESTVALUE` keyword is required. There is no default.

## ***INDEPENDENT Subcommand***

The `INDEPENDENT` subcommand produces independent samples nonparametric tests. The `TEST` and `GROUP` keywords are required, all other keywords are optional. If `INDEPENDENT` is specified with none of the optional keywords, the following tests are performed automatically:

- If the grouping field has only two values, continuous fields are tested using a Mann-Whitney test.

- If the grouping field has more than two values, continuous fields are tested using a Kruskal-Wallis test.
- If there are no continuous fields on the TEST keyword, no tests are performed.

### **TEST Keyword**

The TEST keyword lists the fields that you want to test.

- Specify one or more continuous fields. Independent samples tests are only applicable to continuous fields.

### **GROUP Keyword**

The GROUP keyword specifies the field used to determine groups. The test is performed using the values found in the sample to define the groups. Note that certain tests are not applicable when there are more than two groups. NPTESTS automatically determines which tests are applicable. See the individual keyword descriptions for details.

- Specify a categorical field. A grouping field that does not have nominal or ordinal measurement level will cause an error.
- GROUP applies to all fields that are specified on the TEST keyword.

### **MANN\_WHITNEY Keyword**

The MANN\_WHITNEY keyword produces an independent samples test of whether two samples are from the same population. The test statistic uses the rank of each case to test whether the groups are drawn from the same population.

- A separate test is performed for each and every continuous field specified on the TEST keyword.
- The first value in ascending order of the grouping field defines the first group and the second value defines the second. If the grouping field has more than two values, the test is not performed.

### **WALD\_WOLFOWITZ Keyword**

The WALD\_WOLFOWITZ keyword produces an independent samples test whether the distribution of a field is the same in two independent samples. A runs test is performed with group membership as the criterion.

- A separate test is performed for each and every continuous field specified on the TEST keyword.
- The first value in ascending order of the grouping field defines the first group and the second value defines the second. If the grouping field has more than two values, the test is not performed.

**KOLMOGOROV\_SMIRNOV Keyword**

The KOLMOGOROV\_SMIRNOV keyword produces an independent samples test of whether the distribution of a field is the same in two independent samples that are defined by a grouping field. The test is sensitive to any difference in median, dispersion, skewness, and so forth, between the two distributions.

- A separate test is performed for each and every continuous field specified on the TEST keyword.
- The first value in ascending order of the grouping field defines the first group and the second value defines the second. If the grouping field has more than two values, the test is not performed.

**MOSES Keyword**

```
[MOSES ([TRIMOUTLIERS={SAMPLE**}])]
      {integer }
```

The MOSES keyword produces an independent samples test of whether the range of a continuous field is the same in a control group and a comparison group.

- A separate test is performed for each and every continuous field specified on the TEST keyword.
- The test specifications given on the MOSES keyword apply to all Moses tests performed.
- The first value in ascending order of the grouping field defines the control group and the second defines the comparison group. If the grouping field has more than two values, the test is not performed.

**TRIMOUTLIERS=SAMPLE|integer**

*Records to trim from each end of the range of the control group to remove outliers. SAMPLE trims 5% of the records from each end of the range of the control group to remove outliers, with a minimum of 1 record removed from each end. Specifying an integer causes that exact number of records to be trimmed. SAMPLE is the default.*

**HODGES\_LEHMAN Keyword**

The HODGES\_LEHMAN keyword produces an independent samples estimate and confidence interval for the difference in the medians of two groups.

- A separate confidence interval is computed for each and every continuous field specified on the TEST keyword.
- The first value in ascending order of the grouping field defines the first group and the second value defines the second. If the grouping field has more than two values, the confidence interval is not produced.

**KRUSKAL\_WALLIS Keyword**

```
[KRUSKAL_WALLIS ([COMPARE={PAIRWISE**}])]
                  {STEPWISE  }
                  {NONE      }
```

The `KRUSKAL_WALLIS` keyword produces an independent samples test of whether  $k$  independent samples that are defined by a grouping field are from the same population.

- A separate test is performed for each and every continuous field specified on the `TEST` keyword.
- The test specifications given on the `KRUSKAL_WALLIS` keyword apply to all tests performed.

**COMPARE=PAIRWISE|STEPWISE|NONE**

*Multiple comparisons.* The `COMPARE` keyword controls how and whether multiple comparisons should be performed. `PAIRWISE` produces all pairwise multiple comparisons. `STEPWISE` produces stepwise stepdown comparisons. `NONE` turns off multiple comparisons. By default, all pairwise comparisons are produced.

**MEDIAN Keyword**

```
[MEDIAN ( [TESTVALUE={SAMPLE**} ] [COMPARE={PAIRWISE**}] ) ]
           {value      }           {STEPWISE  }
                                   {NONE      }
```

The `MEDIAN` keyword produces an independent samples test of whether  $k$  independent samples are drawn from populations with the same median.

- A separate test is performed for each and every continuous field specified on the `TEST` keyword.
- The test specifications given on the `MEDIAN` keyword apply to all tests performed.

**TESTVALUE=SAMPLE|value**

*Median value to use in the test.* `SAMPLE` causes the test to be performed using the sample median, calculated from all records that are included in the test. Alternatively, you can specify an exact value. `SAMPLE` is the default.

**COMPARE=PAIRWISE|STEPWISE|NONE**

*Multiple comparisons.* The `COMPARE` keyword controls how and whether multiple comparisons should be performed. `PAIRWISE` produces all pairwise multiple comparisons. `STEPWISE` produces stepwise stepdown comparisons. `NONE` turns off multiple comparisons. By default, all pairwise comparisons are produced.

**JONCKHEERE\_TERPSTRA Keyword**

```
[JONCKHEERE_TERPSTRA (
  [ORDER={ASCENDING** } ] [COMPARE={PAIRWISE**}]
          {DESCENDING  }           {STEPWISE  }
                                   {NONE      }
) ]
```

The `JONCKHEERE-TERPSTRA` keyword produces an independent samples test of whether  $k$  independent samples that are defined by a grouping field are from the same population. This test is particularly powerful when the  $k$  populations have a natural ordering.



- A separate test is performed for each and every continuous field specified on the `TEST` keyword.
- The test specifications given on the `JONCKHEERE_TERPSTRA` keyword apply to all tests performed.

#### **ORDER=ASCENDING|DESCENDING**

*Order of the  $k$  populations.* `ASCENDING` stipulates an alternative hypothesis that the location parameter of the first group is not equal to the second, which in turn is not equal to the third, and so on. This is the default.

`DESCENDING` stipulates an alternative hypothesis that the location parameter of the last group is not equal to the second-to-last, which in turn is not equal to the third-to-last, and so on.

#### **COMPARE=PAIRWISE|STEPWISE|NONE**

*Multiple comparisons.* The `COMPARE` keyword controls how and whether multiple comparisons should be performed. `PAIRWISE` produces all pairwise multiple comparisons. `STEPWISE` produces stepwise stepdown comparisons. `NONE` turns off multiple comparisons. By default, all pairwise comparisons are produced.

## ***RELATED Subcommand***

The `RELATED` subcommand produces related-samples nonparametric tests. The `TEST` keyword is required, all other keywords are optional. If `RELATED` is specified with none of the optional keywords, the following tests are performed automatically:

- If two categorical fields are specified and they have only two values, McNemar's test is performed.
- If two continuous fields are specified, a Wilcoxon Matched-Pairs Signed-Rank test is performed.
- If more than two categorical fields are specified and they each have only two values, Cochran's test is performed.
- If more than two continuous fields are specified, Friedman's Two-Way Analysis of Variance by Ranks is performed.
- Otherwise, no tests are performed.

When fields of differing measurement level are specified, they are first separated by measurement level and then the appropriate test is applied to each group. For example, if you specify:

```
NPTESTS
  /RELATED TEST(c1 n1 o1 c2 o2 n2 o3 c3 o4) .
```

Where  $x_1$ ,  $x_2$ , and  $x_3$  are continuous fields,  $a_1$  and  $a_2$  are nominal fields, and  $o_1$ ,  $o_2$ ,  $o_3$ , and  $o_4$  are ordinal fields, then Friedman's test is applied to the continuous fields, McNemar's test is applied to the nominal fields, and no test is applied to the ordinal fields.

**TEST Keyword**

The `TEST` keyword lists the fields that you want to test.

- Specify one or more fields. Note that certain tests are not applicable to fields of a particular measurement level; for example, the McNemar test is only performed for categorical fields. `NPTESTS` automatically determines which tests are applicable to which fields. See the individual keyword descriptions for details.

**MCNEMAR Keyword**

```
[MCNEMAR ( [SUCCESS={FIRST**          }])]
           {LIST(valuelist)}
```

The `MCNEMAR` keyword produces a related samples test of whether combinations of values between two dichotomous fields are equally likely.

- A separate test is performed for all of the nominal fields and for all of the ordinal fields specified on the `TEST` keyword.
- The test specifications given on the `MCNEMAR` keyword apply to all tests performed.
- If any test has more than two fields, or any field has more than two values, the test is not performed.

**Example**

```
NPTESTS /RELATED TEST(n1 n2 n3 o1 o2) MCNEMAR.
```

If  $n1$ ,  $n2$ ,  $n3$  are nominal fields and  $o1$  and  $o2$  are ordinal fields, then no test is performed for the nominal fields and a test is performed for the ordinal fields so long as  $o1$  and  $o2$  each have only two values.

**SUCCESS=FIRST|LIST(valuelist)**

The `SUCCESS` keyword specifies how “success” is defined for categorical fields.

`FIRST` performs the test using the first value found in the sample to define “success”. This option is only applicable to nominal or ordinal fields with only two values; all other categorical fields specified on a `RELATED` subcommand where `FIRST` is used will not be tested. This is the default.

`LIST` performs the test using the specified list of values to define “success”. Specify a list of string or numeric values. The values in the list do not need to be present in the sample.

**WILCOXON Keyword**

The `WILCOXON` keyword produces a related samples test of whether the distribution of two paired fields in two related samples is the same. The test takes into account the magnitude of the differences between two paired fields.

- A single test is performed for all of the continuous fields specified on the `TEST` keyword.
- If there are more than two continuous fields, the test is not performed.

***SIGN Keyword***

The `SIGN` keyword produces a related samples test of whether the distribution of two paired fields in a two-related-samples test is the same.

- A single test is performed for all of the continuous fields specified on the `TEST` keyword.
- If there are more than two continuous fields, the test is not performed.

***MARGINAL\_HOMOGENEITY Keyword***

The `MARGINAL_HOMOGENEITY` keyword produces a related samples test of whether combinations of values between two paired ordinal fields are equally likely. The marginal homogeneity test is typically used in repeated measures situations. This test is an extension of the McNemar test from binary response to multinomial response.

- A single test is performed for all of the ordinal fields specified on the `TEST` keyword.
- If there are more than two ordinal fields, the test is not performed.

***HODGES\_LEHMAN Keyword***

The `HODGES_LEHMAN` keyword produces a related samples estimate and confidence interval for the median difference between two paired continuous fields.

- A single confidence interval is computed for all of the continuous fields specified on the `TEST` keyword.
- If there are more than two continuous fields, no confidence intervals are created.

***COCHRAN Keyword***

```
[COCHRAN (
  [SUCCESS={FIRST**      } ]
           {LIST(valuelist)}
  [COMPARE={PAIRWISE**} ]
           {STEPWISE  }
           {NONE      }
)]
```

The `COCHRAN` keyword produces a related samples test of whether the distribution of values is the same for  $k$  related dichotomous fields.

- A separate test is performed for all of the nominal fields and for all of the ordinal fields specified on the `TEST` keyword.
- The test specifications given on the `COCHRAN` keyword apply to all tests performed.
- If any field has more than two values, the test is not performed.

***Example***

```
NPTESTS /RELATED TEST(n1 n2 n3 o1 o2) COCHRAN.
```

If  $n1$ ,  $n2$ ,  $n3$  are nominal fields and  $o1$  and  $o2$  are ordinal fields, then a test is performed for the nominal fields and a test is performed for the ordinal fields so long as each field has only two values.

**SUCCESS=FIRST|LIST(valuelist)**

The SUCCESS keyword specifies how “success” is defined for categorical fields.

FIRST performs the test using the first value found in the sample to define “success”. This option is only applicable to nominal or ordinal fields with only two values; all other categorical fields specified on a RELATED subcommand where FIRST is used will not be tested. This is the default.

LIST performs the test using the specified list of values to define “success”. Specify a list of string or numeric values. The values in the list do not need to be present in the sample.

**COMPARE=PAIRWISE|STEPWISE|NONE**

*Multiple comparisons.* The COMPARE keyword controls how and whether multiple comparisons should be performed. PAIRWISE produces all pairwise multiple comparisons. STEPWISE produces stepwise stepdown comparisons. NONE turns off multiple comparisons. By default, all pairwise comparisons are produced.

**FRIEDMAN Keyword**

```
[FRIEDMAN ( [COMPARE={ PAIRWISE** } ] ) ]
                {STEPWISE  }
                {NONE      }
```

The FRIEDMAN subcommand produces a related samples test of whether  $k$  related samples have been drawn from the same population.

A single test is performed for all of the continuous fields specified on the TEST keyword.

**COMPARE=PAIRWISE|STEPWISE|NONE**

*Multiple comparisons.* The COMPARE keyword controls how and whether multiple comparisons should be performed. PAIRWISE produces all pairwise multiple comparisons. STEPWISE produces stepwise stepdown comparisons. NONE turns off multiple comparisons. By default, all pairwise comparisons are produced.

**KENDALL Keyword**

```
[KENDALL ( [COMPARE={ PAIRWISE** } ] ) ]
                {STEPWISE  }
                {NONE      }
```

The KENDALL keyword produces a related samples test of whether  $k$  related samples are from the same population. Kendall's  $W$  is a measure of agreement among judges or raters, where each case is one judge's rating of several items (fields).

A single test is performed for all of the continuous fields specified on the `TEST` keyword.

**COMPARE=PAIRWISE|STEPWISE|NONE**

*Multiple comparisons.* The `COMPARE` keyword controls how and whether multiple comparisons should be performed. `PAIRWISE` produces all pairwise multiple comparisons. `STEPWISE` produces stepwise stepdown comparisons. `NONE` turns off multiple comparisons. By default, all pairwise comparisons are produced.

# NPART TESTS

NPART TESTS is available in the Statistics Base option.

```
NPART TESTS [CHISQUARE=varlist[(lo,hi)]] [/EXPECTED={EQUAL          }]
                                                    {f1,f2,...fn}

[/K-S({UNIFORM [min,max]   })=varlist]
      {NORMAL [mean,stddev]}
      {POISSON [mean]      }
      {EXPONENTIAL [mean]  }

[/RUNS({MEAN   })=varlist]
      {MEDIAN}
      {MODE   }
      {value  }

[/BINOMIAL[({.5})]=varlist[({value1,value2})]]
      { p}          {value  }

[/MCNEMAR=varlist [WITH varlist [(PAIRED)]]]

[/SIGN=varlist [WITH varlist [(PAIRED)]]]

[/WILCOXON=varlist [WITH varlist [(PAIRED)]]]
  | /MH=varlist [WITH varlist [(PAIRED)]]]††

[/COCHRAN=varlist]

[/FRIEDMAN=varlist]

[/KENDALL=varlist]

[/M-W=varlist BY var (value1,value2)]

[/K-S=varlist BY var (value1,value2)]

[/W-W=varlist BY var (value1,value2)]

[/MOSES[(n)]=varlist BY var (value1,value2)]

[/K-W=varlist BY var (value1,value2)]

[/J-T=varlist BY var (value1, value2)]††

[/MEDIAN[(value)]=varlist BY var (value1,value2)]

[/MISSING={ANALYSIS**} [INCLUDE]]
          {LISTWISE  }

[/SAMPLE]

[/STATISTICS={DESCRIPTIVES} [QUARTILES] [ALL]]

  [ /METHOD={MC [CIN({99.0 })] [SAMPLES({10000})] ]}††
    {value}          {value}
    {EXACT [TIMER({5  })]
      {value}
    }
```

\*\*Default if the subcommand is omitted.

††Available only if the Exact Tests option is installed (available only on Windows operating systems).

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 17.0

- Increased limits on number of variables allowed in the analysis.

### **Example**

```
NPAR TESTS K-S (UNIFORM) =V1 /K-S (NORMAL, 0, 1) =V2 .
```

## **Overview**

NPAR TESTS is a collection of nonparametric tests. These tests make minimal assumptions about the underlying distribution of the data. (Siegel and Castellan, 1988) In addition to the nonparametric tests that are available in NPAR TESTS, the  $k$ -sample chi-square and Fisher's exact test are available in procedure CROSSTABS.

The tests that are available in NPAR TESTS can be grouped into three broad categories based on how the data are organized: one-sample tests, related-samples tests, and independent-samples tests. A one-sample test analyzes one variable. A test for related samples compares two or more variables for the same set of cases. An independent-samples test analyzes one variable that is grouped by categories of another variable.

The one-sample tests that are available in procedure NPAR TESTS are:

- BINOMIAL
- CHISQUARE
- K-S (Kolmogorov-Smirnov)
- RUNS

Tests for two related samples are:

- MCNEMAR
- SIGN
- WILCOXON

Tests for  $k$  related samples are:

- COCHRAN
- FRIEDMAN
- KENDALL

Tests for two independent samples are:

- M-W (Mann-Whitney)
- K-S (Kolmogorov-Smirnov)

- W-W (Wald-Wolfowitz)
- MOSES

Tests for  $k$  independent samples are:

- K-W (Kruskal-Wallis)
- MEDIAN

Tests are described below in alphabetical order.

### **Options**

**Statistical Display.** In addition to the tests, you can request univariate statistics, quartiles, and counts for all variables that are specified on the command. You can also control the pairing of variables in tests for two related samples.

**Random Sampling.** NPAR TESTS must store cases in memory when computing tests that use ranks. You can use random sampling when there is not enough space to store all cases.

### **Basic Specification**

The basic specification is a single test subcommand and a list of variables to be tested. Some tests require additional specifications. CHISQUARE has an optional subcommand.

### **Subcommand Order**

Subcommands can be used in any order.

### **Syntax Rules**

- The STATISTICS, SAMPLE, and MISSING subcommands are optional. Each subcommand can be specified only once per NPAR TESTS command.
- You can request any or all tests, and you can specify a test subcommand more than once on a single NPAR TESTS command.
- If you specify a variable more than once on a test subcommand, only the first variable is used.
- Keyword ALL in any variable list refers to all user-defined variables in the active dataset.
- Keyword WITH controls pairing of variables in two-related-samples tests.
- Keyword BY introduces the grouping variable in two- and  $k$ -independent-samples tests.
- Keyword PAIRED can be used with keyword WITH on the MCNEMAR, SIGN, and WILCOXON subcommands to obtain sequential pairing of variables for two related samples.

### **Operations**

- If a string variable is specified on any subcommand, NPAR TESTS will stop executing.
- When ALL is used, requests for tests of variables with themselves are ignored and a warning is displayed.



**Limitations**

- A maximum of 100 subcommands is allowed.
- A maximum of 48,000 variables total per NPAR TESTS command is allowed.
- A maximum of 200 values for subcommand CHISQUARE is allowed.

**BINOMIAL Subcommand**

```
NPAR TESTS BINOMIAL [({.5})]=varlist[({value,value})]
                    {p }           {value      }
```

BINOMIAL tests whether the observed distribution of a dichotomous variable is the same as what is expected from a specified binomial distribution. By default, each named variable is assumed to have only two values, and the distribution of each named variable is compared to a binomial distribution with  $p$  (the proportion of cases expected in the first category) equal to 0.5. The default output includes the number of valid cases in each group, the test proportion, and the two-tailed probability of the observed proportion.

**Syntax**

- The minimum specification is a list of variables to be tested.
- To change the default 0.5 test proportion, specify a value in parentheses immediately after keyword BINOMIAL.
- A single value in parentheses following the variable list is used as a cutting point. Cases with values that are equal to or less than the cutting point form the first category; the remaining cases form the second category.
- If two values appear in parentheses after the variable list, cases with values that are equal to the first value form the first category, and cases with values that are equal to the second value form the second category.
- If no values are specified, the variables must be dichotomous. The first value encountered in the dataset defines the first category, and the other value defines the second category.

**Operations**

- The proportion observed in the first category is compared to the test proportion. The probability of the observed proportion occurring given the test proportion and a binomial distribution is then computed. A test statistic is calculated for each variable specified.
- If the test proportion is the default (0.5), a two-tailed probability is displayed. For any other test proportion, a one-tailed probability is displayed. The direction of the one-tailed test depends on the observed proportion in the first category. If the observed proportion is more than the test proportion, the significance of observing that many or more in the first category is reported. If the observed proportion is less than or equal to the test proportion, the significance of observing that many or fewer in the first category is reported. In other words, the test is always done in the observed direction.

**Example**

```
NPAR TESTS BINOMIAL(.667)=V1(0,1).
```

- NPAR TESTS displays the Binomial Test table, showing the number of cases, observed proportion, test proportion (0.667), and the one-tailed significance for each category.
- If more than 0.667 of the cases have value 0 for *VI*, BINOMIAL gives the probability of observing that many or more values of 0 in a binomial distribution with probability 0.667. If fewer than 0.667 of the cases are 0, the test will be of observing that many or fewer values.

## CHISQUARE Subcommand

```
NPAR TESTS CHISQUARE=varlist [(lo,hi)] [/EXPECTED={EQUAL**      }]
                                         {f1,f2,..., fn}
```

The CHISQUARE (alias CHI-SQUARE) one-sample test computes a chi-square statistic based on the differences between the observed and expected frequencies of categories of a variable. By default, equal frequencies are expected in each category. The output includes the frequency distribution, expected frequencies, residuals, chi-square, degrees of freedom, and probability.

### Syntax

- The minimum specification is a list of variables to be tested. Optionally, you can specify a value range in parentheses following the variable list. You can also specify expected proportions with the EXPECTED subcommand.
- If you use the EXPECTED subcommand to specify unequal expected frequencies, you must specify a value greater than 0 for each observed category of the variable. The expected frequencies are specified in ascending order of category value. You can use the notation  $n*f$  to indicate that frequency  $f$  is expected for  $n$  consecutive categories.
- Specifying keyword EQUAL on the EXPECTED subcommand has the same effect as omitting the EXPECTED subcommand.
- EXPECTED applies to all variables that are specified on the CHISQUARE subcommand. Use multiple CHISQUARE and EXPECTED subcommands to specify different expected proportions for variables.

### Operations

- If no range is specified for the variables that are to be tested, a separate Chi-Square Frequency table is produced for each variable. Each distinct value defines a category.
- If a range is specified, integer-valued categories are established for each value within the range. Non-integer values are truncated before classification. Cases with values that are outside the specified range are excluded. One combined Chi-Square Frequency table is produced for all specified variables.
- Expected values are interpreted as proportions, not absolute values. Values are summed, and each value is divided by the total to calculate the proportion of cases expected in the corresponding category.
- A test statistic is calculated for each specified variable.

### Example

```
NPAR TESTS CHISQUARE=V1 (1,5) /EXPECTED= 12, 3*16, 18.
```

- This example requests the chi-square test for values 1 through 5 of variable *V1*.
- The observed frequencies for variable *V1* are compared with the hypothetical distribution of 12/78 occurrences of value 1; 16/78 occurrences each of values 2, 3, and 4; and 18/78 occurrences of value 5.

## **COCHRAN Subcommand**

```
NPAR TESTS COCHRAN=varlist
```

COCHRAN calculates Cochran's  $Q$ , which tests whether the distribution of values is the same for  $k$  related dichotomous variables. The output shows the frequency distribution for each variable in the Cochran Frequencies table and the number of cases, Cochran's  $Q$ , degrees of freedom, and probability in the Test Statistics table.

### **Syntax**

- The minimum specification is a list of two variables.
- The variables must be dichotomous and must be coded with the same two values.

### **Operations**

- A  $k \times 2$  contingency table (variables by categories) is constructed for dichotomous variables, and the proportions for each variable are computed. A single test is calculated, comparing all variables.
- Cochran's  $Q$  statistic has approximately a chi-square distribution.

### **Example**

```
NPAR TESTS COCHRAN=RV1 TO RV3.
```

- This example tests whether the distribution of values 0 and 1 for *RV1*, *RV2*, and *RV3* is the same.

## **FRIEDMAN Subcommand**

```
NPAR TESTS FRIEDMAN=varlist
```

FRIEDMAN tests whether  $k$  related samples have been drawn from the same population. The output shows the mean rank for each variable in the Friedman Ranks table and the number of valid cases, chi-square, degrees of freedom, and probability in the Test Statistics table.

### **Syntax**

- The minimum specification is a list of two variables.
- Variables should be at least at the ordinal level of measurement.

**Operations**

- The values of  $k$  variables are ranked from 1 to  $k$  for each case, and the mean rank is calculated for each variable over all cases.
- The test statistic has approximately a chi-square distribution. A single test statistic is calculated, comparing all variables.

**Example**

```
NPAR TESTS FRIEDMAN=V1 V2 V3
  /STATISTICS=DESCRIPTIVES.
```

- This example tests variables  $V1$ ,  $V2$ , and  $V3$ , and the example requests univariate statistics for all three variables.

**J-T Subcommand**

```
NPAR TESTS /J-T=varlist BY variable(value1,value2)
```

J-T (alias JONCKHEERE-TERPSTRA) performs the Jonckheere-Terpstra test, which tests whether  $k$  independent samples that are defined by a grouping variable are from the same population. This test is particularly powerful when the  $k$  populations have a natural ordering. The output shows the number of levels in the grouping variable; the total number of cases; observed, standardized, mean, and standard deviation of the test statistic; the two-tailed asymptotic significance; and, if a /METHOD subcommand is specified, one-tailed and two-tailed exact or Monte Carlo probabilities. This subcommand is available only if the Exact Tests option is installed.

**Syntax**

- The minimum specification is a test variable, the keyword BY, a grouping variable, and a pair of values in parentheses.
- Every value in the range defined by the pair of values for the grouping variable forms a group.
- If the /METHOD subcommand is specified, and the number of populations,  $k$ , is greater than 5, the  $p$  value is estimated by using the Monte Carlo sampling method. The exact  $p$  value is not available when  $k$  exceeds 5.

**Operations**

- Cases from the  $k$  groups are ranked in a single series, and the rank sum for each group is computed. A test statistic is calculated for each variable that is specified before BY.
- The Jonckheere-Terpstra statistic has approximately a normal distribution.
- Cases with values other than values in the range that is specified for the grouping variable are excluded.
- The direction of a one-tailed inference is indicated by the sign of the standardized test statistic.

**Example**

```
NPAR TESTS /J-T=V1 BY V2(0,4)
```

```
/METHOD=EXACT.
```

- This example performs the Jonckheere-Terpstra test for groups that are defined by values 0 through 4 of *V2*. The exact *p* values are calculated.

## ***K-S Subcommand (One-Sample)***

```
NPAR TESTS K-S({NORMAL [mean, stddev]})=varlist
              {POISSON [mean]      }
              {UNIFORM [min,max]   }
              {EXPONENTIAL [mean] }
```

The *K-S* (alias *KOLMOGOROV-SMIRNOV*) one-sample test compares the cumulative distribution function for a variable with a uniform, normal, Poisson, or exponential distribution, and the test tests whether the distributions are homogeneous. The parameters of the test distribution can be specified; the defaults are the observed parameters. The output shows the number of valid cases, parameters of the test distribution, most-extreme absolute, positive, and negative differences, Kolmogorov-Smirnov *Z*, and two-tailed probability for each variable.

### ***Syntax***

The minimum specification is a distribution keyword and a list of variables. The distribution keywords are *NORMAL*, *POISSON*, *EXPONENTIAL*, and *UNIFORM*.

- The distribution keyword and its optional parameters must be enclosed within parentheses.
- The distribution keyword must be separated from its parameters by blanks or commas.

<b>NORMAL</b> [mean, stdev]	<i>Normal distribution.</i> The default parameters are the observed mean and standard deviation.
<b>POISSON</b> [mean]	<i>Poisson distribution.</i> The default parameter is the observed mean.
<b>UNIFORM</b> [min,max]	<i>Uniform distribution.</i> The default parameters are the observed minimum and maximum values.
<b>EXPONENTIAL</b> [mean]	<i>Exponential distribution.</i> The default parameter is the observed mean.

### ***Operations***

- The Kolmogorov-Smirnov *Z* is computed from the largest difference in absolute value between the observed and test distribution functions.
- The *K-S* probability levels assume that the test distribution is specified entirely in advance. The distribution of the test statistic and resulting probabilities are different when the parameters of the test distribution are estimated from the sample. No correction is made. The power of the test to detect departures from the hypothesized distribution may be seriously diminished. For testing against a normal distribution with estimated parameters, consider the adjusted *K-S* Lilliefors test that is available in the *EXAMINE* procedure.
- For a mean of 100,000 or larger, a normal approximation to the Poisson distribution is used.
- A test statistic is calculated for each specified variable.

**Example**

```
NPAR TESTS K-S (UNIFORM) =V1 /K-S (NORMAL, 0, 1) =V2 .
```

- The first `K-S` subcommand compares the distribution of *V1* with a uniform distribution that has the same range as *V1*.
- The second `K-S` subcommand compares the distribution of *V2* with a normal distribution that has a mean of 0 and a standard deviation of 1.

***K-S Subcommand (Two-Sample)***

```
NPAR TESTS K-S=varlist BY variable(value1,value2)
```

`K-S` (alias `KOLMOGOROV-SMIRNOV`) tests whether the distribution of a variable is the same in two independent samples that are defined by a grouping variable. The test is sensitive to any difference in median, dispersion, skewness, and so forth, between the two distributions. The output shows the valid number of cases in each group in the Frequency table. The output also shows the largest absolute, positive, and negative differences between the two groups, the Kolmogorov-Smirnov *Z*, and the two-tailed probability for each variable in the Test Statistics table.

**Syntax**

- The minimum specification is a test variable, the keyword `BY`, a grouping variable, and a pair of values in parentheses.
- The test variable should be at least at the ordinal level of measurement.
- Cases with the first value form one group, and cases with the second value form the other group. The order in which values are specified determines which difference is the largest positive and which difference is the largest negative.

**Operations**

- The observed cumulative distributions are computed for both groups, as are the maximum positive, negative, and absolute differences. A test statistic is calculated for each variable that is named before `BY`.
- Cases with values other than values that are specified for the grouping variable are excluded.

**Example**

```
NPAR TESTS K-S=V1 V2 BY V3 (0,1) .
```

- This example specifies two tests. The first test compares the distribution of *V1* for cases with value 0 for *V3* with the distribution of *V1* for cases with value 1 for *V3*.
- A parallel test is calculated for *V2*.

***K-W Subcommand***

```
NPAR TESTS K-W=varlist BY variable(value1,value2)
```

K-W (alias KRUSKAL-WALLIS) tests whether  $k$  independent samples that are defined by a grouping variable are from the same population. The output shows the number of valid cases and the mean rank of the variable in each group in the Ranks table. The output also shows the chi-square, degrees of freedom, and probability in the Test Statistics table.

### **Syntax**

- The minimum specification is a test variable, the keyword BY, a grouping variable, and a pair of values in parentheses.
- Every value in the range defined by the pair of values for the grouping variable forms a group.

### **Operations**

- Cases from the  $k$  groups are ranked in a single series, and the rank sum for each group is computed. A test statistic is calculated for each variable that is specified before BY.
- Kruskal-Wallis  $H$  has approximately a chi-square distribution.
- Cases with values other than values in the range that is specified for the grouping variable are excluded.

### **Example**

```
NPAR TESTS K-W=V1 BY V2 (0,4) .
```

- This example tests  $V1$  for groups that are defined by values 0 through 4 of  $V2$ .

## **KENDALL Subcommand**

```
NPAR TESTS KENDALL=varlist
```

KENDALL tests whether  $k$  related samples are from the same population.  $W$  is a measure of agreement among judges or raters, where each case is one judge's rating of several items (variables). The output includes the mean rank for each variable in the Ranks table and the valid number of cases, Kendall's  $W$ , chi-square, degrees of freedom, and probability in the Test Statistics table.

### **Syntax**

The minimum specification is a list of two variables.

### **Operations**

- The values of the  $k$  variables are ranked from 1 to  $k$  for each case, and the mean rank is calculated for each variable over all cases. Kendall's  $W$  and a corresponding chi-square statistic are calculated, correcting for ties. In addition, a single test statistic is calculated for all variables.
- $W$  ranges between 0 (no agreement) and 1 (complete agreement).

**Example**

```
DATA LIST /V1 TO V5 1-10.
BEGIN DATA
2 5 4 5 1
3 3 4 5 3
3 4 4 6 2
2 4 3 6 2
END DATA.
NPAR TESTS KENDALL=ALL.
```

- This example tests four judges (cases) on five items (variables *V1* through *V5*).

**M-W Subcommand**

```
NPAR TESTS M-W=varlist BY variable(value1,value2)
```

M-W (alias MANN-WHITNEY) tests whether two independent samples that are defined by a grouping variable are from the same population. The test statistic uses the rank of each case to test whether the groups are drawn from the same population. The output shows the number of valid cases of each group; the mean rank of the variable within each group and the sum of ranks in the Ranks table and the Mann-Whitney *U*; Wilcoxon *W* (the rank sum of the smaller group); *Z* statistic; and probability in the Test Statistics table.

**Syntax**

- The minimum specification is a test variable, the keyword **BY**, a grouping variable, and a pair of values in parentheses.
- Cases with the first value form one group and cases with the second value form the other group. The order in which the values are specified is unimportant.

**Operations**

- Cases are ranked in order of increasing size, and test statistic *U* (the number of times that a score from group 1 precedes a score from group 2) is computed.
- An exact significance level is computed if there are 40 or fewer cases. For more than 40 cases, *U* is transformed into a normally distributed *Z* statistic, and a normal approximation *p* value is computed.
- A test statistic is calculated for each variable that is named before **BY**.
- Cases with values other than values that are specified for the grouping variable are excluded.

**Example**

```
NPAR TESTS M-W=V1 BY V2(1,2).
```

- This example tests *V1* based on the two groups that are defined by values 1 and 2 of *V2*.

**MCNEMAR Subcommand**

```
NPAR TESTS MCNEMAR=varlist [WITH varlist [(PAIRED)]]
```



MCNEMAR tests whether combinations of values between two dichotomous variables are equally likely. The output includes a Crosstabulation table for each pair and a Test Statistics table for all pairs, showing the number of valid cases, chi-square, and probability for each pair.

### **Syntax**

- The minimum specification is a list of two variables. Variables must be dichotomous and must have the same two values.
- If keyword WITH is not specified, each variable is paired with every other variable in the list.
- If WITH is specified, each variable before WITH is paired with each variable after WITH. If PAIRED is also specified, the first variable before WITH is paired with the first variable after WITH, the second variable before WITH is paired with the second variable after WITH, and so on. PAIRED cannot be specified without WITH.
- With PAIRED, the number of variables that are specified before and after WITH must be the same. PAIRED must be specified in parentheses after the second variable list.

### **Operations**

- For the purposes of computing the test statistics, only combinations for which the values for the two variables are different are considered.
- If fewer than 25 cases change values from the first variable to the second variable, the binomial distribution is used to compute the probability.

### **Example**

```
NPAR TESTS MCNEMAR=V1 V2 V3 .
```

- This example performs the MCNEMAR test on variable pairs *V1* and *V2*, *V1* and *V3*, and *V2* and *V3*.

## **MEDIAN Subcommand**

```
NPAR TESTS MEDIAN [(value)]=varlist BY variable(value1,value2)
```

MEDIAN determines whether  $k$  independent samples are drawn from populations with the same median. The independent samples are defined by a grouping variable. For each variable, the output shows a table of the number of cases that are greater than and less than or equal to the median in each category in the Frequency table. The output also shows the number of valid cases, the median, chi-square, degrees of freedom, and probability in the Test Statistics table.

### **Syntax**

- The minimum specification is a single test variable, the keyword BY, a grouping variable, and two values in parentheses.
- If the first grouping value is less than the second value, every value in the range that is defined by the pair of values forms a group, and a  $k$ -sample test is performed.
- If the first value is greater than the second value, two groups are formed by using the two values, and a two-sample test is performed.

- By default, the median is calculated from all cases that are included in the test. To override the default, specify a median value in parentheses following the `MEDIAN` subcommand keyword.

### Operations

- A  $2 \times k$  contingency table is constructed with counts of the number of cases that are greater than the median and less than or equal to the median for the  $k$  groups.
- Test statistics are calculated for each variable that is specified before `BY`.
- For more than 30 cases, a chi-square statistic is computed. For 30 or fewer cases, Fisher's exact procedure (two-tailed) is used instead of chi-square.
- For a two-sample test, cases with values other than the two specified values are excluded.

### Example

```
NPAR TESTS MEDIAN(8.4)=V1 BY V2(1,2) /MEDIAN=V1 BY V2(1,2)
/MEDIAN=V1 BY V3(1,4) /MEDIAN=V1 BY V3(4,1).
```

- The first two `MEDIAN` subcommands test variable  $V1$  grouped by values 1 and 2 of variable  $V2$ . The first test specifies a median of 8.4, and the second test uses the observed median.
- The third `MEDIAN` subcommand requests a four-samples test, dividing the sample into four groups based on values 1, 2, 3, and 4 of variable  $V3$ .
- The last `MEDIAN` subcommand requests a two-samples test, grouping cases based on values 1 and 4 of  $V3$  and ignoring all other cases.

## MH Subcommand

```
NPAR TESTS /MH=varlist [WITH varlist [(PAIRED)]]
```

`MH` performs the marginal homogeneity test, which tests whether combinations of values between two paired ordinal variables are equally likely. The marginal homogeneity test is typically used in repeated measures situations. This test is an extension of the McNemar test from binary response to multinomial response. The output shows the number of distinct values for all test variables; the number of valid off-diagonal cell counts; mean; standard deviation; observed and standardized values of the test statistics; the asymptotic two-tailed probability for each pair of variables; and, if a `/METHOD` subcommand is specified, one-tailed and two-tailed exact or Monte Carlo probabilities.

This subcommand is available only if the Exact Tests option is installed (available only on Windows operating systems).

### Syntax

- The minimum specification is a list of two variables. Variables must be polychotomous and must have more than two values. If the variables contain only two values, the McNemar test is performed.
- If keyword `WITH` is not specified, each variable is paired with every other variable in the list.

- If `WITH` is specified, each variable before `WITH` is paired with each variable after `WITH`. If `PAIRED` is also specified, the first variable before `WITH` is paired with the first variable after `WITH`, the second variable before `WITH` is paired with the second variable after `WITH`, and so on. `PAIRED` cannot be specified without `WITH`.
- With `PAIRED`, the number of variables that are specified before and after `WITH` must be the same. `PAIRED` must be specified in parentheses after the second variable list.

### Operations

- The data consist of paired, dependent responses from two populations. The marginal homogeneity test tests the equality of two multinomial  $c \times 1$  tables, and the data can be arranged in the form of a square  $c \times c$  contingency table. A  $2 \times c$  table is constructed for each off-diagonal cell count. The marginal homogeneity test statistic is computed for cases with different values for the two variables. Only combinations for which the values for the two variables are different are considered. The first row of each  $2 \times c$  table specifies the category that was chosen by population 1, and the second row specifies the category that was chosen by population 2. The test statistic is calculated by summing the first row scores across all  $2 \times c$  tables.

### Example

```
NPAR TESTS /MH=V1 V2 V3
/METHOD=MC.
```

- This example performs the marginal homogeneity test on variable pairs *V1* and *V2*, *V1* and *V3*, and *V2* and *V3*. The exact *p* values are estimated by using the Monte Carlo sampling method.

## MOSES Subcommand

```
NPAR TESTS MOSES[(n)]=varlist BY variable(value1,value2)
```

The `MOSES` test of extreme reactions tests whether the range of an ordinal variable is the same in a control group and a comparison group. The control and comparison groups are defined by a grouping variable. The output includes a Frequency table, showing, for each variable before `BY`, the total number of cases and the number of cases in each group. The output also includes a Test Statistics table, showing the number of removed outliers, span of the control group before and after outliers are removed, and one-tailed probability of the span with and without outliers.

### Syntax

- The minimum specification is a test variable, the keyword `BY`, a grouping variable, and two values in parentheses.
- The test variable must be at least at the ordinal level of measurement.
- The first value of the grouping variable defines the control group, and the second value defines the comparison group.
- By default, 5% of the cases are trimmed from each end of the range of the control group to remove outliers. You can override the default by specifying a value in parentheses following the `MOSES` subcommand keyword. This value represents an actual number of cases, not a percentage.

**Operations**

- Values from the groups are arranged in a single ascending sequence. The span of the control group is computed as the number of cases in the sequence containing the lowest and highest control values.
- No adjustments are made for tied cases.
- Cases with values other than values that are specified for the grouping variable are excluded.
- Test statistics are calculated for each variable that is named before BY.

**Example**

```
NPAR TESTS MOSES=V1 BY V3 (0,1) /MOSES=V1 BY V3 (1,0) .
```

- The first MOSES subcommand tests *V1* by using value 0 of *V3* to define the control group and value 1 for the comparison group. The second MOSES subcommand reverses the comparison and control groups.

**RUNS Subcommand**

```
NPAR TESTS RUNS({MEAN  })=varlist
                {MEDIAN}
                {MODE  }
                {value }
```

RUNS tests whether the sequence of values of a dichotomized variable is random. The output includes a Run Test table, showing the test value (cut point that is used to dichotomize the variable tested), number of runs, number of cases that are below the cut point, number of cases that are greater than or equal to the cut point, and test statistic *Z* with its two-tailed probability for each variable.

**Syntax**

- The minimum specification is a cut point in parentheses followed by a test variable.
- The cut point can be specified by an exact value or one of the keywords MEAN, MEDIAN, or MODE.

**Operations**

- All tested variables are treated as dichotomous: cases with values that are less than the cut point form one category, and cases with values that are greater than or equal to the cut point form the other category.
- Test statistics are calculated for each specified variable.

**Example**

```
NPAR TESTS RUNS (MEDIAN) =V2 /RUNS (24.5) =V2 /RUNS (1) =V3 .
```

- This example performs three runs tests. The first test tests variable *V2* by using the median as the cut point. The second test also tests *V2* by using 24.5 as the cut point. The third test tests variable *V3*, with value 1 specified as the cut point.

## ***SIGN Subcommand***

```
NPAR TESTS SIGN=varlist [WITH varlist [(PAIRED)] ]
```

*SIGN* tests whether the distribution of two paired variables in a two-related-samples test is the same. The output includes a Frequency table, showing, for each pair, the number of positive differences, number of negative differences, number of ties, and the total number. The output also includes a Test Statistics table, showing the *Z* statistic and two-tailed probability.

### ***Syntax***

- The minimum specification is a list of two variables.
- Variables should be at least at the ordinal level of measurement.
- If keyword *WITH* is not specified, each variable in the list is paired with every other variable in the list.
- If keyword *WITH* is specified, each variable before *WITH* is paired with each variable after *WITH*. If *PAIRED* is also specified, the first variable before *WITH* is paired with the first variable after *WITH*, the second variable before *WITH* is paired with the second variable after *WITH*, and so on. *PAIRED* cannot be specified without *WITH*.
- With *PAIRED*, the number of variables that are specified before and after *WITH* must be the same. *PAIRED* must be specified in parentheses after the second variable list.

### ***Operations***

- The positive and negative differences between the pair of variables are counted. Ties are ignored.
- The probability is taken from the binomial distribution if 25 or fewer differences are observed. Otherwise, the probability comes from the *Z* distribution.
- Under the null hypothesis for large sample sizes, *Z* is approximately normally distributed with a mean of 0 and a variance of 1.

### ***Example***

```
NPAR TESTS SIGN=N1,M1 WITH N2,M2 (PAIRED) .
```

- *N1* is tested with *N2*, and *M1* is tested with *M2*.

## ***W-W Subcommand***

```
NPAR TESTS W-W=varlist BY variable(value1,value2)
```

*W-W* (alias *WALD-WOLFOWITZ*) tests whether the distribution of a variable is the same in two independent samples. A runs test is performed with group membership as the criterion. The output includes a Frequency table, showing the total number of valid cases for each variable that is specified before *BY* and the number of valid cases in each group. The output also includes a Test Statistics table, showing the number of runs, *Z*, and one-tailed probability of *Z*. If ties are present, the minimum and maximum number of possible runs, their *Z* statistics, and one-tailed probabilities are displayed.

**Syntax**

- The minimum specification is a single test variable, the keyword `BY`, a grouping variable, and two values in parentheses.
- Cases with the first value form one group, and cases with the second value form the other group. The order in which values are specified is unimportant.

**Operations**

- Cases are combined from both groups and ranked from lowest to highest, and a runs test is performed, using group membership as the criterion. For ties involving cases from both groups, both the minimum and maximum number of possible runs are calculated. Test statistics are calculated for each variable that is specified before `BY`.
- For a sample size of 30 or less, the exact one-tailed probability is calculated. For a sample size that is greater than 30, the normal approximation is used.
- Cases with values other than values that are specified for the grouping variable are excluded.

**Example**

```
NPAR TESTS W-W=V1 BY V3 (0,1) .
```

- This example ranks cases from lowest to highest based on their values for `V1`, and a runs test is performed. Cases with value 0 for `V3` form one group, and cases with value 1 form the other group.

**WILCOXON Subcommand**

```
NPAR TESTS WILCOXON=varlist [WITH varlist [(PAIRED)] ]
```

`WILCOXON` tests whether the distribution of two paired variables in two related samples is the same. This test takes into account the magnitude of the differences between two paired variables. The output includes a Ranks table, showing, for each pair, the number of valid cases, positive and negative differences, their respective mean and sum of ranks, and the number of ties. The output also includes a Test Statistics table, showing  $Z$  and probability of  $Z$ .

**Syntax**

- The minimum specification is a list of two variables.
- If keyword `WITH` is not specified, each variable is paired with every other variable in the list.
- If keyword `WITH` is specified, each variable before `WITH` is paired with each variable after `WITH`. If `PAIRED` is also specified, the first variable before `WITH` is paired with the first variable after `WITH`, the second variable before `WITH` is paired with the second variable after `WITH`, and so on. `PAIRED` cannot be specified without `WITH`.
- With `PAIRED`, the number of variables that are specified before and after `WITH` must be the same. `PAIRED` must be specified in parentheses after the second variable list.

**Operations**

- The differences between the pair of variables are counted, the absolute differences are ranked, the positive and negative ranks are summed, and the test statistic  $Z$  is computed from the positive and negative rank sums.
- Under the null hypothesis for large sample sizes,  $Z$  is approximately normally distributed with a mean of 0 and a variance of 1.

**Example**

NPAR TESTS WILCOXON=A B WITH C D (PAIRED) .

- This example pairs  $A$  with  $C$  and  $B$  with  $D$ . If PAIRED were not specified, the example would also pair  $A$  with  $D$  and  $B$  with  $C$ .

**STATISTICS Subcommand**

STATISTICS requests summary statistics for variables that are named on the NPAR TESTS command. Summary statistics are displayed in the Descriptive Statistics table before all test output.

- If STATISTICS is specified without keywords, univariate statistics (keyword DESCRIPTIVES) are displayed.

<b>DESCRIPTIVES</b>	<i>Univariate statistics.</i> The displayed statistics include the mean, maximum, minimum, standard deviation, and number of valid cases for each variable named on the command.
<b>QUARTILES</b>	<i>Quartiles and number of cases.</i> The 25th, 50th, and 75th percentiles are displayed for each variable that is named on the command.
<b>ALL</b>	<i>All statistics available on NPAR TESTS.</i>

**MISSING Subcommand**

MISSING controls the treatment of cases with missing values.

- ANALYSIS and LISTWISE are alternatives. However, each of those commands can be specified with INCLUDE.

<b>ANALYSIS</b>	<i>Exclude cases with missing values on a test-by-test basis.</i> Cases with missing values for a variable that is used for a specific test are omitted from that test. On subcommands that specify several tests, each test is evaluated separately. This setting is the default.
<b>LISTWISE</b>	<i>Exclude cases with missing values listwise.</i> Cases with missing values for any variable that is named on any subcommand are excluded from all analyses.
<b>INCLUDE</b>	<i>Include user-missing values.</i> User-missing values are treated as valid values.

## **SAMPLE Subcommand**

NPAR TESTS must store cases in memory. SAMPLE allows you to select a random sample of cases when there is not enough space on your computer to store all cases. SAMPLE has no additional specifications.

- Because sampling would invalidate a runs test, this option is ignored when the RUNS subcommand is used.

## **METHOD Subcommand**

METHOD displays additional results for each requested statistic. If no METHOD subcommand is specified, the standard asymptotic results are displayed. If fractional weights have been specified, results for all methods will be calculated on the weight rounded to the nearest integer. This subcommand is available only if you have the Exact Tests add-on option installed, which is only available on Windows operating systems.

<b>MC</b>	Displays an unbiased point estimate and confidence interval, based on the Monte Carlo sampling method, for all statistics. Asymptotic results are also displayed. When exact results can be calculated, they will be provided instead of the Monte Carlo results. See Exact Tests for situations under which exact results are provided instead of Monte Carlo results.
<b>CIN(n)</b>	Controls the confidence level for the Monte Carlo estimate. CIN is available only when /METHOD=MC is specified. CIN has a default value of 99.0. You can specify a confidence interval between 0.01 and 99.9, inclusive.
<b>SAMPLES</b>	Specifies the number of tables that were sampled from the reference set when calculating the Monte Carlo estimate of the exact $p$ value. Larger sample sizes lead to narrower confidence limits but also take longer to calculate. You can specify any integer between 1 and 1,000,000,000 as the sample size. SAMPLES has a default value of 10,000.
<b>EXACT</b>	Computes the exact significance level for all statistics, in addition to the asymptotic results. If both the EXACT and MC keywords are specified, only exact results are provided. Calculating the exact $p$ value can be memory-intensive. If you have specified /METHOD=EXACT and find that you have insufficient memory to calculate results, close any other applications that are currently running. You can also enlarge the size of your swap file (see your Windows manual for more information). If you still cannot obtain exact results, specify /METHOD=MC to obtain the Monte Carlo estimate of the exact $p$ value. An optional TIMER keyword is available if you choose /METHOD=EXACT.
<b>TIMER(n)</b>	Specifies the maximum number of minutes during which the exact analysis for each statistic can run. If the time limit is reached, the test is terminated, no exact results are provided, and the program begins to calculate the next test in the analysis. TIMER is available only when /METHOD=EXACT is specified. You can specify any integer value for TIMER. Specifying a value of 0 for TIMER turns the timer off completely. TIMER has a default value of 5 minutes. If a test exceeds a time limit of 30 minutes, it is recommended that you use the Monte Carlo method, rather than the exact method.

## **References**

Siegel, S., and N. J. Castellan. 1988. *Nonparametric statistics for the behavioral sciences*. New York: McGraw-Hill, Inc..



# NUMERIC

```
NUMERIC varlist[(format)] [/varlist...]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
NUMERIC V1.
```

## **Overview**

NUMERIC declares new numeric variables that can be referred to in the transformation language before they are assigned values. Commands such as COMPUTE, IF, RECODE, and COUNT can be used to assign values to the new numeric variables.

### **Basic Specification**

The basic specification is the name of the new variables. By default, variables are assigned a format of F8.2 (or the format that is specified on the SET command).

### **Syntax Rules**

- A FORTRAN-like format can be specified in parentheses following a variable or variable list. Each specified format applies to all variables in the list. To specify different formats for different groups of variables, separate each format group with a slash.
- Keyword TO can be used to declare multiple numeric variables. The specified format applies to each variable that is named and implied by the TO construction.
- NUMERIC can be used within an input program to predetermine the order of numeric variables in the dictionary of the active dataset. When used for this purpose, NUMERIC must precede DATA LIST in the input program.

### **Operations**

- NUMERIC takes effect as soon as it is encountered in the command sequence. Special attention should be paid to the position of NUMERIC among commands. [For more information, see the topic Command Order on p. 41.](#)
- The specified formats (or the defaults) are used as both print and write formats.
- Permanent or temporary variables are initialized to the system-missing value. Scratch variables are initialized to 0.
- Variables that are named on NUMERIC are added to the working file in the order in which they are specified. The order in which they are used in transformations does not affect their order in the active dataset.

## Examples

### Declaring Multiple Numeric Variables

```
NUMERIC V1 V2 (F4.0) / V3 (F1.0).
```

- NUMERIC declares variables *V1* and *V2* with format *F4.0* and declares variable *V3* with format *F1.0*.

```
NUMERIC V1 TO V6 (F3.1) / V7 V10 (F6.2).
```

- NUMERIC declares variables *V1*, *V2*, *V3*, *V4*, *V5*, and *V6* with format *F3.1* and declares variables *V7* and *V10* with format *F6.2*.

### Specifying Variable Order in the Active Dataset

```
NUMERIC SCALE85 IMPACT85 SCALE86 IMPACT86 SCALE87 IMPACT87
      SCALE88 IMPACT88.
```

- Variables *SCALE85* to *IMPACT88* are added to the active dataset in the order that is specified on NUMERIC. The order in which they are used in transformations does not affect their order in the active dataset.

```
INPUT PROGRAM.
STRING CITY (A24).
NUMERIC POP81 TO POP83 (F9)/ REV81 TO REV83(F10).
DATA LIST FILE=POPDATA RECORDS=3
  /1 POP81 22-30 REV81 31-40
  /2 POP82 22-30 REV82 31-40
  /3 POP83 22-30 REV83 31-40
  /4 CITY 1-24(A).
END INPUT PROGRAM.
```

- STRING and NUMERIC are specified within an input program to predetermine variable order in the active dataset. Though data in the file are in a different order, the working file dictionary uses the order that is specified on STRING and NUMERIC. Thus, *CITY* is the first variable in the dictionary, followed by *POP81*, *POP82*, *POP83*, *REV81*, *REV82*, and *REV83*.
- Formats are specified for the variables on NUMERIC. Otherwise, the program uses the default numeric format (*F8.2*) from the NUMERIC command for the dictionary format, even though it uses the format on DATA LIST to read the data. In other words, the dictionary uses the first formats specified, even though DATA LIST may use different formats to read cases.

# OLAP CUBES

OLAP CUBES is available in the Statistics Base option.

```
OLAP CUBES {varlist} BY varlist [BY...]  
  
[/CELLS= [MEAN**] [COUNT**] [STDDEV**]  
[NPCT**] [SPCT**] [SUM**]  
[MEDIAN] [GMEDIAN] [SEMEAN] [MIN] [MAX] [RANGE]  
[VARIANCE] [KURT] [SEKURT] [SKEW] [SESKEW] [FIRST]  
[LAST] [NPCT(var)] [SPCT(var)] [HARMONIC] [GEOMETRIC]  
[DEFAULT] [ALL] [NONE] ]  
  
[/CREATE [ {'catname'}... ] = {GAC } (gvarname {(gvarvalue gvarvalue) }  
{DEFAULT } {GPC } [ {(gvarvalue gvarvalue)...} ] ]  
{GAC GPC}  
  
--or--  
  
{VAC } {(svarname svarname)}  
{VPC } {(svarname svarname)...}  
{VAC VPC}  
  
[/TITLE = 'string'] [FOOTNOTE= 'string']  
[/HIDESMALLCOUNTS [COUNT = {5 } ]  
{integer}]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 19.0

- HIDESMALLCOUNTS subcommand introduced.

## Example

```
OLAP CUBES sales BY quarter by region.
```

## Overview

OLAP CUBES produces summary statistics for continuous, quantitative variables within categories defined by one or more categorical grouping variables.

## Basic Specification

The basic specification is the command name, OLAP CUBES, with a summary variable, the keyword BY, and one or more grouping variables.

- The minimum specification is a summary variable, the keyword `BY`, and a grouping variable.
- By default, `OLAP CUBES` displays a Case Processing Summary table showing the number and percentage of cases included, excluded, and their total, and a Layered Report showing means, standard deviations, sums, number of cases for each category, percentage of total N, and percentage of total sum.

### **Syntax Rules**

- Both numeric and string variables can be specified. String variables can be short or long. Summary variables must be numeric.
- String specifications for `TITLE` and `FOOTNOTE` cannot exceed 255 characters. Values must be enclosed in quotes. When the specification breaks on multiple lines, enclose each line in quotes and separate the specifications for each line by at least one blank. To specify line breaks in titles and footnotes, use the `\n` specification.
- Each subcommand can be specified only once. Multiple use results in a warning, and the last specification is used.
- When a variable is specified more than once, only the first occurrence is honored. The same variables specified after different `BY` keywords will result in an error.

### **Limitations**

- Up to 10 `BY` keywords can be specified.
- See [SET TABLERENDER](#) for information on how multiple layer dimensions are rendered in lightweight tables.

### **Operations**

- The data are processed sequentially. It is not necessary to sort the cases before processing. If a `BY` keyword is used, the output is always sorted.
- A Case Processing Summary table is always generated, showing the number and percentage of the cases included, excluded, and the total.
- For each combination of grouping variables specified after different `BY` keywords, `OLAP CUBES` produces a group in the report.

## **Examples**

```
OLAP CUBES
  tenure age income services monthly BY churn BY region
  /CELLS=COUNT MEAN STDDEV MEDIAN
  /TITLE='Descriptive Statistics' /FOOTNOTE 'By Customer Churn and Geographic Region'.
```

- The procedure produces summaries of *tenure*, *age*, *income*, *services*, and *monthly*, grouped by values of *churn* and *region*.
- `CELLS` requests cell counts, means, standard deviations, and medians.
- `TITLE` produces the requested title for the OLAP cube and `FOOTNOTE` produces the specified table footnote.

## Options

**Cell Contents.** By default, OLAP CUBES displays means, standard deviations, cell counts, sums, percentage of total N, and percentage of total sum. Optionally, you can request any combination of available statistics.

**Group Differences.** You can display arithmetic and/or percentage differences between categories of a grouping variable or between different variables with the CREATE subcommand.

**Format.** You can specify a title and a caption for the report using the TITLE and FOOTNOTE subcommands.

## TITLE and FOOTNOTE Subcommands

TITLE and FOOTNOTE provide a title and a caption for the Layered Report.

- TITLE and FOOTNOTE are optional and can be placed anywhere.
- The specification on TITLE or FOOTNOTE is a string within quotes. To specify a multiple-line title or footnote, enclose each line in quotes and separate the specifications for each line by at least one blank.
- To insert line breaks in the displayed title or footnote, use the \n specification.
- The string you specify cannot exceed 255 characters.

## CELLS Subcommand

By default, OLAP CUBES displays the means, standard deviations, number of cases, sum, percentage of total cases, and percentage of total sum.

- If CELLS is specified without keywords, OLAP CUBES displays the default statistics.
- If any keywords are specified on CELLS, only the requested information is displayed.

<b>DEFAULT</b>	<i>Means, standard deviations, cell counts, sum, percentage of total N, and percentage of total sum. This is the default if CELLS is omitted.</i>
<b>MEAN</b>	<i>Cell means.</i>
<b>STDDEV</b>	<i>Cell standard deviations.</i>
<b>COUNT</b>	<i>Cell counts.</i>
<b>MEDIAN</b>	<i>Cell median.</i>
<b>GMEDIAN</b>	<i>Grouped median.</i>
<b>SEMEAN</b>	<i>Standard error of cell mean.</i>
<b>SUM</b>	<i>Cell sums.</i>
<b>MIN</b>	<i>Cell minimum.</i>
<b>MAX</b>	<i>Cell maximum.</i>
<b>RANGE</b>	<i>Cell range.</i>
<b>VARIANCE</b>	<i>Variances.</i>
<b>KURT</b>	<i>Cell kurtosis.</i>
<b>SEKURT</b>	<i>Standard error of cell kurtosis.</i>
<b>SKEW</b>	<i>Cell skewness.</i>

<b>SESKEW</b>	<i>Standard error of cell skewness.</i>
<b>FIRST</b>	<i>First value.</i>
<b>LAST</b>	<i>Last value.</i>
<b>SPCT</b>	<i>Percentage of total sum.</i>
<b>NPCT</b>	<i>Percentage of total number of cases.</i>
<b>SPCT(var)</b>	<i>Percentage of total sum within specified variable. The specified variable must be one of the grouping variables.</i>
<b>NPCT(var)</b>	<i>Percentage of total number of cases within specified variable. The specified variable must be one of the grouping variables.</i>
<b>HARMONIC</b>	<i>Harmonic mean.</i>
<b>GEOMETRIC</b>	<i>Geometric mean.</i>
<b>ALL</b>	<i>All cell information.</i>

## CREATE Subcommand

CREATE allows you to calculate and display arithmetic and percentage differences between groups or between variables. You can also define labels for these difference categories.

**GAC (gvar(cat1 cat2))** *Arithmetic difference (change) in the summary variable(s) statistics between each specified pair of grouping variable categories.* The keyword must be followed by a grouping variable name specified in parentheses, and the variable name must be followed by one or more pairs of grouping category values. Each pair of values must be enclosed in parentheses inside the parentheses that contain the grouping variable name. String values must be enclosed in single or double quotation marks. You can specify multiple pairs of category values, but you can only specify one grouping variable, and the grouping variable must be one of the grouping variables specified at the beginning of the OLAP CUBES command, after the BY keyword.

The difference calculated is the summary statistic value for the second category specified minus the summary statistic value for the first category specified:  
 $cat2 - cat1.$

**GPC (gvar(cat1 cat2))** *Percentage difference (change) in the summary variable(s) statistics between each specified pair of grouping variable categories.* The keyword must be followed by a grouping variable name enclosed in parentheses, and the variable name must be followed by one or more pairs of grouping category values. Each pair of values must be enclosed in parentheses inside the parentheses that contain the grouping variable name. String values must be enclosed in single or double quotation marks. You can specify multiple pairs of category values, but you can only specify one grouping variable, and the grouping variable must be one of the grouping variables specified at the beginning of the OLAP CUBES command, after the BY keyword.

The percentage difference calculated is the summary statistic value for the second category specified minus the summary statistic value for the first category specified, divided by the summary statistic value for the first category specified:  $(cat2 - cat1)/cat1.$

**VAC(svar1 svar2)** *Arithmetic difference (change) in summary statistics between each pair of specified summary variables.* Each pair of variables must be enclosed in parentheses, and all specified variables must be specified as summary variables at the beginning of the OLAP CUBES command.

The difference calculated is the summary statistic value for the second variable specified minus the summary statistic value for the first variable specified:  
 $svar2 - svar1.$

<b>VPC(svar1 svar2)</b>	<p><i>Percentage difference (change) in summary statistics between each pair of specified summary variables. Each pair of variables must be enclosed in parentheses, and all specified variables must be specified as summary variables at the beginning of the OLAP CUBES command.</i></p> <p>The percentage difference calculated is the summary statistic value for the second variable specified minus the summary statistic value for the first variable specified: <math>(svar2 - svar1)/svar1</math>.</p>
<b>'category label'</b>	<p><i>Optional label for each difference category created. These labels must be the first specification in the CREATE subcommand. Each label must be enclosed in single or double quotation marks. If no labels are specified, defined value or variable labels are used. If no labels are defined, data values or variable names are displayed. If multiple differences are created, the order of the labels corresponds to the order the differences are specified. To mix custom labels with default labels, use the keyword DEFAULT for the difference categories without custom labels.</i></p>

Both arithmetic and percentage differences can be specified in the same command, but you cannot specify both grouping variable differences (GAC/GPC) and summary variable differences (VAC/VPC) in the same command.

#### Example

```
OLAP CUBES
sales96 BY region
/CELLS=SUM NPCT
/CREATE GAC GPC (region (1 3) (2 3)).
```

- Both the arithmetic (GAC) and percentage (GPC) differences will be calculated.
- Differences will be calculated for two different pairs of categories of the grouping variable *region*.
- The grouping variable specified in the CREATE subcommand, *region*, is also specified as a grouping variable at the beginning of the OLAP CUBES command.

#### Example

```
OLAP CUBES
sales95 sales96 BY region
/CELLS=SUM NPCT
/CREATE VAC VPC (sales95 sales96).
```

- Both the arithmetic (VAC) and percentage (VPC) differences will be calculated.
- The difference calculated will be *sales96 - sales95*.
- The percentage difference calculated will be  $(sales96 - sales95)/sales95$ .
- The two variables, *sales95* and *sales96* are also specified as summary variables at the beginning of the OLAP CUBES command.

#### Example

```
OLAP CUBES
sales96 BY region
/CELLS=SUM NPCT
/CREATE DEFAULT 'West-East GPC' DEFAULT 'West-Central % Difference'
GAC GPC (region (1 3) (2 3)).
```

- Four labels are specified, corresponding to the four difference categories that will be created: arithmetic and percentage differences between regions 3 and 1 and between regions 3 and 2.
- The two `DEFAULT` labels will display the defined value labels or values if there aren't any value labels for the two arithmetic (GAC) difference categories.

## ***HIDESMALLCOUNTS Subcommand***

`HIDESMALLCOUNTS` allows you to hide counts displayed in tables for count values that are less than a specified integer. Hidden values are displayed as <N, where N is the specified integer.

### ***COUNT Keyword***

The `COUNT` keyword specifies the cutoff, below which count values are hidden. Values equal to the cutoff are not hidden. The value of `COUNT` must be an integer and greater than or equal to 2. The default is 5.



# OMS

*Note:* Square brackets used in the OMS syntax chart are required parts of the syntax and are not used to indicate optional elements. Any equals signs (=) displayed in the syntax chart are required. All subcommands except DESTINATION are optional.

```
OMS

/SELECT CHARTS HEADINGS LOGS MODELS
        TABLES TEXTS TREES WARNINGS
or
/SELECT ALL EXCEPT = [list]

/IF  COMMANDS = ["expression" "expression"...]
     SUBTYPES = ["expression" "expression"...]
     LABELS   = ["expression" "expression"...]
     INSTANCES = [n n... LAST]

/EXCEPTIF (same keywords as IF, except for INSTANCES)

/DESTINATION
FORMAT=
    {DOC      }
    {HTML    }
    {OXML    }
    {PDF     ]
    {TABTEXT }
    {TEXT    }
    {SAV     }
    {SPV     }
    {SPW     ]
    {XLS     }

{OUTFILE = "outfile expression"          }
{XMLWORKSPACE="name"                    }
{OUTPUTSET = {SUBTYPES} FOLDER = "dirspec"}
    {LABELS   }

VIEWER={YES|NO}
NUMBERED = 'varname'
IMAGES={YES|NO}
IMAGEFORMAT={PNG|JPG|EMF|BMP|VML}
IMAGEROOT='rootname'
CHARTSIZE=percent
IMAGEMAP={NO|YES}
TREEFORMAT={XML|IMAGE}
CHARTFORMAT={XML|IMAGE}
MODELFORMAT={XML|IMAGE}
TABLES={PIVOTABLE|STATIC}

/COLUMNS DIMNAMES = ["dimension1" "dimension2" ...]
or
/COLUMNS SEQUENCE = [R1 R2 ... RALL C1 C2... CALL L1 L2... LALL]

/TAG = "string"

/NOWARN
```

- NUMBERED only applies to FORMAT=SAV.
- IMAGES applies only to FORMAT=OXML, HTML, and SPV.
- IMAGEFORMAT applies only to FORMAT=OXML, HTML, SPV, and SPW.
- IMAGEROOT and CHARTSIZE only apply to FORMAT=OXML and HTML.

- `IMAGEMAP` only applies to `FORMAT=HTML`.
- `TREEFORMAT` only applies to `FORMAT=OXML` and `SPV`.
- `CHARTFORMAT` only applies to `FORMAT=OXML`.
- `TABLES` only applies to `FORMAT=SPW` files used in IBM® SPSS® Collaboration and Deployment Services.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

#### Release 13.0

- `TREES` keyword introduced on `SELECT` subcommand.
- `IMAGES`, `IMAGEROOT`, `CHARTSIZE`, and `IMAGEFORMAT` keywords introduced on `DESTINATION` subcommand.

#### Release 14.0

- `XMLWORKSPACE` keyword introduced on `DESTINATION` subcommand.

#### Release 16.0

- `IMAGEFORMAT=VML` introduced for `FORMAT=HTML` on `DESTINATION` subcommand.
- `IMAGEMAP` keyword introduced for `FORMAT=HTML` on `DESTINATION` subcommand.
- `FORMAT=SPV` introduced for saving output in Viewer format.
- `CHARTFORMAT` keyword introduced.
- `TREEFORMAT` keyword introduced.
- `TABLES` keyword introduced.
- `FORMAT=SVWSOXML` is no longer supported.

#### Release 17.0

- `MODELS` keyword introduced on `SELECT` subcommand.
- `FORMAT=DOC`, `XLS`, `PDF`, and `SPW` introduced.
- `MODELFORMAT` keyword introduced.

#### Release 19.0

- `IMAGEFORMAT=VML` introduced for `FORMAT=OXML` on `DESTINATION` subcommand.

### **Example**

```
OMS
  /DESTINATION FORMAT = OXML OUTFILE = '/mydir/myfile.xml'
  VIEWER = NO.
OMS
  /SELECT TABLES
  /IF COMMANDS = ['Regression']
    SUBTYPES = ['Coefficients']
  /DESTINATION FORMAT = SAV
```

```
OUTFILE = '/mydir/regression_coefficients.sav'.  
REGRESSION  
  /STATISTICS COEFF OUTS R ANOVA  
  /DEPENDENT income  
  /METHOD=ENTER age address ed employ.  
OMSEND.
```

## Overview

The OMS command controls the routing and format of output from IBM® SPSS® Statistics to files and can suppress Viewer output. Output formats include: Word, Excel, PDF, SPSS Statistics data file format (SAV), Viewer file format (SPV), web report format (SPW), XML, HTML, and text.

### Basic Specification

The basic specification is the command name followed by a `DESTINATION` subcommand that contains a `FORMAT` and/or a `VIEWER` specification. For `FORMAT`, an `OUTFILE` or `OUTPUTSET` specification is also required.

### Syntax Rules

- All subcommands except `DESTINATION` are optional. No subcommand may occur more than once in each OMS command.
- Multiple OMS commands are allowed. For more information, see [Basic Operation below](#).
- Subcommands can appear in any order.
- If duplicates are found in a list, they are ignored except in `/COLUMNS SEQUENCE` where they cause an error.
- When a keyword takes a square-bracketed list, the brackets are required even if the list contains only a single item.

### Limitations

- For OXML format, the `HEADINGS` keyword has no effect. If any output from a procedure is included, the procedure title output is included.
- If the OMS specification results in nothing other than Headings objects or a Notes tables being included for a procedure, then nothing is included for that procedure.

## Basic Operation

- Once an OMS command is executed, it remains in effect until the end of the session or until ended by an `OMSEND` command.
- A destination file specified on an OMS command is unavailable to other commands and other applications until the OMS command is ended by an `OMSEND` command or the end of the session.
- While an OMS command is in effect, the specified destination files are stored in memory (RAM), so active OMS commands that write a large amount of output to external files may consume a large amount of memory.

- Multiple OMS commands are independent of each other (except as noted below). The same output can be routed to different locations in different formats based on the specifications in different OMS commands.
- Display of output objects in the Viewer is determined by the most recent OMS command that includes the particular output type. For example, if an OMS command includes all tables from the `FREQUENCIES` command and also contains a `VIEWER = YES` specification, and a subsequent OMS command includes all tables of the subtype 'Statistics' with `VIEWER = NO`, Statistics tables for subsequent `FREQUENCIES` commands will *not* be displayed in the Viewer.
- The `COLUMNS` subcommand has no effect on pivot tables displayed in the Viewer.
- The order of the output objects in any particular destination is the order in which they were created, which is determined by the order and operation of the commands that generate the output.

## ***SELECT Subcommand***

`SELECT` specifies the types of output objects to be routed to the specified destination(s). You can select multiple types. You can also specify `ALL` with `EXCEPT` to exclude specified types. If there is no `SELECT` subcommand, all supported output types are selected.

**ALL.** *All output objects.* This is the default.

**CHARTS.** *Chart objects.* This includes charts created by the commands such as `GRAPH` and `GGRAPH` and charts created by statistical procedures (for example, the `BARChart` subcommand of the `FREQUENCIES` command). It does not include tree diagrams produced by the `TREE` procedure or charts in models displayed in the Model Viewer.

**HEADINGS.** *Text objects labeled Title in the outline pane of the Viewer.* For destination format `OXML`, the `HEADINGS` keyword has no effect. If any output for the procedure is included in `OXML` output, the procedure title is included as the `text` attribute value of the `command` element.

**LOGS.** *Log text objects.* Log objects contain certain types of error and warning messages. With `SET PRINTBACK=ON`, log objects also contain the command syntax executed during the session. Log objects are labeled *Log* in the outline pane of the Viewer.

**MODELS.** *Output objects displayed in the Model Viewer.* A single model object can contain multiple views of the model, including both tables and charts.

**TABLES.** *Output objects that are pivot tables in the Viewer.* This includes Notes tables. It does not include tables displayed in the Model Viewer. Pivot tables are the only output objects that can be routed to the destination format `SAV`.

**TEXTS.** *Text objects that aren't logs or headings.* This includes objects labeled *Text Output* in the outline pane of the Viewer.

**TREES.** *Tree model diagrams produced by the TREE procedure (Decision Tree option).*

**WARNINGS.** *Warnings objects.* Warnings objects contain certain types of error and warning messages.

**Example**

```

OMS /SELECT TABLES LOGS TEXTS WARNINGS HEADINGS
/DESTINATION FORMAT = HTML OUTFILE = '/mypath/myfile1.htm'.
OMS /SELECT ALL EXCEPT = [CHARTS]
/DESTINATION FORMAT = HTML OUTFILE = '/mypath/myfile2.htm'.

```

The two SELECT subcommands are functionally equivalent. The first one explicitly lists all types but CHARTS, and the second one explicitly excludes only CHARTS.

Figure 159-1  
Output object types in the Viewer

The screenshot shows the SPSS Viewer window with a menu bar (File, Edit, View, Data, Transform, Insert, Format, Analyze, Graphs, Utilities, Add-ons, Window, Help) and a tree view on the left. The main area displays several output objects:

- Heading:** A box labeled "Heading" points to the title "Frequencies" above the table.
- Table:** A box labeled "Table" points to the table below the heading.
- Chart:** A box labeled "Chart" points to a bar chart below the table.
- Log:** A box labeled "Log" points to the error message text below the chart.
- Warning:** A box labeled "Warning" points to the warning message text at the bottom.

**Table Data:**

		Frequency	Percent	Valid Percent	Cumulative Percent
Valid	Clerical	363	76.6	76.6	76.6
	Custodial	27	5.7	5.7	82.3
	Manager	84	17.7	17.7	100.0
	Total	474	100.0	100.0	

**Bar Chart Data:**

Employment Category	Frequency
Clerical	363
Custodial	27
Manager	84

**Log:**

```

-Error # 1. Command name: comptue
-The first word in the line is not recognized as an SPSS command.
-This command not executed.

```

**Warnings:**

```

Text: newvar
An undefined variable name, or a scratch or system variable was specified in a variable list
which accepts only standard variables. Check spelling and verify the existence of this variable.
This command not executed.

```

**Notes Table Limitation**

An OMS command that selects only tables will not select a Notes table if the Notes tables is the only table produced by a procedure. This can occur if the command contains syntax errors that result in a Notes table and a warning object, but no other tables. For example:

```
DATA LIST FREE /var1 var2.
BEGIN DATA
1 2
END DATA.
OMS SELECT TABLES
  /DESTINATION FORMAT=HTML
  OUTFILE=' /temp/htmltest.htm' .
FREQUENCIES VARIABLES=var1.
DESCRIPTIVES VARIABLES=var02.
OMSEND.
```

The DESCRIPTIVES command refers to a variable that doesn't exist, causing an error that results in the creation of a Notes table and a warning object, but the HTML file will not include this Notes table. To make sure Notes tables are selected when no other tables are created by a procedure, include WARNINGS in the SELECT subcommand, as in:

```
OMS SELECT TABLES WARNINGS
  /DESTINATION FORMAT=HTML
  OUTFILE=' /temp/htmltest.htm' .
```

**IF Subcommand**

The IF subcommand specifies particular output objects of the types determined by SELECT. Without an IF subcommand, all objects of the specified types are selected. If you specify multiple conditions, only those objects that meet *all* conditions will be selected.

**Example**

```
OMS
  /SELECT TABLES
  /IF COMMANDS = ['Regression']
    SUBTYPES = ['Coefficients']
  /DESTINATION FORMAT = SAV
  OUTFILE = '/mydir/regression_coefficients.sav'.
```

This OMS command specifies that only coefficient tables from the REGRESSION command will be selected.

**COMMANDS Keyword**

The COMMANDS keyword restricts the selection to the specified command(s). The keyword COMMANDS must be followed by an equals sign (=) and a list of quoted command identifiers enclosed in square bracket, as in:

```
OMS
  /SELECT TABLES
  /IF COMMANDS = ['Frequencies' 'Factor Analysis']
  /DESTINATION...
```

Command identifiers are:

- Unique. No two commands have the same identifier.
- Not case-sensitive.
- Not subject to translation, which means they are the same for all language versions and output languages.
- Often not exactly the same or even similar to the command name. You can obtain the identifier for a particular command from the OMS Control Panel (Utilities menu) or by generating output from the command in the Viewer and then right-clicking the command heading in the outline pane and selecting Copy OMS Command Identifier from the context menu.

Command identifiers are available for all statistical and charting procedures and any other commands that produce blocks of output with their own identifiable heading in the outline pane of the Viewer. For example, CASESTOVARS and VARSTOCASES have corresponding identifiers ('Cases to Variables' and 'Variables to Cases') because they produce their own output blocks (with command headings in the outline pane that happen to match the identifiers), but FLIP does not because any output produced by FLIP is included in a generic Log text object.

## ***SUBTYPES Keyword***

The SUBTYPES keyword restricts the selection to the specified table types. The keyword SUBTYPES must be followed by an equals sign (=) and a list of quoted subtype identifiers enclosed in square bracket, as in:

```
OMS
/SELECT TABLES
/IF SUBTYPES = ['Descriptive Statistics' 'Coefficients']
/DESTINATION...
```

- Subtypes apply only to tables that would be displayed as pivot tables in the Viewer.
- Like command identifiers, subtype identifiers are not case-sensitive and are not subject to translation.
- Unlike command identifiers, subtype identifiers are not necessarily unique. For example, multiple commands produce a table with the subtype identifier "Descriptive Statistics," but not all of those tables share the same structure. If you want only a particular table type for a particular command, use both the COMMANDS and SUBTYPES keywords.
- The OMS Control Panel (Utilities menu) provides a complete list of subtypes. You can also obtain the identifier for a particular table by generating output from the command in the Viewer and then right-clicking outline item for the Table in the outline pane of the Viewer and selecting Copy OMS Table Subtype from the context menu. The identifiers are generally fairly descriptive of the particular table type.

## ***LABELS Keyword***

The LABELS keyword selects particular output objects according to the text displayed in the outline pane of the Viewer. The keyword LABELS must be followed by an equals sign (=) and a list of quoted label text enclosed in square brackets, as in:

```
OMS
/SELECT TABLES
/IF LABELS = ['Job category * Gender Crosstabulation']
/DESTINATION...
```

The LABELS keyword is useful for differentiating between multiple graphs or multiple tables of the same type in which the outline text reflects some attribute of the particular output object such as the variable names or labels. There are, however, a number of factors that can affect the label text:

- If split file processing is on, split file group identification may be appended to the label.
- Labels that include information about variables or values are affected by the OVARS and ONUMBERS settings on the SET command.
- Labels are affected by the current output language setting (SET OLANG).

## ***INSTANCES Keyword***

The INSTANCES subcommand selects the *n*th instance of an object matching the other criteria on the IF subcommand within a single command execution. The keyword INSTANCES must be followed by an equals sign (=) and a list of positive integers and/or the keyword LAST enclosed in square brackets.

### ***Example***

```
OMS
/SELECT TABLES
/IF COMMANDS = ['Frequencies']
    SUBTYPES = ['Frequencies']
    INSTANCES = [1 LAST]
/DESTINATION...
OMS
/SELECT TABLES
/IF COMMANDS = ['Frequencies']
    INSTANCES = [1 LAST]
/DESTINATION...
```

- The first OMS command will select the first and last frequency tables from each FREQUENCIES command.
- The second OMS command, in the absence of a SUBTYPES or LABELS specification, will select the first and last tables of *any* kind from the selected command. For the FREQUENCIES command (and most other statistical and charting procedures), the first table would be the Notes table.



## Wildcards

For `COMMANDS`, `SUBTYPES`, and `LABELS`, you can use an asterisk (\*) as a wildcard indicator at the end of a quoted string to include all commands, tables, and/or charts that start with that quoted string, as in:

```
OMS
/SELECT TABLES
/IF SUBTYPES = ['Correlation*']
/DESTINATION...
```

In this example, all table subtypes that begin with “Correlation” will be selected.

The values of `LABELS` can contain asterisks as part of the value as in “First variable \* Second variable Crosstabulation,” but only an asterisk as the last character in the quoted string is interpreted as a wildcard, so:

```
OMS
/SELECT TABLES
/IF LABELS = ['First Variable **']
/DESTINATION...
```

will select all tables with labels that start with “First Variable \*”.

## EXCEPTIF Subcommand

The `EXCEPTIF` subcommand excludes specified output object types. It has the same keywords and syntax as `IF`, with the exception of `INSTANCES`, which will cause an error if used with `EXCEPTIF`.

### Example

```
OMS
/SELECT TABLES
/IF COMMANDS = ['Regression']
/EXCEPTIF SUBTYPES = ['Notes' 'Case Summar*']
/DESTINATION...
```

## DESTINATION Subcommand

The `DESTINATION` subcommand is the only required subcommand. It specifies the format and location for the routed output. You can also use this subcommand to control what output is displayed in the Viewer.

- Output continues to flow to a specified destination until its OMS specification is ended, at which point the file is closed. [For more information, see the topic Basic Operation on p. 1405.](#)
- Different OMS commands may refer to the same destination file as long as the `FORMAT` is the same. When a request becomes active, it starts contributing to the appropriate output stream. If the `FORMAT` differs, an error results. When multiple requests target the same destination, the output is written in the order in which it is created, not the order of OMS commands.

### Example

```
OMS
```

```
/DESTINATION FORMAT = OXML OUTFILE = '/mydir/myfile.xml'.
```

## **FORMAT Keyword**

The `DESTINATION` subcommand must include either a `FORMAT` or `VIEWER` specification (or both). The `FORMAT` keyword specifies the format for the routed output. The keyword must be followed by an equals sign (=) and one of the following alternatives:

**DOC.** *Word/RTF format.* Pivot tables are exported as Word tables with all formatting attributes intact—for example, cell borders, font styles, and background colors. Text output is exported as formatted RTF. Charts, tree diagrams, and model views are included in PNG format.

**HTML.** *HTML 4.0.* Output objects that would be pivot tables in the Viewer are converted to simple HTML tables. No TableLook attributes (font characteristics, border styles, colors, etc.) are supported. Text output objects are tagged <PRE> in the HTML. Charts, tree diagrams, and model views are exported as separate files in the selected graphics format and are embedded by reference. Image file names use the HTML file name as the root name, followed by a sequential integer, starting with 0.

**OXML.** *Output XML.* XML that conforms to the spss-output schema ([xml.spss.com/spss/oms](http://xml.spss.com/spss/oms)). [For more information, see the topic OXML Table Structure on p. 1430.](#) You can specify the version of the spss-output schema that generates the XML output and you can specify whether language invariant versions of text attributes are included in the generated XML. These options are specified with the `XVERSION` and `OATTRS` subcommands of the `SET` command. [For more information, see the topic XVERSION and OATTRS Subcommands on p. 1882.](#)

**PDF.** *Portable Document Format.* Output is exported as it would appear in Print Preview, with all formatting attributes intact. The PDF file includes bookmarks that correspond to the entries in the Viewer outline pane.

**SAV.** *IBM® SPSS® Statistics format data file.* This format is a binary file format. All output object types other than tables are excluded. Each column of a table becomes a variable in the data file. To use a data file that is created with OMS in the same session, you must end the active OMS request before you can open the data file. [For more information, see the topic Routing Output to SAV Files on p. 1422.](#) In Unicode mode, the file encoding is UTF-8; in code page mode, the file encoding is the code page determined by the current locale. See [SET command, UNICODE subcommand](#) for more information.

**SPV.** *Viewer file format.* This is the same format used when you save the contents of a Viewer window.

**SPW.** *Web report format.* This output file format is designed for use with Predictive Enterprise Services. It is essentially the same as the SPSS Statistics Viewer format except that tree diagrams are saved as static images.

**TEXT.** *Space-separated text.* Output is written as text, with tabular output aligned with spaces for fixed-pitch fonts. Charts, tree diagrams, and model views are excluded. In Unicode mode, the file encoding is UTF-8; in code page mode, the file encoding is the code page determined by the current locale. See [SET command, UNICODE subcommand](#) for more information.

**TABTEXT.** *Tab-delimited text.* For output that is displayed as pivot tables in the Viewer, tabs delimit table column elements. Text block lines are written as is; no attempt is made to divide them with tabs at useful places. Charts, tree diagrams, and model views are excluded. In Unicode mode, the file encoding is UTF-8; in code page mode, the file encoding is the code page determined by the current locale. See [SET command](#), [UNICODE subcommand](#) for more information.

**XLS.** *Excel 97-2003 format.* Pivot table rows, columns, and cells are exported as Excel rows, columns, and cells, with all formatting attributes intact — for example, cell borders, font styles, and background colors. Text output is exported with all font attributes intact. Each line in the text output is a row in the Excel file, with the entire contents of the line contained in a single cell. Charts, tree diagrams, and model views are included in PNG format.

## **NUMBERED Keyword**

For `FORMAT=SAV`, you can also specify the `NUMBERED` keyword to identify the source tables, which can be useful if the data file is constructed from multiple tables. This creates an additional variable in the data file. The value of the variable is a positive integer that indicates the sequential table number. The default variable name is `TableNumber_`. You can override the default with an equals sign (=) followed by a valid variable name in quotes after the `NUMBERED` keyword.

### **Example**

```
OMS
/SELECT TABLES
/IF COMMANDS = ['Regression'] SUBTYPES = ['Coefficients']
/DESTINATION = SAV NUMBERED = 'Table_number'
OUTFILE = 'data.sav'.
```

## **IMAGES and IMAGEFORMAT Keywords**

For HTML and OXML document format, you can include charts, tree diagrams, and model views in a number of different graphic formats with the `IMAGES` and `IMAGEFORMAT` keywords. For SPV and SPW formats, these settings only apply to tree diagrams, not charts or model views.

### **IMAGES Keyword**

- `IMAGES=YES` produces a separate image file for each chart and/or tree diagram. Image files are saved in a separate subdirectory (folder). The subdirectory name is the name of the destination file without any extension and with “\_files” appended to the end. This is the default setting.

*Note:* When the destination is the XML workspace (`XMLWORKSPACE` keyword), images are stored in the workspace rather than in a physical file. [For more information, see the topic XMLWORKSPACE Keyword on p. 1417.](#)

- For HTML document format, standard `<IMG SRC='filename'>` tags are included in the HTML document for each image file.
- For charts in OXML document format, the XML file contains `chart` elements with an `ImageFile` attribute of the general form `<chart imageFile="filepath/filename"/>` for each image file. `IMAGES=YES` has no effect on `FORMAT=OXML` unless you also specify

CHARTFORMAT=IMAGE. For more information, see the topic [CHARTFORMAT Keyword](#) on p. 1416.

- For model views in OXML document format, the XML file contains `modelview` elements with an `ImageFile` attribute of the general form `<modelview imageFile="filepath/filename"/>` for each image file. `IMAGES=YES` has no effect on `FORMAT=OXML` unless you also specify `MODELFORMAT=IMAGE`. For more information, see the topic [MODELFORMAT Keyword](#) on p. 1416.
- For HTML format, `IMAGES=NO` excludes charts, tree diagrams, and modelviews.
- For OXML format, `IMAGES=NO` causes charts, tree diagrams, and model views to be included in the document in XML format instead of as separate image files.

#### **IMAGEFORMAT Keyword**

- PNG is the default image format.
- For HTML document format, the available image formats are PNG, JPG, EMF, BMP, and VML.
- For OXML document format, the available image formats are PNG, JPG, EMF, BMP, and VML.
- For SPV and SPW document format, the available image formats are PNG, JPG, and BMP.
- EMF (enhanced metafile) format is available only on Windows operating systems.
- VML image format does not create separate image files. The VML code that renders the image is embedded in the HTML.
- VML image format does not include tree diagrams.

Note: Not all browsers support VML.

#### **Example**

```
OMS SELECT TABLES CHARTS
  /DESTINATION FORMAT=HTML
                IMAGES=YES
                IMAGEFORMAT=JPG
                OUTFILE=' /htmloutput/julydata.htm' .
```

### **CHARTSIZE and IMAGEROOT Keywords**

For HTML and OXML document formats, you can control the relative size and root name of chart and tree diagram images, if charts and/or tree diagrams are saved as separate image files.

- **CHARTSIZE=n**. Specifies the scaling, expressed as a percentage value between 10 and 200. The default is `CHARTSIZE=100`.
- **IMAGEROOT='rootname'**. User-specified rootname for image files. Image files are constructed from the rootname, an underscore, and a sequential three-digit number. The rootname should be specified in quotes, as in: `IMAGEROOT='julydata'`. When the destination is the XML workspace (`XMLWORKSPACE` keyword), the rootname applies to the name associated with the image in the workspace.

**Example**

```
OMS SELECT TABLES CHARTS
/DESTINATION FORMAT=HTML
IMAGES=YES
CHARTSIZE=50
IMAGEROOT='rutabaga'
OUTFILE='/htmloutput/julydata.htm'.
```

**IMAGEMAP Keyword**

For HTML document format, you can use the `IMAGEMAP` keyword to create image map tooltips that display information for some chart elements, such as the value of the selected point on a line chart or bar on a bar chart. The default is `IMAGEMAP=NO`. To include image map tooltips, use `IMAGEMAP=YES`.

`IMAGEMAP` has no effect on tree diagram images or document formats other than HTML.

**Example**

```
OMS SELECT TABLES CHARTS
/DESTINATION FORMAT=HTML
IMAGES=YES
IMAGEFORMAT=PNG
IMAGEMAP=YES
OUTFILE='/htmloutput/julydata.htm'.
```

**TREEFORMAT Keyword**

For OXML and SPV document formats, `TREEFORMAT` controls the format of tree diagrams (produced by the `TREE` command). The keyword is followed by and equals sign and one of the following alternatives:

- **XML.** Tree diagrams are included as XML that conforms to the pmml schema ([www.dmg.org](http://www.dmg.org)). For SPV format, this is the format required to activate and edit trees in the Viewer window. This is the default.
- **IMAGE.** For SPV format, tree diagrams are included in the Viewer document as static images in the selected format. For OXML format, image files are saved in a separate folder. [For more information, see the topic IMAGES and IMAGEFORMAT Keywords on p. 1413.](#)

*Note:* When the destination is the XML workspace (`XMLWORKSPACE` keyword), images are stored in the workspace rather than in a physical file. [For more information, see the topic XMLWORKSPACE Keyword on p. 1417.](#)

`IMAGES=NO` overrides the `TREEFORMAT` setting and includes tree diagrams in XML format. [For more information, see the topic IMAGES and IMAGEFORMAT Keywords on p. 1413.](#)

**Example**

```
OMS SELECT TABLES TREES
/DESTINATION FORMAT=SPV
IMAGES=YES
IMAGEFORMAT=PNG
```

```
TREEFORMAT=IMAGE  
OUTFILE='/viewerdocs/results.spv'.
```

### ***CHARTFORMAT Keyword***

For OXML document format, `CHARTFORMAT` controls the format of charts. The keyword is followed by and equals sign and one of the following alternatives:

- **XML.** Charts are included as XML that conforms to the vizml schema ([xml.spss.com/spss/visualization](http://xml.spss.com/spss/visualization)). This is the default.
- **IMAGE.** For SPV format, charts are included in the Viewer document as static images in the selected format. Image files are saved in the selected format in a separate folder. [For more information, see the topic IMAGES and IMAGEFORMAT Keywords on p. 1413.](#)

*Note:* When the destination is the XML workspace (`XMLWORKSPACE` keyword), images are stored in the workspace rather than in a physical file. [For more information, see the topic XMLWORKSPACE Keyword on p. 1417.](#)

`IMAGES=NO` overrides the `CHARTFORMAT` setting and includes charts in XML format. [For more information, see the topic IMAGES and IMAGEFORMAT Keywords on p. 1413.](#)

### ***MODELFORMAT Keyword***

For OXML document format, `MODELFORMAT` controls the format of models. The keyword is followed by and equals sign and one of the following alternatives:

- **XML.** Models are included as XML that conforms to the PMML schema ([www.dmg.org](http://www.dmg.org)), plus additional XML that specifies data and styles. This is the default.
- **IMAGE.** The OXML contains `modelview` elements that specify the image file names. The image files are saved in the selected format in a separate folder. One model may produce multiple image files. [For more information, see the topic IMAGES and IMAGEFORMAT Keywords on p. 1413.](#)

*Note:* When the destination is the XML workspace (`XMLWORKSPACE` keyword), images are stored in the workspace rather than in a physical file. [For more information, see the topic XMLWORKSPACE Keyword on p. 1417.](#)

`IMAGES=NO` overrides the `MODELFORMAT` setting and includes models in XML format. [For more information, see the topic IMAGES and IMAGEFORMAT Keywords on p. 1413.](#)

XML output is not available for models created by the `LINEAR` and `GENLINMIXED` commands.

### ***TABLES Keyword***

For SPW format, the `TABLES` keyword controls the format of pivot tables. The keyword is followed by an equals sign and one of the following alternatives:

- **PIVOTABLE.** Tables are included as dynamic pivot tables. This is the default.
- **STATIC.** Tables cannot be pivoted.

## **OUTFILE Keyword**

If a `FORMAT` is specified, the `DESTINATION` subcommand must also include either an `OUTFILE`, `XMLWORKSPACE`, or `OUTPUTSET` specification. `OUTFILE` specifies an output file. The keyword must be followed by an equals sign (=) and a file specification in quotes or a previously defined file handle defined with the `FILE HANDLE` command. With `FORMAT=SAV`, you can specify a previously defined dataset name instead of a file.

### **Example**

```
OMS
/DESTINATION FORMAT = OXML OUTFILE = '/mydir/myfile.xml'.
```

## **XMLWORKSPACE Keyword**

For `FORMAT=OXML`, you can route the output to an area in memory referred to as the **XML workspace** where it can be used in flow control and other programming features available with `BEGIN PROGRAM-END PROGRAM`. The keyword must be followed by an equals sign (=) and a quoted string specifying a name that identifies this output in the workspace.

### **Example**

```
OMS SELECT TABLES
/IF COMMANDS=['Frequencies'] SUBTYPES=['Frequencies']
/DESTINATION FORMAT=OXML XMLWORKSPACE='freq_table'.
```

For more information, see the documentation for the `EvaluateXPath` function in the Help system.

## **OUTPUTSET Keyword**

`OUTPUTSET` is an alternative to `OUTFILE` that allows you to route each output object to a separate file. The keyword must be followed by an equals sign (=) and one of the following alternatives:

- |                 |   |
|-----------------|---|
| <b>LABELS</b>   | <i>Output file names based on output object label text.</i> Label text is the text that appears in the outline pane of the Viewer. <a href="#">For more information, see the topic LABELS Keyword on p. 1410.</a> |
| <b>SUBTYPES</b> | <i>Output file names based on subtype identifiers.</i> Subtypes apply only to tables. <a href="#">For more information, see the topic SUBTYPES Keyword on p. 1409.</a>  |

### **Example**

```
OMS
/SELECT TABLES
/DESTINATION FORMAT = OXML OUTPUTSET = SUBTYPES.
```

- `OUTPUTSET` will not overwrite existing files. If a specified file name already exists, an underscore and a sequential integer will be appended to the file name.

## **FOLDER Keyword**

With *OUTPUTSET*, you can also use the *FOLDER* keyword to specify the location for the routed output. Since you may not know what is considered to be the “current” directory, it’s probably a good idea to explicitly specify the location. The keyword must be followed by an equals sign (=) and a valid location specification in quotes.

### **Example**

```
OMS
/SELECT TABLES
/IF SUBTYPES = ['Frequencies' 'Descriptive Statistics']
/DESTINATION FORMAT = OXML OUTPUTSET = SUBTYPES
  FOLDER = '/maindir/nextdir/newdir'.
```

- If the last folder (directory) specified on the path does not exist, it will be created.
- If any folders prior to the last folder on the path do not already exist, the specification is invalid, resulting in an error.

## **VIEWER Keyword**

By default, output is displayed in the Viewer as well as being routed to other formats and destinations specified with the *FORMAT* keyword. You can use *VIEWER = NO* to suppress the Viewer display of output for the specified output types. The *VIEWER* keyword can be used without the *FORMAT* keyword (and associated *OUTFILE* or *OUTPUTSET* keywords) to simply control what output is displayed in the Viewer.

### **Example**

```
OMS
/SELECT TABLES
/IF SUBTYPES = ['Correlations*']
/DESTINATION FORMAT SAV OUTFILE = '/mydir/myfile.sav'
  VIEWER = NO.
OMS
/SELECT TABLES
/IF SUBTYPES = ['NOTES']
/DESTINATION VIEWER = NO.
```

- The first OMS command will route tables with subtype names that start with “Correlation” to data file in IBM® SPSS® Statistics format and will not display those tables in the Viewer. All other output will be displayed in the Viewer
- The second OMS command simply suppresses the Viewer display of all Notes tables, without routing the Notes table output anywhere else.

Note: The *VIEWER* keyword has no effect on OMS output saved to external formats or files, including the Viewer SPV and SPW formats. It also has no effect on output saved to SPV format in a batch job executed with the Batch Facility (available with SPSS Statistics Server).



## ***COLUMNS Subcommand***

You can use the `COLUMNS` subcommand to specify the dimension elements that should appear in the columns. All other dimension elements appear in the rows.

- This subcommand applies only to tables that would be displayed as pivot tables in the Viewer and is ignored without warning if the `OMS` command does not include any tables.
- With `DESTINATION FORMAT = SAV`, columns become variables in the data file. If you specify multiple dimension elements on the `COLUMNS` subcommand, then variable names will be constructed by combining nested element and column labels. [For more information, see the topic Routing Output to SAV Files on p. 1422.](#)
- The `COLUMNS` subcommand has no effect on pivot tables displayed in the Viewer.
- If you specify multiple dimension elements, they are nested in the columns in the order in which they are listed on the `COLUMNS` subcommand. For example: `COLUMNS DIMNAMES=['Variables' 'Statistics']` will nest statistics within variables in the columns.
- If a table doesn't contain any of the dimension elements listed, then all dimension elements for that table will appear in the rows.

## ***DIMNAMES Keyword***

The `COLUMNS` subcommand must be followed by either the `DIMNAMES` or `SEQUENCE` keyword. Each dimension of a table may contain zero or more elements. For example, a simple two-dimensional crosstabulation contains a single row dimension element and a single column dimension element, each with labels based on the variables in those dimensions, plus a single layer dimension element labeled *Statistics* (if English is the output language). These element labels may vary based on the output language (`SET OLANG`) and/or settings that affect the display of variable names and/or labels in tables (`SET TVARS`). The keyword `DIMNAMES` must be followed by an equals sign (=) and a list of quoted dimension element labels enclosed in square brackets.

### ***Example***

```
OMS
/SELECT TABLES
/IF COMMANDS = ['Correlations' 'Frequencies']
/DESTINATION FORMAT = SAV OUTPUTSET = SUBTYPES
/COLUMNS DIMNAMES = ['Statistics'].
```

The labels associated with the dimension elements may not always be obvious. To see all the dimension elements and their labels for a particular pivot table:

- ▶ Activate (double-click) the table in the Viewer.
  - ▶ From the menus choose `View > Show All`.
- and/or
- ▶ If the pivoting trays aren't displayed, from the menus choose `Pivot > Pivoting Trays`.
  - ▶ Hover over each icon in the pivoting trays for a ToolTip pop-up that displays the label.

Figure 159-2  
Displaying table dimension element labels

**Table without Dimension Element Labels Displayed**

		Current Salary	Beginning Salary
Female	Mean	\$26,031.92	\$13,091.97
	Median	\$24,300.00	\$12,375.00
Male	Mean	\$41,441.78	\$20,301.40
	Median	\$32,850.00	\$15,750.00
Total	Mean	\$34,419.57	\$17,016.09
	Median	\$28,875.00	\$15,000.00

**Table with Dimension Element Labels Displayed**

		Variables	
Gender	Statistics	Current Salary	Beginning Salary
Female	Mean	\$26,031.92	\$13,091.97
	Median	\$24,300.00	\$12,375.00
Male	Mean		
	Median		
Total	Mean		
	Median		

## SEQUENCE Keyword

SEQUENCE is an alternative to DIMNAMES that uses positional arguments. These positional arguments do not vary based on output language or output display settings. The SEQUENCE keyword must be followed by an equals sign (=) and a list of positional arguments enclosed in square brackets.

- The general form of a positional argument is a letter indicating the default position of the element—C for column, R for row, or L for layer—followed by a positive integer indicating the default position within that dimension. For example, R1 would indicate the outermost row dimension element.
- A letter indicating the default dimension followed by ALL indicates all elements in that dimension in their default order. For example, RALL would indicate all row dimension elements, and CALL by itself would be unnecessary since it would not alter the default arrangement of the table. ALL cannot be combined with positional sequence numbers in the same dimension.
- SEQUENCE=[CALL RALL LALL] will put all dimension elements in the columns. With FORMAT=SAV, this will result in one case per table in the data file.

**Example**

```

OMS
/SELECT TABLES
/IF COMMANDS = ['Regression'] SUBTYPES = ['Coefficient Correlations']
/DESTINATION FORMAT = SAV OUTFILE = '/mydir/myfile.sav'
/COLUMNS SEQUENCE = [R1 R2].

```

Figure 159-3  
Positional arguments for dimension elements

Model	R1	R2	R3	Previous Experience (months)	Months since Hire	Beginning Salary	Date of Birth
1	Correlations	Previous Experience (months)		1.000	.067	-.087	.805
		Months since Hire		.067	1.000	.012	.085
		Beginning Salary		-.087	.012	1.000	-.075
		Date of Birth		.805	.085	-.075	1.000
	Covariances	Previous Experience (months)		31.307	12.940	-.022	7.096E-06
		Months since Hire		12.940	1205.248	.019	4.635E-06
		Beginning Salary		-.022	.019	.002	-5.236E-09
		Date of Birth		7.096E-06	4.635E-06	-5.236E-09	2.485E-12

**TAG Subcommand**

OMS commands remain in effect until the end of the session or until you explicitly end them with the OMSEND command, and you can have multiple OMS commands in effect at the same time. You can use the TAG subcommand to assign an ID value to each OMS command, which allows you to selectively end particular OMS commands with a corresponding TAG keyword on the OMSEND command. The ID values assigned on the TAG subcommand are also used to identify OMS commands in the log created by the OMSLOG command.

**Example**

```

OMS
/DESTINATION FORMAT = OXML OUTFILE = '/mydir/myfile.xml'
/TAG = 'oxmlout'.

```

- The TAG subcommand must be followed by an equals sign (=) and a quoted ID value.
- The ID value cannot start with a dollar sign.
- Multiple active OMS commands cannot use the same TAG value.

See OMSEND and OMSLOG for more information.

**NOWARN Subcommand**

The NOWARN subcommand suppresses all warnings from OMS. The NOWARN subcommand applies only to the current OMS command. It has no additional specifications.

## ***Routing Output to SAV Files***

IBM® SPSS® Statistics data files consist of variables in the columns and cases in the rows, and that's essentially how pivot tables are converted to data files:

- Columns in the table are variables in the data file. Valid variable names are constructed from the column labels.
- Row labels in the table become variables with generic variable names (*Var1*, *Var2*, *Var3*...) in the data file. The values of these variables are the row labels in the table.
- Three table-identifier variables are automatically included in the data file: *Command\_*, *Subtype\_*, and *Label\_*. All three are string variables. The first two are the command and subtype identifiers. *Label\_* contains the table title text.
- Rows in the table become cases in the data file.

### ***Data File Created from One Table***

Data files can be created from one or more tables. There are two basic variations for data files created from a single table:

- Data file created from a two-dimensional table with no layers.
- Data file created from a three-dimension table with one or more layers.

#### ***Example***

In the simplest case—a single, two-dimensional table—the table columns become variables and the rows become cases in data file.

Figure 159-4  
Single two-dimensional table

**Report**

Gender		Current Salary	Beginning Salary
Female	Mean	\$26,031.92	\$13,091.97
	Median	\$24,300.00	\$12,375.00
	Minimum	\$15,750	\$9,000
	Maximum	\$58,125	\$30,000
Male	Mean	\$41,441.78	\$20,301.40
Total			

	Command	Subtype	Label	Var1	Var2	CurrentSalary	BeginningSal
1	Means	Report	Report	Female	Mean	\$26,031.92	\$13,091.97
2	Means	Report	Report	Female	Median	\$24,300.00	\$12,375.00
3	Means	Report	Report	Female	Minimum	\$15,750	\$9,000
4	Means	Report	Report	Female	Maximum	\$58,125	\$30,000
5	Means	Report	Report	Male	Mean	\$41,441.78	\$20,301.40
6	Means	Report	Report	Male	Median	\$32,850.00	\$15,750.00
7	Means	Report	Report	Male	Minimum	\$19,650	\$9,000
8	Means	Report	Report	Male	Maximum	\$135,000	\$79,980
9	Means	Report	Report	Total	Mean	\$34,419.57	\$17,016.09
10	Means	Report	Report	Total	Median	\$28,875.00	\$15,000.00

- The first three variables identify the source table by command, subtype, and label.
- The two elements that defined the rows in the table—values of the variable *Gender* and statistical measures—are assigned the generic variable names *Var1* and *Var2*. These are both string variables.
- The column labels from the table are used to create valid variable names. In this case, those variable names are based on the variable labels of the three scale variables summarized in the table. If the variables didn't have defined variable labels or you chose to display variable names instead of variable labels as the column labels in the table, then the variable names in the new data file would be the same as in the source data file.

### Example

If the default table display places one or more elements in layers, additional variables are created to identify the layer values.

Figure 159-5  
Table with layers

Layer	Minority Classification	Gender	
		Female	Male
Employment Category	Clerical	40	47
	Custodial/Manager		

Layer	Minority Classification	Gender	
		Female	Male
Employment Category	Clerical	166	110
	Custodial	0	14
	Manager	10	70

	Var1	Var2	Var3	Var4	Female	Male
1	Employment Category	Clerical	Minority Classification	No	166	110
2	Employment Category	Clerical	Minority Classification	Yes	40	47
3	Employment Category	Custodial	Minority Classification	No	0	14
4	Employment Category	Custodial	Minority Classification	Yes	0	13
5	Employment Category	Manager	Minority Classification	No	10	70
6	Employment Category	Manager	Minority Classification	Yes	0	4

- In the table, the variable labeled *Minority Classification* defines the layers. In the data file, this creates two additional variables: one that identifies the layer element, and one that identifies the categories of the layer element.
- As with the variables created from the row elements, the variables created from the layer elements are string variables with generic variable names (the prefix *Var* followed by a sequential number).

### Data Files Created from Multiple Tables

When multiple tables are routed to the same data file, each table is added to the data file in a fashion similar to the `ADD FILES` command.

- Each subsequent table will always add cases to the data file.
- If column labels in the tables differ, each table may also add variables to the data file, with missing values for cases from other tables that don't have an identically labeled column.

#### Example

Multiple tables that contain the same column labels will typically produce the most immediately useful data files (data files that don't require additional manipulation).

Figure 159-6  
Multiple tables with the same column labels

Gender					
		Frequency	Percent	Valid Percent	Cumulative Percent
Valid	Female	216	45.6	45.6	45.6
	Male	258	54.4	54.4	100.0

Employment Category					
		Frequency	Percent	Valid Percent	Cumulative Percent
Valid	Clerical	363	76.6	76.6	76.6
	Custodial	27	5.7	5.7	82.3
	Manager	84	17.7	17.7	100.0
	Total	474	100.0	100.0	

The screenshot shows the SPSS Data Editor window titled "temp.sav - Data Editor". The menu bar includes File, Edit, View, Data, Transform, Analyze, Graphs, Utilities, Add-ons, Window, and Help. The toolbar contains icons for file operations, editing, and analysis. The main window displays a table with the following structure:

1: Command_							
	Label_	Var1	Var2	Frequency	Percent	ValidPercent	CumulativePercent
1	Gender	Valid	Female	216	45.6	45.6	45.6
2	Gender	Valid	Male	258	54.4	54.4	100.0
3	Gender	Valid	Total	474	100.0	100.0	.
4	Employment Category	Valid	Clerical	363	76.6	76.6	76.6
5	Employment Category	Valid	Custodial	27	5.7	5.7	82.3
6	Employment Category	Valid	Manager	84	17.7	17.7	100.0
7	Employment Category	Valid	Total	474	100.0	100.0	.

The bottom of the window shows "Data View" and "Variable View" tabs, with "Data View" selected.

- The second table contributes additional cases (rows) to the data file but no new variables because the column labels are exactly the same; so there are no large patches of missing data.
- Although the values for *Command\_* and *Subtype\_* are the same, the *Label\_* value identifies the source table for each group of cases because the two frequency tables have different titles.

### Example

A new variable is created in the data file for each unique column label in the tables routed to the data file, which will result in blocks of missing values if the tables contain different column labels.

Figure 159-7  
Multiple tables with different column labels

Report			
Gender		Current Salary	Beginning Salary
Female	Mean	\$26,031.92	\$13,091.97
	Median	\$24,300.00	\$12,375.00
Male	Mean	\$41,441.78	\$20,301.40
Total	Mean	\$34,420.00	\$17,016.10

Report			
Gender		Educational Level (years)	Months since Hire
Female	Mean	12.37	80.38
	Median	12.00	81.00
Male	Mean	14.43	81.72
Total	Mean	13.49	81.11
Total	Median	12.00	81.00

temp.sav - Data Editor

File Edit View Data Transform Analyze Graphs Utilities Add-ons Window Help

16: Command\_

	Label_	Var1	Var2	Current Salary	Beginning Salary	Educational Levelyears	Monthssince Hire
1	Report	Female	Mean	\$26032	\$13092.0	.	.
2	Report	Female	Median	\$24300	\$12375.0	.	.
3	Report	Male	Mean	\$41442	\$20301.4	.	.
4	Report	Male	Median	\$32850	\$15750.0	.	.
5	Report	Total	Mean	\$34420	\$17016.1	.	.
6	Report	Total	Median	\$28875	\$15000.0	.	.
7	Report	Female	Mean	.	.	12.37	80.38
8	Report	Female	Median	.	.	12.00	81.00
9	Report	Male	Mean	.	.	14.43	81.72
10	Report	Male	Median	.	.	15.00	82.00
11	Report	Total	Mean	.	.	13.49	81.11
12	Report	Total	Median	.	.	12.00	81.00

Data View Variable View

- The first table has columns labeled *Beginning Salary* and *Current Salary*, which are not present in the second table, resulting in missing values for those variables for cases from the second table.
- Conversely, the second table has columns labeled *Education level* and *Months since hire*, which are not present in the first table, resulting in missing values for those variables for cases from the first table.
- Mismatched variables, such as those in this example, can occur even with tables of the same subtype. In fact, in this example, both tables are of the same subtype.

### Data Files Not Created from Multiple Tables

If any tables do not have the same number of row elements as the other tables, no data file will be created. The number of rows doesn't have to be the same; the number of row *elements* that become variables in the data file must be the same.

For example, a two-variable crosstabulation and a three-variable crosstabulation from CROSSTABS contain different numbers of row elements, since the "layer" variable is actually nested within the row variable in the default three-variable crosstabulation display.



**Figure 159-8**  
*Tables with different numbers of row elements*

**Employment Category \* Gender Crosstabulation**

		Gender		Total
		Female	Male	
Employment Category	Clerical	206	157	363
	Custodial	0	27	27
	Manager	10	74	84
Total		216	258	474

**Employment Category \* Gender \* Minority Classification Crosstabulation**

Minority Classification			Gender		Total
			Female	Male	
No	Employment Category	Clerical	166	110	276
		Custodial	0	14	14
		Manager	10	70	80
	Total	176	194	370	
Yes	Employment Category	Clerical	40	47	87
		Custodial	0	13	13
		Manager	0	4	4
	Total	40	64	104	

In general, the less specific the subtype selection in the OMS command, the less likely you are to get sensible data files, or any data files at all. For example:

```
OMS /SELECT TABLES /DESTINATION FORMAT=SAV OUTFILE='mydata.sav' .
```

will probably fail to create a data file more often than not, since it will select all tables, including Notes tables, which have a table structure that is incompatible with most other table types.

## ***Controlling Column Elements to Control Variables in the Data File***

You can use the COLUMNS subcommand to specify which dimension elements should be in the columns and therefore are used to create variables in the generated data file. This is equivalent to pivoting the table in the Viewer.

### ***Example***

The DESCRIPTIVES command produces a table of descriptive statistics with variables in the rows and statistics in the columns. A data file created from that table would therefore use the statistics as variables and the original variables as cases. If you want the original variables to be variables in the generated data file and the statistics to be cases:

```
OMS
/SELECT TABLES
/IF COMMANDS=['Descriptives'] SUBTYPES=['Descriptive Statistics']
/DESTINATION FORMAT=SAV OUTFILE='/temp/temp.sav'
/COLUMNS DIMNAMES=['Variables'].
DESCRIPTIVES VARIABLES=salary salbegin.
OMSEND.
```

- When you use the `COLUMNS` subcommand, any dimension elements not listed on the subcommand will become rows (cases) in the generated data file.
- Since the descriptive statistics table has only two dimension elements, the syntax `COLUMNS DIMNAMES=['Variables']` will put the variables in the columns *and* will put the statistics in the row. So this is equivalent to swapping the positions of the original row and column elements.

Figure 159-9

Default and pivoted table and generated data file

**Default Descriptive Statistics Table**

	N	Minimum	Maximum	Mean	Std. Deviation
Current Salary	474	\$15,750	\$135,000	\$34,419.57	\$17,075.661
Beginning Salary	474	\$9,000	\$79,980	\$17,016.09	\$7,870.638
Valid N (listwise)	474				

**Descriptive Statistics Table with Variables in Columns and Statistics in Rows**

	Current Salary	Beginning Salary	Valid N (listwise)
N	474	474	474
Minimum	\$15,750	\$9,000	
Maximum	\$135,000	\$79,980	
Mean	\$34,419.57	\$17,016.09	
Std. Deviation	\$17,075.661	\$7,870.638	

### Example

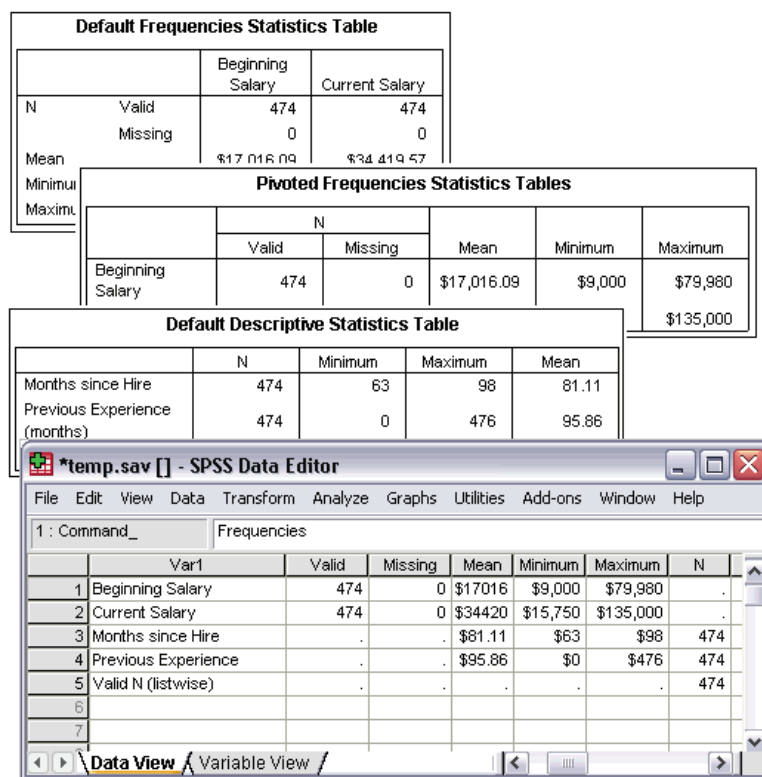
The `FREQUENCIES` command produces a descriptive statistics table with statistics in the rows, while the `DESCRIPTIVES` command produces a descriptive statistics table with statistics in the columns. To include both table types in the same data file in a meaningful fashion, you need to change the column dimension for one of them.

```
OMS
/SELECT TABLES
/IF COMMANDS=['Frequencies' 'Descriptives']
SUBTYPES=['Statistics' 'Descriptive Statistics']
/DESTINATION FORMAT=SAV OUTFILE='/temp/temp.sav'
/COLUMNS DIMNAMES=['Statistics'].
FREQUENCIES
VARIABLES=salbegin salary
/FORMAT=NOTABLE
/STATISTICS=MINIMUM MAXIMUM MEAN.
DESCRIPTIVES
```

```
VARIABLES=jobtime prevexp
/STATISTICS=MEAN MIN MAX.
OMSEND.
```

- The COLUMNS subcommand will be applied to all selected table types that have a *Statistics* dimension element.
- Both table types have a *Statistics* dimension element, but since it's already in the column dimension for the table produced by the DESCRIPTIVES command, the COLUMNS subcommand has no effect on the structure of the data from that table type.
- For the *FREQUENCIES* statistics table, COLUMNS DIMNAMES=['Statistics'] is equivalent to pivoting the *Statistics* dimension element into the columns and pivoting the *Variables* dimension element into the rows.
- Some of the variables will have missing values, since the table structures still aren't exactly the same with statistics in the columns.

Figure 159-10  
Combining different table types in same data file



## Variable Names

OMS constructs valid, unique variable names from column labels.

- Row and layer elements are assigned generic variable names: the prefix `var` followed by a sequential number.

- Characters that aren't allowed in variable names, such as spaces and parentheses, are removed. For example, "This (Column) Label" would become a variable named *ThisColumnLabel*.
- If the label begins with a character that is allowed in variable names but not allowed as the first character (for example, a number), "@" is inserted as a prefix. For example "2nd" would become a variable named *@2nd*.
- Underscores or periods at the end of labels are removed from the resulting variable names. (The underscores at the end of the automatically generated variables *Command\_*, *Subtype\_*, and *Label\_* are not removed.)
- If more than one element is in the column dimension, variable names are constructed by combining category labels with underscores between category labels. Group labels are not included. For example, if *VarB* is nested under *VarA* in the columns, you would get variables like *CatA1\_CatB1*, not *VarA\_CatA1\_VarB\_CatB1*.

Figure 159-11  
Variable names in SAV files

Layer Variable		Column Variable			Total
		In	(Out)		
High	Row Variable	Yes	1	1	2
		No	0	1	1
	Total		1	2	3
Low	Row Variable	Yes	1	0	1
		No	1	1	2
	Total		2	1	3

Layer Variable		High			Low		
		Column Variable			Column Variable		
		In	(Out)	Total	In	(Out)	Total
Row Variable	Yes	1	1	2	1	0	1
	No	0	1	1	1	1	2
Total		1	2	3	2	1	3

## OXML Table Structure

OXML is XML that conforms to the spss-output schema.

- OMS command and subtype identifiers are used as values of the `command` and `subType` attributes in OXML. For example:

```
<command text="Frequencies" command="Frequencies"...>
<pivotTable text="Gender" label="Gender" subType="Frequencies"...>
```

These attribute values are not affected by output language (`SET OLANG`) or display settings for variable names/labels or values/value labels (`SET TVARS` and `SET TNUMBERS`).

- XML is case-sensitive. The element name `pivotTable` is considered a different element from one named "pivottable" or "Pivottable" (the latter two don't exist in OXML).

- Command and subtype identifiers generated by the OMS Control Panel or the OMS Identifiers dialog box (both on the Utilities menu) use the same case as that used for values of the `command` and `subType` OXML attributes.
- All of the information displayed in a table is contained in attribute values in OXML. At the individual cell level, OXML consists of “empty” elements that contain attributes but no “content” other than that contained in attribute values.
- Table structure in OXML is represented row by row; elements that represent columns are nested within the rows, and individual cells are nested within the column elements:

```

<pivotTable...>
<dimension axis="row"...>
  <dimension axis="column"...>
    <category...>
      <cell text="..." number="..." decimals="..."/>
    </category>
    <category...>
      <cell text="..." number="..." decimals="..."/>
    </category>
  </dimension>
</dimension>
...
</pivotTable>

```

The preceding example is a simplified representation of the structure that shows the descendant/ancestor relationships of these elements, but not necessarily the parent/child relationships, because there are typically intervening nested element levels. The following figures show a simple table as displayed in the Viewer and the OXML that represents that table.

Figure 159-12  
*Simple frequency table*

		Gender			
		Frequency	Percent	Valid Percent	Cumulative Percent
Valid	Female	216	45.6	45.6	45.6
	Male	258	54.4	54.4	100.0
Total		474	100.0	100.0	

Figure 159-13  
*OXML for simple frequency table*

```

<?xml version="1.0" encoding="UTF-8" ?>
<outputTree>
  <command text="Frequencies" command="Frequencies"
    displayTableValues="label" displayOutlineValues="label"
    displayTableVariables="label" displayOutlineVariables="label">
    <pivotTable text="Gender" label="Gender" subType="Frequencies"
      varName="gender" variable="true">
      <dimension axis="row" text="Gender" label="Gender"
        varName="gender" variable="true">
        <group text="Valid">
          <group hide="true" text="Dummy">
            <category text="Female" label="Female" string="f"
              varName="gender">
              <dimension axis="column" text="Statistics">
                <category text="Frequency">

```

```

    <cell text="216" number="216"/>
  </category>
  <category text="Percent">
    <cell text="45.6" number="45.569620253165" decimals="1"/>
  </category>
  <category text="Valid Percent">
    <cell text="45.6" number="45.569620253165" decimals="1"/>
  </category>
  <category text="Cumulative Percent">
    <cell text="45.6" number="45.569620253165" decimals="1"/>
  </category>
</dimension>
</category>
<category text="Male" label="Male" string="m" varName="gender">
  <dimension axis="column" text="Statistics">
    <category text="Frequency">
      <cell text="258" number="258"/>
    </category>
    <category text="Percent">
      <cell text="54.4" number="54.430379746835" decimals="1"/>
    </category>
    <category text="Valid Percent">
      <cell text="54.4" number="54.430379746835" decimals="1"/>
    </category>
    <category text="Cumulative Percent">
      <cell text="100.0" number="100" decimals="1"/>
    </category>
  </dimension>
</category>
</group>
<category text="Total">
  <dimension axis="column" text="Statistics">
    <category text="Frequency">
      <cell text="474" number="474"/>
    </category>
    <category text="Percent">
      <cell text="100.0" number="100" decimals="1"/>
    </category>
    <category text="Valid Percent">
      <cell text="100.0" number="100" decimals="1"/>
    </category>
  </dimension>
</category>
</group>
</dimension>
</pivotTable>
</command>
</outputTree>

```

As you may notice, a simple, small table produces a substantial amount of XML. That's partly because the XML contains some information not readily apparent in the original table, some information that might not even be available in the original table, and a certain amount of redundancy.

- The table contents as they are (or would be) displayed in a pivot table in the Viewer are contained in `text` attributes. For example:

```
<command text="Frequencies" command="Frequencies"...>
```

These text attributes can be affected by both output language (`SET OLANG`) and settings that affect the display of variable names/labels and values/value labels (`SET TVARS` and `SET TNUMBERS`). In this example, the `text` attribute value will differ depending on the

output language, whereas the `command` attribute value remains the same regardless of output language.

- Wherever variables or values of variables are used in row or column labels, the XML will contain a `text` attribute and one or more additional attribute values. For example:

```
<dimension axis="row" text="Gender" label="Gender" varName="gender">  
...  
<category text="Female" label="Female" string="f" varName="gender">
```

For a numeric variable, there would be a `number` attribute instead of a `string` attribute. The `label` attribute is present only if the variable or values have defined labels.

- The `<cell>` elements that contain cell values for numbers will contain the `text` attribute and one or more additional attribute values. For example:

```
<cell text="45.6" number="45.569620253165" decimals="1"/>
```

The `number` attribute is the actual, unrounded numeric value, and the `decimals` attribute indicates the number of decimal positions displayed in the table.

- Since columns are nested within rows, the `category` element that identifies each column is repeated for each row. For example, since the statistics are displayed in the columns, the element `<category text="Frequency">` appears three times in the XML—once for the male row, once for the female row, and once for the total row.

Examples of using XSLT to transform OXML are provided in the Help system.

## ***Command and Subtype Identifiers***

The OMS Control Panel (Utilities menu) provides a complete list of command and subtype identifiers. For any command or table displayed in the Viewer, you can find out the command or subtype identifier by right-clicking the item in the Viewer outline pane.

# OMSEND

*Note:* Square brackets used in the OMSSEND syntax chart are required parts of the syntax and are not used to indicate optional elements. Any equals signs (=) displayed in the syntax chart are required. All specifications other than the command name OMSSEND are optional.

```
OMSEND  
TAG = {[ 'idvalue' 'idvalue'... ]}  
      {ALL  
      }  
FILE = [ 'filespec' 'filespec'... ]  
LOG
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
OMS  
/DESTINATION FORMAT = OXML OUTFILE = '/mydir/myfile.xml'.  
[some commands that produce output]  
OMSEND.  
[some more commands that produce output]
```

## **Overview**

OMSEND ends active OMS commands. The minimum specification is the command name OMSSEND. In the absence of any other specifications, this ends all active OMS commands and logging.

## **TAG Keyword**

The optional TAG keyword identifies specific OMS commands to end, based on the ID value assigned on the OMS TAG subcommand or automatically generated if there is no TAG subcommand. To display the automatically generated ID values for active OMS commands, use the OMSINFO command

The TAG keyword must be followed by an equals sign (=) and a list of quoted ID values or the keyword ALL enclosed in square brackets.

## **Example**

```
OMSEND TAG = [ 'reg_tables_to_sav' 'freq_tables_to_html' ].
```

A warning is issued if any of the specified values don't match any active OMS commands.



## **FILE Keyword**

The optional `FILE` keyword ends specific OMS commands based on the filename specified with the `OUTFILE` keyword of the `DESTINATION` subcommand of the `OMS` command. The `FILE` keyword must be followed by an equals sign (=), and a list of quoted file specifications must be enclosed in square brackets.

### **Example**

```
OMSEND  
FILE = ['/mydir/mysavfile.sav' '/otherdir/myhtmlfile.htm'].
```

- If the specified file doesn't exist or isn't associated with a currently running OMS command, a warning is issued.
- The `FILE` keyword specification has no effect on OMS commands that use `OUTPUTSET` instead of `OUTFILE`.

## **LOG Keyword**

IF OMS logging is in effect (`OMSLOG` command), the `LOG` keyword ends logging.

### **Examples**

```
OMSEND LOG.
```

In this example, the `OMSEND` command ends logging without ending any active OMS commands.

# ***OMSINFO***

OMSINFO .

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Example***

OMSINFO .

## ***Overview***

The OMSINFO command displays a table of all active OMS commands. It has no additional specifications.

# ***OMSLOG***

```
OMSLOG FILE = 'filespec'  
  [/APPEND = [NO ]]  
           [YES]  
  [/FORMAT = [XML ]]  
           [TEXT]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Example***

```
OMSLOG FILE = '/mydir/mylog.xml'.
```

## ***Overview***

OMSLOG creates a log file in either XML or text form for subsequent OMS commands during a session.

- The log contains one line or main XML element for each destination file and contains the event name, filename, and location, the ID tag value, and a timestamp. The log also contains an entry when an OMS command is started and stopped.
- The log file remains open, and OMS activity is appended to the log, unless logging is turned off by an OMSEND command or the end of the session.
- A subsequent OMSLOG command that specifies a different log file ends logging to the file specified on the previous OMSLOG command.
- A subsequent OMSLOG file that specifies the same log file will overwrite the current log file for the default `FORMAT = XML` or in the absence of `APPEND = YES` for `FORMAT = TEXT`.
- OMS activity for any OMS commands executed *before* the first OMSLOG command in the session is not recorded in any log file.

## ***Basic Specification***

The basic specification is the command name OMSLOG followed by a FILE subcommand that specifies the log filename and location.

## ***Syntax Rules***

- The FILE subcommand is required. All other specifications are optional.
- Equals signs (=) shown in the command syntax chart and examples are required, not optional.

## **FILE Subcommand**

The `FILE` subcommand specifies the log filename and location. The subcommand name must be followed by an equals sign (=) and a file specification in quotes. If the file specification includes location information (drive, directory/folder), the location must be a valid, existing location; otherwise an error will result.

### **Example**

```
OMSLOG FILE = '/mydir/mylog.xml'.
```

## **APPEND Subcommand**

If the `FILE` subcommand specifies an existing file, by default the file is overwritten. For text format log files, you can use the `APPEND` subcommand to append new logging information to the file instead of overwriting.

### **Example**

```
OMSLOG FILE = '/mydir/mylog.txt'  
/APPEND = YES  
/FORMAT = TEXT.
```

- `APPEND = YES` is only valid with `FORMAT = TEXT`. For XML log files, the `APPEND` subcommand is ignored.
- `APPEND = YES` with `FORMAT = TEXT` will append to an existing file, even if the existing file contains XML-format log information. (An XML file *is* a text file, and OMSLOG does not differentiate based on file extension or content.)
- If the specified file does not exist, `APPEND` has no effect.

## **FORMAT Subcommand**

The `FORMAT` subcommand specifies the format of the log file. The default format is XML. You can use `FORMAT = TEXT` to write the log in simple text format.

# ONEWAY

ONEWAY is available in the Statistics Base option.

```
ONEWAY varlist BY varname

[/POLYNOMIAL=n] [/CONTRAST=coefficient list] [/CONTRAST=... ]

[/POSTHOC={ [SNK] [TUKEY] [BTUKEY] [DUNCAN] [SCHEFFE] [DUNNETT[refcat]]
            [DUNNETTL(refcat)] [DUNNETTR(refcat)] [BONFERRONI] [LSD]
            [SIDAK] [GT2] [GABRIEL] [FREGW] [QREGW] [T2] [T3] [GH] [C]
            [WALLER({100** })] } [ALPHA({0.05**})]
            {Kratio} {α }

[/RANGES={LSD } [{0.05**}] ] [/RANGES=... ]
            {DUNCAN } {α }
            {SNK }
            {TUKEYB }
            {TUKEY }
            {MODLSD }
            {SCHEFFE }

[/STATISTICS={NONE**} [DESCRIPTIVES] [EFFECTS] [HOMOGENEITY] [ALL] ]
            [WELCH] [BROWNFORSYTHE]

[/PLOT MEANS ]

[/MISSING={ANALYSIS**} [{EXCLUDE**}] ]
            {LISTWISE } {INCLUDE }

[/MATRIX = [IN({* })] [OUT({* })] [NONE] ]
            {'savfile'|'dataset'} {'savfile'|'dataset'}

[/TEMPLATE='filename']
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 19.0

- TEMPLATE subcommand introduced.

## Example

```
ONEWAY V1 BY V2.
```

## Overview

ONEWAY produces a one-way analysis of variance for an interval-level dependent variable by one numeric independent variable that defines the groups for the analysis. Other procedures that perform an analysis of variance are SUMMARIZE, UNIANOVA, and GLM (GLM is available in the Advanced Statistics option). Some tests not included in the other procedures are available as options in ONEWAY.

### **Options**

**Trend and Contrasts.** You can partition the between-groups sums of squares into linear, quadratic, cubic, and higher-order trend components using the `POLYNOMIAL` subcommand. You can specify up to 10 contrasts to be tested with the  $t$  statistic on the `CONTRAST` subcommand.

**Post Hoc Tests.** You can specify 20 different post hoc tests for comparisons of all possible pairs of group means or multiple comparisons using the `POSTHOC` subcommand.

**Statistical Display.** In addition to the default display, you can obtain means, standard deviations, and other descriptive statistics for each group using the `STATISTICS` subcommand. Fixed- and random-effects statistics as well as Levene's test for homogeneity of variance are also available.

**Matrix Input and Output.** You can write means, standard deviations, and category frequencies to a matrix data file that can be used in subsequent `ONEWAY` procedures using the `MATRIX` subcommand. You can also read matrix materials consisting of means, category frequencies, pooled variance, and degrees of freedom for the pooled variance.

**Templates.** You can specify a template, using the `TEMPLATE` subcommand, to override the default chart attribute settings on your system.

### **Basic Specification**

The basic specification is a dependent variable, keyword `BY`, and an independent variable. `ONEWAY` produces an ANOVA table displaying the between- and within-groups sums of squares, mean squares, degrees of freedom, the  $F$  ratio, and the probability of  $F$  for each dependent variable by the independent variable.

### **Subcommand Order**

- The variable list must be specified first.
- The remaining subcommands can be specified in any order.

### **Operations**

- All values of the independent variable are used. Each different value creates one category.
- If a string variable is specified as an independent or dependent variable, `ONEWAY` is not executed.

### **Limitations**

- Maximum 100 dependent variables and 1 independent variable.
- An unlimited number of categories for the independent variable. However, post hoc tests are not performed if the number of nonempty categories exceeds 50. Contrast tests are not performed if the total of empty and nonempty categories exceeds 50.
- Maximum 1 `POLYNOMIAL` subcommand.
- Maximum 1 `POSTHOC` subcommand.
- Maximum 10 `CONTRAST` subcommands.

## Example

### Testing equality of group variances

```
ONEWAY perform BY group  
/MISSING ANALYSIS  
/STATISTICS HOMOGENEITY.
```

- The procedure specifies *perform* as the dependent variable and *group* as the factor.
- `STATISTICS` requests the homogeneity of variance test.
- All other options are set to their default values.

## Analysis List

The analysis list consists of a list of dependent variables, keyword `BY`, and an independent (grouping) variable.

- Only one analysis list is allowed, and it must be specified before any of the optional subcommands.
- All variables named must be numeric.

## POLYNOMIAL Subcommand

`POLYNOMIAL` partitions the between-groups sums of squares into linear, quadratic, cubic, or higher-order trend components. The display is an expanded analysis-of-variance table that provides the degrees of freedom, sums of squares, mean square,  $F$ , and probability of  $F$  for each partition.

- The value specified on `POLYNOMIAL` indicates the highest-degree polynomial to be used.
- The polynomial value must be a positive integer less than or equal to 5 and less than the number of groups. If the polynomial specified is greater than the number of groups, the highest-degree polynomial possible is assumed.
- Only one `POLYNOMIAL` subcommand can be specified per `ONEWAY` command. If more than one is used, only the last one specified is in effect.
- `ONEWAY` computes the sums of squares for each order polynomial from weighted polynomial contrasts, using the category of the independent variable as the metric. These contrasts are orthogonal.
- With unbalanced designs and equal spacing between groups, `ONEWAY` also computes sums of squares using the unweighted polynomial contrasts. These contrasts are not orthogonal.
- The deviation sums of squares are always calculated from the weighted sums of squares (Speed, 1976).

### Example

```
ONEWAY WELL BY EDUC6  
/POLYNOMIAL=2.
```

- ONEWAY requests an analysis of variance of *WELL* by *EDUC6* with second-order (quadratic) polynomial contrasts.
- The ANOVA table is expanded to include both linear and quadratic terms.

## ***CONTRAST Subcommand***

*CONTRAST* specifies a priori contrasts to be tested by the *t* statistic. The specification on *CONTRAST* is a vector of coefficients, where each coefficient corresponds to a category of the independent variable. The Contrast Coefficients table displays the specified contrasts for each group and the Contrast Tests table displays the value of the contrast and its standard error, the *t* statistic, and the degrees of freedom and two-tailed probability of *t* for each variable. Both pooled- and separate-variance estimates are displayed.

- A contrast coefficient must be specified or implied for every group defined for the independent variable. If the number of contrast values is not equal to the number of groups, the contrast test is not performed.
- The contrast coefficients for a set should sum to 0. If they do not, a warning is issued. ONEWAY will still give an estimate of this contrast.
- Coefficients are assigned to groups defined by ascending values of the independent variable.
- The notation *n\*c* can be used to indicate that coefficient *c* is repeated *n* times.

### ***Example***

```
ONEWAY V1 BY V2
  /CONTRAST = -1 -1 1 1
  /CONTRAST = -1 0 0 1
  /CONTRAST = -1 0 .5 .5.
```

- *V2* has four levels.
- The first *CONTRAST* subcommand contrasts the combination of the first two groups with the combination of the last two groups.
- The second *CONTRAST* subcommand contrasts the first group with the last group.
- The third *CONTRAST* subcommand contrasts the first group with the combination of the third and fourth groups.

### ***Example***

```
ONEWAY V1 BY V2
  /CONTRAST = -1 1 2*0
  /CONTRAST = -1 1 0 0
  /CONTRAST = -1 1.
```

- The first two *CONTRAST* subcommands specify the same contrast coefficients for a four-group analysis. The first group is contrasted with the second group in both cases.
- The first *CONTRAST* uses the *n\*c* notation.
- The last *CONTRAST* does not work because only two coefficients are specified for four groups.



## POSTHOC Subcommand

POSTHOC produces post hoc tests for comparisons of all possible pairs of group means or multiple comparisons. In contrast to a priori analyses specified on the CONTRAST subcommand, post hoc analyses are usually not planned at the beginning of the study but are suggested by the data in the course of the study.

- Twenty post hoc tests are available. Some detect homogeneity subsets among the groups of means, some produce pairwise comparisons, and others perform both. POSTHOC produces a Multiple Comparison table showing up to 10 test categories. Nonempty group means are sorted in ascending order, with asterisks indicating significantly different groups. In addition, homogeneous subsets are calculated and displayed in the Homogeneous Subsets table if the test is designed to detect homogeneity subsets.
- When the number of valid cases in the groups varies, the harmonic mean of the group sizes is used as the sample size in the calculation for homogeneity subsets except for QREGW and FREGW. For QREGW and FREGW and tests for pairwise comparison, the sample sizes of individual groups are always used.
- You can specify only one POSTHOC subcommand per ONEWAY command. If more than one is specified, the last specification takes effect.
- You can specify one alpha value used in all POSTHOC tests using keyword ALPHA. The default is 0.05.

<b>SNK</b>	<i>Student-Newman-Keuls procedure based on the Studentized range test.</i> Used for detecting homogeneity subsets.
<b>TUKEY</b>	<i>Tukey's honestly significant difference.</i> This test uses the Studentized range statistic to make all pairwise comparisons between groups. Used for pairwise comparison and for detecting homogeneity subsets.
<b>BTUKEY</b>	<i>Tukey's b.</i> Multiple comparison procedure based on the average of Studentized range tests. Used for detecting homogeneity subsets.
<b>DUNCAN</b>	<i>Duncan's multiple comparison procedure based on the Studentized range test.</i> Used for detecting homogeneity subsets.
<b>SCHEFFE</b>	<i>Scheffé's multiple comparison t test.</i> Used for pairwise comparison and for detecting homogeneity subsets.
<b>DUNETT(refcat)</b>	<i>Dunnett's two-tailed t test.</i> Used for pairwise comparison. Each group is compared to a reference category. You can specify a reference category in parentheses. The default is the last category. This keyword must be spelled out in full.
<b>DUNETTTL(refcat)</b>	<i>Dunnett's one-tailed t test.</i> Used for pairwise comparison. This test indicates whether the mean of each group (except the reference category) is <i>smaller</i> than that of the reference category. You can specify a reference category in parentheses. The default is the last category. This keyword must be spelled out in full.
<b>DUNETTR(refcat)</b>	<i>Dunnett's one-tailed t test.</i> Used for pairwise comparison. This test indicates whether the mean of each group (except the reference category) is <i>larger</i> than that of the reference category. You can specify a reference category in parentheses. The default is the last category. This keyword must be spelled out in full.
<b>BONFERRONI</b>	<i>Bonferroni t test.</i> This test is based on Student's <i>t</i> statistic and adjusts the observed significance level for the fact that multiple comparisons are made. Used for pairwise comparison.

<b>LSD</b>	<i>Least significant difference t test.</i> Equivalent to multiple <i>t</i> tests between all pairs of groups. Used for pairwise comparison. This test does not control the overall probability of rejecting the hypotheses that some pairs of means are different, while in fact they are equal.
<b>SIDAK</b>	<i>Sidak t test.</i> Used for pairwise comparison. This test provides tighter bounds than the Bonferroni test.
<b>GT2</b>	<i>Hochberg's GT2.</i> Used for pairwise comparison and for detecting homogeneity subsets. This test is based on the Studentized maximum modulus test. Unless the cell sizes are extremely unbalanced, this test is fairly robust even for unequal variances.
<b>GABRIEL</b>	<i>Gabriel's pairwise comparisons test based on the Studentized maximum modulus test.</i> Used for pairwise comparison and for detecting homogeneity subsets.
<b>FREGW</b>	<i>Ryan-Einot-Gabriel-Welsch's multiple stepdown procedure based on an F test.</i> Used for detecting homogeneity subsets.
<b>QREGW</b>	<i>Ryan-Einot-Gabriel-Welsch's multiple stepdown procedure based on the Studentized range test.</i> Used for detecting homogeneity subsets.
<b>T2</b>	<i>Tamhane's T2.</i> Used for pairwise comparison. This test is based on a <i>t</i> test and can be applied in situations where the variances are unequal.
<b>T3</b>	<i>Tamhane's T3.</i> Used for pairwise comparison. This test is based on the Studentized maximum modulus test and can be applied in situations where the variances are unequal.
<b>GH</b>	<i>Games and Howell's pairwise comparisons test based on the Studentized range test.</i> Used for pairwise comparison. This test can be applied in situations where the variances are unequal.
<b>C</b>	<i>Dunnett's C.</i> Used for pairwise comparison. This test is based on the weighted average of Studentized ranges and can be applied in situations where the variances are unequal.
<b>WALLER(kratio)</b>	<i>Waller-Duncan t test.</i> Used for detecting homogeneity subsets. This test uses a Bayesian approach. The <i>k</i> -ratio is the Type 1/Type 2 error seriousness ratio. The default value is 100. You can specify an integer greater than 1 within parentheses.

**Example**

```
ONEWAY WELL BY EDUC6
  /POSTHOC=SNK SCHEFFE ALPHA=.01.
```

- ONEWAY requests two different post hoc tests. The first uses the Student-Newman-Keuls test and the second uses Scheffé's test. Both tests use an alpha of 0.01.

**RANGES Subcommand**

RANGES produces results for some post hoc tests. It is available only through syntax. You can always produce the same results using the POSTHOC subcommand.

- Up to 10 RANGE subcommands are allowed. The effect is cumulative. If you specify more than one alpha value for different range tests, the last specified value takes effect for all tests. The default is 0.05.
- Keyword MODLSD on the RANGE subcommand is equivalent to keyword BONFERRONI on the POSTHOC subcommand. Keyword LSDMOD is an alias for MODLSD.

## **PLOT MEANS Subcommand**

PLOT MEANS produces a chart that plots the subgroup means (the means for each group defined by values of the factor variable).

## **STATISTICS Subcommand**

By default, ONEWAY displays the ANOVA table showing between- and within-groups sums of squares, mean squares, degrees of freedom,  $F$  ratio, and probability of  $F$ . Use STATISTICS to obtain additional statistics.

<b>BROWNFORSYTHE</b>	<i>Brown-Forsythe statistic.</i> The Brown-Forsythe statistic, degrees of freedom, and the significance level are computed for each dependent variable.
<b>WELCH</b>	<i>Welch statistic.</i> The Welch statistic, degrees of freedom, and the significance level are computed for each dependent variable.
<b>DESCRIPTIVES</b>	<i>Group descriptive statistics.</i> The statistics include the number of cases, mean, standard deviation, standard error, minimum, maximum, and 95% confidence interval for each dependent variable for each group.
<b>EFFECTS</b>	<i>Fixed- and random-effects statistics.</i> The statistics include the standard deviation, standard error, and 95% confidence interval for the fixed-effects model, and the standard error, 95% confidence interval, and estimate of between-components variance for the random-effects model.
<b>HOMOGENEITY</b>	<i>Homogeneity-of-variance tests.</i> The statistics include Levene statistic, degrees of freedom, and the significance level displayed in the Test of Homogeneity-of-Variations table.
<b>NONE</b>	<i>No optional statistics.</i> This is the default.
<b>ALL</b>	<i>All statistics available for ONEWAY.</i>

## **MISSING Subcommand**

MISSING controls the treatment of missing values.

- Keywords ANALYSIS and LISTWISE are alternatives. Each can be used with INCLUDE or EXCLUDE. The default is ANALYSIS and EXCLUDE.
- A case outside of the range specified for the grouping variable is not used.

<b>ANALYSIS</b>	<i>Exclude cases with missing values on a pair-by-pair basis.</i> A case with a missing value for the dependent or grouping variable for a given analysis is not used for that analysis. This is the default.
<b>LISTWISE</b>	<i>Exclude cases with missing values listwise.</i> Cases with missing values for any variable named are excluded from all analyses.
<b>EXCLUDE</b>	<i>Exclude cases with user-missing values.</i> User-missing values are treated as missing. This is the default.
<b>INCLUDE</b>	<i>Include user-missing values.</i> User-missing values are treated as valid values.

## **MATRIX Subcommand**

MATRIX reads and writes matrix data files.

- Either `IN` or `OUT` and a matrix file in parentheses are required.
- You cannot specify both `IN` and `OUT` on the same `ONEWAY` procedure.
- Use `MATRIX=NONE` to explicitly indicate that a matrix data file is not being written or read.

**OUT ('savfile'|'dataset')**      *Write a matrix data file or dataset.* Specify either a filename, a previously declared dataset name, or an asterisk, enclosed in parentheses. Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. If you specify an asterisk (\*), the matrix data file replaces the active dataset. If you specify an asterisk or a dataset name, the file is not stored on disk unless you use `SAVE` or `XSAVE`.

**IN ('savfile'|'dataset')**      *Read a matrix data file or dataset.* Specify either a filename, dataset name created during the current session, or an asterisk enclosed in parentheses. An asterisk reads the matrix data from the active dataset. Filenames should be enclosed in quotes and are read from the working directory unless a path is included as part of the file specification.

### **Matrix Output**

- `ONEWAY` writes means, standard deviations, and frequencies to a matrix data file that can be used by subsequent `ONEWAY` procedures. For a description of the file, see [Format of the Matrix Data File](#) below.

### **Matrix Input**

- `ONEWAY` can read the matrices it writes, and it can also read matrix materials that include the means, category frequencies, pooled variance, and degrees of freedom for the pooled variance. The pooled variance has a `ROWTYPE_` value `MSE`, and the vector of degrees of freedom for the pooled variance has the `ROWTYPE_` value `DFE`.
- The dependent variables named on `ONEWAY` can be a subset of the dependent variables in the matrix data file.
- `MATRIX=IN` cannot be specified unless an active dataset has already been defined. To read an existing matrix data file at the beginning of a session, use `GET` to retrieve the matrix file and then specify `IN(*)` on `MATRIX`.

### **Format of the Matrix Data File**

- The matrix data file includes two special variables created by the program: `ROWTYPE_` and `VARNAME_`.
- `ROWTYPE_` is a short string variable with values `MEAN`, `STDDEV`, and `N`.
- `VARNAME_` is a short string variable that never has values for procedure `ONEWAY`. `VARNAME_` is included with the matrix materials so that matrices written by `ONEWAY` can be read by procedures that expect to read a `VARNAME_` variable.
- The independent variable is between variables `ROWTYPE_` and `VARNAME_`.
- The remaining variables in the matrix file are the dependent variables.

## Split Files

- When split-file processing is in effect, the first variables in the matrix data file are the split variables, followed by *ROWTYPE\_*, the independent variable, *VARNAME\_*, and the dependent variables.
- A full set of matrix materials is written for each split-file group defined by the split variable(s).
- A split variable cannot have the same variable name as any other variable written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split file must be in effect when that matrix is read by any procedure.
- Generally, matrix rows, independent variables, and dependent variables can be in any order in the matrix data file read by keyword *IN*. However, all split-file variables must precede variable *ROWTYPE\_*, and all split-group rows must be consecutive. *ONEWAY* ignores unrecognized *ROWTYPE\_* values.

## Missing Values

Missing-value treatment affects the values written to an matrix data file. When reading a matrix data file, be sure to specify a missing-value treatment on *ONEWAY* that is compatible with the treatment that was in effect when the matrix materials were generated.

## Example

```
GET FILE=GSS80.
ONEWAY WELL BY EDUC6
  /MATRIX=OUT(ONEMTX) .
```

- *ONEWAY* reads data from file *GSS80* and writes one set of matrix materials to the file *ONEMTX*.
- The active dataset is still *GSS80*. Subsequent commands are executed on *GSS80*.

## Example

```
GET FILE=GSS80.
ONEWAY WELL BY EDUC6
  /MATRIX=OUT(*) .
LIST.
```

- *ONEWAY* writes the same matrix as in the example above. However, the matrix data file replaces the active dataset. The *LIST* command is executed on the matrix file, not on the *GSS80* file.

## Example

```
GET FILE=PRSNL.
FREQUENCIES VARIABLE=AGE.
ONEWAY WELL BY EDUC6
  /MATRIX=IN(ONEMTX) .
```

- This example performs a frequencies analysis on *PRSNL* and then uses a different file for *ONEWAY*. The file is an existing matrix data file.

---

*ONEWAY*

- `MATRIX=IN` specifies the matrix data file.
- `ONEMTX` does not replace `PRSNL` as the active dataset.

**Example**

```
GET FILE=ONEMTX.  
ONEWAY WELL BY EDUC6  
/MATRIX=IN(*).
```

- The `GET` command retrieves the matrix data file `ONEMTX`.
- `MATRIX=IN` specifies an asterisk because the active dataset is the matrix data file `ONEMTX`. If `MATRIX=IN(ONEMTX)` is specified, the program issues an error message, since `ONEMTX` is already open.
- If the `GET` command is omitted, the program issues an error message.

**TEMPLATE Subcommand**

`TEMPLATE` uses an existing file as a template and applies it to the means plot (`/PLOT MEANS`), if requested, for the current `ONEWAY` command. The template overrides the default settings that are used to create the plot. Templates are created in the Chart Editor by saving an existing chart as a template.

**Example**

```
ONEWAY perform BY group  
/MISSING ANALYSIS  
/STATISTICS HOMOGENEITY  
/PLOT MEANS  
/TEMPLATE='/templates/mytemplate.sgt'.
```

**References**

Speed, M. F. 1976. Response curves in the one way classification with unequal numbers of observations per cell. In: *Proceedings of the Statistical Computing Section*, Alexandria, VA: AmericanStatistical Association, 270–272.

# OPTIMAL BINNING

OPTIMAL BINNING is available in the Data Preparation option.

```
OPTIMAL BINNING
/VARIABLES [GUIDE = variable] BIN = varlist [SAVE = {NO**
                                                {YES [(INTO = new varlist)]}]
[/CRITERIA
  [PREPROCESS = {EQUALFREQ**[(BINS = {1000**})]}]
                {n          }
                {NONE          }
  [METHOD = {MDLP**
            {EQUALFREQ [(BINS = {10**})]}
            {n          }
            }
  [LOWEREND = {UNBOUNDED**}] [UPPEREND = {UNBOUNDED**}]
              {OBSERVED  }           {OBSERVED  }
  [LOWERLIMIT = {INCLUSIVE**}]
                {EXCLUSIVE  }
  [FORCEMERGE = {0**  }]
                {value}
[/MISSING [SCOPE = {PAIRWISE**}]
          {LISTWISE  }
[/OUTFILE RULES = filespec]
[/PRINT [ENDPOINTS**] [DESCRIPTIVES] [ENTROPY] [NONE]]
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 15.0

- Command introduced.

## Example

```
OPTIMAL BINNING
/VARIABLES GUIDE = guide-variable BIN = binning-input-variable
```

## Overview

The OPTIMAL BINNING procedure discretizes one or more scale variables (referred to henceforth as **binning input** variables) by distributing the values of each variable into bins. Bins can then be used instead of the original data values of the binning input variables for further analysis. OPTIMAL BINNING is useful for reducing the number of distinct values in the given binning input variables.

**Options**

**Methods.** The OPTIMAL BINNING procedure offers the following methods of discretizing binning input variables.

- Unsupervised binning via the equal frequency algorithm discretizes the binning input variables. A guide variable is not required.
- Supervised binning via the MDLP (Minimal Description Length Principle) algorithm discretizes the binning input variables without any preprocessing. It is suitable for datasets with a small number of cases. A guide variable is required.

**Output.** The OPTIMAL BINNING procedure displays every binning input variable's end point set in pivot table output and offers an option for suppressing this output. In addition, the procedure can save new binned variables corresponding to the binning input variables and can save a command syntax file with commands corresponding to the binning rules.

**Basic Specification**

The basic specification is the OPTIMAL BINNING command and a VARIABLES subcommand. VARIABLES provides the binning input variables and, if applicable, the guide variable.

- For unsupervised binning via the equal frequency algorithm, a guide variable is not required.
- For supervised binning via the MDLP algorithm and hybrid binning, a guide variable must be specified.

**Syntax Rules**

- When a supervised binning method is used, a guide variable must be specified on the VARIABLES subcommand.
- Subcommands may be specified only once.
- An error occurs if a variable or keyword is specified more than once within a subcommand.
- Parentheses, slashes, and equals signs shown in the syntax chart are required.
- Empty subcommands are not honored.
- The command name, subcommand names, and keywords must be spelled in full.

**Case Frequency**

- If a WEIGHT variable is specified, then its values are used as frequency weights by the OPTIMAL BINNING procedure.
- Weight values are rounded to the nearest whole numbers before use. For example, 0.5 is rounded to 1, and 2.4 is rounded to 2.
- The WEIGHT variable may not be specified on any subcommand in the OPTIMAL BINNING procedure.
- Cases with missing weights or weights less than 0.5 are not used in the analyses.



### Limitations

The number of distinct values in a guide variable should be less than or equal to 256, irrespective of the platform on which IBM® SPSS® Statistics is running. If the number is greater than 256, this results in an error.

## Examples

```
* Optimal Binning.
OPTIMAL BINNING
/VARIABLES GUIDE=default BIN=age employ address income debtinc creddebt
othdebt SAVE=YES (INTO=age_bin employ_bin address_bin income_bin debtinc_bin
creddebt_bin othdebt_bin)
/CRITERIA METHOD=MDLP
PREPROCESS=EQUALFREQ (BINS=1000)
FORCEMERGE=0
LOWERLIMIT=INCLUSIVE
LOWEREND=UNBOUNDED
UPPEREND=UNBOUNDED
/MISSING SCOPE=PAIRWISE
/OUTFILE RULES='/bankloan_binning-rules.sps'
/PRINT ENDPOINTS DESCRIPTIVES ENTROPY.
```

- The procedure will discretize the binning input variables *age*, *employ*, *address*, *income*, *debtinc*, *creddebt*, and *othdebt* using MDLP binning with the guide variable *default*.
- The discretized values for these variables will be stored in the new variables *age\_bin*, *employ\_bin*, *address\_bin*, *income\_bin*, *debtinc\_bin*, *creddebt\_bin*, and *othdebt\_bin*.
- If a binning input variable has more than 1000 distinct values, then the equal frequency method will reduce the number to 1000 before performing MDLP binning.
- Command syntax representing the binning rules is saved to the file */bankloan\_binning-rules.sps*.
- Bin endpoints, descriptive statistics, and model entropy values are requested for binning input variables.
- Other binning criteria are set to their default values.

## VARIABLES Subcommand

The `VARIABLES` subcommand specifies the guide variable (if applicable) and one or more binning input variables. It can also be used to save new variables containing the binned values.

### GUIDE=variable

*Guide variable.* The bins formed by supervised binning methods are “optimal” with respect to the specified guide variable. You must specify a guide variable to perform MDLP (`CRITERIA METHOD = MDLP`) or the hybrid method (`CRITERIA PREPROCESS = EQUALFREQ METHOD = MDLP`). This option is silently ignored if it is specified when the equal frequency method (`CRITERIA METHOD = EQUALFREQ`) is in effect. The guide variable may be numeric or string.

### BIN=varlist

*Binning input variable list.* These are the variables to be binned. The variable list must include at least one variable. Binning input variables must be numeric.

`SAVE = NO | YES (INTO = new varlist)`

*Create new variables containing binned values.* By default, the procedure does not create any new variables (NO). If YES is specified, variables containing the binned values are saved to the active dataset. Optionally, specify the names of the new variables using the INTO keyword.

The number of variables specified on the INTO list must equal the number of variables on the BIN list. All specified names must be valid variable names. Violation of either of these rules results in an error.

If INTO is omitted, new variable names are created by concatenating the guide variable name (if applicable) and an underscore ‘\_’, followed by the binning input variable name and an underscore, followed by ‘bin’. For example, /VARIABLES GUIDE=E BIN=F G SAVE=YES will generate two new variables: *E\_F\_bin* and *E\_G\_bin*.

## **CRITERIA Subcommand**

The CRITERIA subcommand specifies bin creation options.

### **PREPROCESS=EQUALFREQ(BINS=n) | NONE**

*Preprocessing method when MDLP binning is used.* PREPROCESS = EQUALFREQ creates preliminary bins using the equal frequency method before performing MDLP binning. These preliminary bins—rather than the original data values of the binning input variables—are input to the MDLP binning method.

EQUALFREQ may be followed by parentheses containing the BINS keyword, an equals sign, and an integer greater than 1. The BINS value serves as a preprocessing threshold and specifies the number of bins to create. The default value is EQUALFREQ (BINS = 1000).

If the number of distinct values in a binning input variable is greater than the BINS value, then the number of bins created is no more than the BINS value. Otherwise, no preprocessing is done for the input variable.

NONE requests no preprocessing.

### **METHOD=MDLP | EQUALFREQ(BINS=n)**

*Binning method.* The MDLP option performs supervised binning via the MDLP algorithm. If METHOD = MDLP is specified, then a guide variable must be specified on the VARIABLES subcommand.

Alternatively, METHOD = EQUALFREQ performs unsupervised binning via the equal frequency algorithm. EQUALFREQ may be followed by parentheses containing the BINS keyword, an equals sign, and an integer greater than 1. The BINS value specifies the number of bins to create. The default value of the BINS argument is 10.

If the number of distinct values in a binning input variable is greater than the BINS value, then the number of bins created is no more than the BINS value. Otherwise, BINS gives an upper bound on the number of bins created. Thus, for example, if BINS = 10 is specified but a binning input variable has at most 10 distinct values, then the number of bins created will equal the number of distinct values in the input variable.

If EQUALFREQ is specified, then the VARIABLES subcommand GUIDE keyword and the CRITERIA subcommand PREPROCESS keyword are silently ignored.

The default METHOD option depends on the presence of a GUIDE specification on the VARIABLES subcommand. If GUIDE is specified, then METHOD = MDLP is the default. If GUIDE is not specified, then METHOD = EQUALFREQ is the default.

### **LOWEREND = UNBOUNDED | OBSERVED**

*Specifies how the minimum end point for each binning input variable is defined.* Valid option values are UNBOUNDED or OBSERVED. If UNBOUNDED, then the minimum end point extends to negative infinity. If OBSERVED, then the minimum observed data value is used.

**UPPEREND = UNBOUNDED | OBSERVED**

*Specifies how the maximum end point for each binning input variable is defined.* Valid option values are UNBOUNDED or OBSERVED. If UNBOUNDED, then the maximum end point extends to positive infinity. If OBSERVED, then the maximum of the observed data is used.

**LOWERLIMIT = INCLUSIVE | EXCLUSIVE**

*Specifies how the lower limit of an interval is defined.* Valid option values are INCLUSIVE or EXCLUSIVE. Suppose the start and end points of an interval are  $p$  and  $q$ , respectively. If LOWERLIMIT = INCLUSIVE, then the interval contains values greater than or equal to  $p$  but less than  $q$ . If LOWERLIMIT = EXCLUSIVE, then the interval contains values greater than  $p$  and less than or equal to  $q$ .

**FORCEMERGE = value**

*Small bins threshold.* Occasionally, the procedure may produce bins with very few cases. The following strategy deletes these pseudo cut points:

- ▶ For a given variable, suppose that the algorithm found  $n_{\text{final}}$  cut points, and thus  $n_{\text{final}}+1$  bins. For bins  $i = 2, \dots, n_{\text{final}}$  (the second lowest-valued bin through the second highest-valued bin), compute

$$\frac{\text{sizeof}(b_i)}{\min(\text{sizeof}(b_{i-1}), \text{sizeof}(b_{i+1}))}$$

where  $\text{sizeof}(b)$  is the number of cases in the bin.

- ▶ When this value is less than the specified merging threshold,  $b_i$  is considered sparsely populated and is merged with  $b_{i-1}$  or  $b_{i+1}$ , whichever has the lower class information entropy.

The procedure makes a single pass through the bins.

The default value of FORCEMERGE is 0; by default, forced merging of very small bins is not performed.

## **MISSING Subcommand**

The MISSING subcommand specifies whether missing values are handled using listwise or pairwise deletion.

- User-missing values are always treated as invalid. When recoding the original binning input variable values into a new variable, user-missing values are converted into system-missing values.

**SCOPE = PAIRWISE | LISTWISE**

*Missing value handling method.* LISTWISE provides a consistent case base. It operates across all variables specified on the VARIABLES subcommand. If any variable is missing for a case, then the entire case is excluded.

PAIRWISE makes use of as many valid values as possible. When METHOD = MDLP, it operates on each guide and binning input variable pair. The procedure will make use of all cases with nonmissing values on the guide and binning input variable. When METHOD = EQUALFREQ, it uses all cases with nonmissing values for each binning input variable. PAIRWISE is the default.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand writes syntax to an external command syntax file.

### **RULES=filespec**

*Rules file specification.* The procedure can generate command syntax that can be used to bin other datasets. The recoding rules are based on the end points determined by the binning algorithm. Specify an external file to contain the saved syntax.

Note that saved variables (see the `SAVE` keyword in the [VARIABLES subcommand](#)) are generated using end points exactly as computed by the algorithm, while the bins created via saved syntax rules use end points converted to and from a decimal representation. Conversion errors in this process can, in certain cases, cause the end points read from syntax to differ from the original ones. The syntax precision of end points is 17 digits.

## ***PRINT Subcommand***

The `PRINT` subcommand controls the display of the output results. If the `PRINT` subcommand is not specified, then the default output is the end point set for each binning input variable.

### **ENDPOINTS**

*Display the binning interval end points for each input variable.* This is the default output.

### **DESCRIPTIVES**

*Display descriptive information for all binning input variables.* For each binning input variable, this option displays the number of cases with valid values, the number of cases with missing values, the number of distinct valid values, and the minimum and maximum values. For the guide variable, this option displays the class distribution for each related binning input variable.

### **ENTROPY**

*Display the model entropy for each binning input variable interval when MDLP binning is used.* The `ENTROPY` keyword is ignored with a warning if `METHOD = EQUALFREQ` is specified or implied on the `CRITERIA` subcommand.

### **NONE**

*Suppress all displayed output except the notes table and any warnings.* Specifying `NONE` with any other keywords results in an error.

# ORTHOPLAN

ORTHOPLAN is available in the Conjoint option.

```
ORTHOPLAN
[/FACTORS=varlist ['labels'] (values ['labels'])...]

[{/REPLACE      }]
  {/OUTFILE='savfile'|'dataset'}

[/MINIMUM=value]

[/HOLDOUT=value]  [/MIXHOLD={YES}]
                  {NO }
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example:**

```
ORTHOPLAN FACTORS=SPEED 'Highest possible speed'
          (70 '70 mph' 100 '100 mph' 130 '130mph')
WARRANTY 'Length of warranty'
          ('1 year' '3 year' '5 year')
SEATS (2, 4)
/MINIMUM=9 /HOLDOUT=6.
```

## **Overview**

ORTHOPLAN generates an orthogonal main-effects plan for a full-concept conjoint analysis. ORTHOPLAN can append or replace an existing active dataset, or it can build an active dataset (if one does not already exist). The generated plan can be listed in full-concept profile, or card, format using PLANCARDS. The file that is created by ORTHOPLAN can be used as the plan file for CONJOINT.

### **Options**

**Number of Cases.** You can specify the minimum number of cases to be generated in the plan.

**Holdout and Simulation Cases.** In addition to the experimental main-effects cases, you can generate a specified number of holdout cases and identify input data as simulation cases.

### **Basic Specification**

- The basic specification is ORTHOPLAN followed by FACTORS, a variable list, and a value list in parentheses. ORTHOPLAN will generate cases in the active dataset, with each case representing a profile in the conjoint experimental plan and consisting of a new combination of the factor values. By default, the smallest possible orthogonal plan is generated.
- If you are appending to an existing active dataset that has previously defined values, the FACTORS subcommand is optional.

**Subcommand Order**

- Subcommands can be named in any order.

**Operations**

- ORTHOPLAN builds an active dataset (if one does not already exist) by using the variable and value information on the FACTORS subcommand.
- When ORTHOPLAN appends to an active dataset and FACTORS is not used, the factor levels (values) must be defined on a previous ORTHOPLAN or VALUE LABELS command.
- New variables *STATUS\_* and *CARD\_* are created and added to the active dataset by ORTHOPLAN if they do not already exist. *STATUS\_*=0 for experimental cases, 1 for holdout cases, and 2 for simulation cases. Holdout cases are judged by the subjects but are not used when CONJOINT estimates utilities. Instead, the cases are used as a check on the validity of the estimated utilities. Simulation cases are entered by the user. They are factor-level combinations that are not rated by the subjects but are estimated by CONJOINT based on the ratings of the experimental cases. *CARD\_* contains the case identification numbers in the generated plan.
- Duplication between experimental cases and simulation cases is reported.
- If a user-entered experimental case (*STATUS\_*=0) is duplicated by ORTHOPLAN, only one copy of the case is kept.
- Occasionally, ORTHOPLAN may generate duplicate experimental cases. One way to handle these duplicates is to edit or delete them, in which case the plan is no longer orthogonal. Alternatively, you can try running ORTHOPLAN again. With a different *seed*, ORTHOPLAN might produce a plan without duplicates. See the SEED subcommand on SET for more information about the random seed generator.
- The SPLIT FILE and WEIGHT commands are ignored by ORTHOPLAN.

**Limitations**

- Missing data are not allowed.
- A maximum of 9 levels can be specified for each factor.
- The total number of levels for all factors cannot exceed 90.
- A maximum of 81 cases can be generated by ORTHOPLAN.

**Examples**

```
ORTHOPLAN FACTORS=SPEED 'Highest possible speed'
          (70 '70 mph' 100 '100 mph' 130 '130mph')
          WARRANTY 'Length of warranty'
          ('1 year' '3 year' '5 year')
          SEATS (2, 4) /MINIMUM=9 /HOLDOUT=6 /OUTFILE='CARPLAN.SAV'.
```

- The FACTORS subcommand defines the factors and levels to be used in building the file. Labels for some of the factors and some of the levels of each factor are also supplied.
- The MINIMUM subcommand specifies that the orthogonal plan should contain at least nine full-concept cases.

- `HOLDOUT` specifies that six holdout cases should be generated. A new variable, `STATUS_`, is created by `ORTHOPLAN` to distinguish these holdout cases from the regular experimental cases. Another variable, `CARD_`, is created to assign identification numbers to the plan cases.
- The `OUTFILE` subcommand saves the plan that is generated by `ORTHOPLAN` as a data file so that it can be used at a later date with `CONJOINT`.

**Example: Appending Plan to the Working File**

```
DATA LIST FREE /SPEED WARRANTY SEATS.
VALUE LABELS speed 70 '70 mph' 100 '100 mph' 130 '130 mph'
/WARRANTY 1 '1 year' 3 '3 year' 5 '5 year'
/SEATS 2 '2 seats' 4 '4 seats'.
BEGIN DATA
130 5 2
130 1 4
END DATA.
ORTHOPLAN
/OUTFILE='CARPLAN.SAV' .
```

- In this example, `ORTHOPLAN` appends the plan to the active dataset and uses the variables and values that were previously defined in the active dataset as the factors and levels of the plan.
- The data between `BEGIN DATA` and `END DATA` are assumed to be simulation cases and are assigned a value of 2 on the newly created `STATUS_` variable.
- The `OUTFILE` subcommand saves the plan that is generated by `ORTHOPLAN` as a data file so that it can be used at a later date with `CONJOINT`.

## **FACTORS Subcommand**

`FACTORS` specifies the variables to be used as factors and the values to be used as levels in the plan.

- `FACTORS` is required for building a new active dataset or replacing an existing one. `FACTORS` is optional for appending to an existing file.
- The keyword `FACTORS` is followed by a variable list, an optional label for each variable, a list of values for each variable, and optional value labels.
- The list of values and the value labels are enclosed in parentheses. Values can be numeric or they can be strings enclosed in apostrophes.
- The optional variable and value labels are enclosed in apostrophes.
- If the `FACTORS` subcommand is not used, every variable in the active dataset (other than `STATUS_` and `CARD_`) is used as a factor, and level information is obtained from the value labels that are defined in the active dataset. `ORTHOPLAN` *must* be able to find value information either from a `FACTORS` subcommand or from a `VALUE LABELS` command. (See the `VALUE LABELS` command for more information.)

**Example**

```
ORTHOPLAN FACTORS=SPEED 'Highest possible speed'
(70 '70 mph' 100 '100 mph' 130 '130mph')
WARRANTY 'Length of warranty'
(1 '1 year' 3 '3 year' 5 '5 year')
SEATS 'Number of seats' (2 '2 seats' 4 '4 seats')
EXCOLOR 'Exterior color'
```

```
INCOLOR 'Interior color' ('RED' 'BLUE' 'SILVER').
```

- *SPEED*, *WARRANTY*, *SEATS*, *EXCOLOR*, and *INCOLOR* are specified as the factors. They are given the labels *Highest possible speed*, *Length of warranty*, *Number of seats*, *Exterior color*, and *Interior color*.
- Following each factor and its label are the list of values and the value labels in parentheses. Note that the values for two of the factors, *EXCOLOR* and *INCOLOR*, are the same and thus need to be specified only once after both factors are listed.

## **REPLACE Subcommand**

REPLACE can be specified to indicate that the active dataset, if present, should be replaced by the generated plan. There is no further specification after the REPLACE keyword.

- By default, the active dataset is not replaced. Any new variables that are specified on a FACTORS subcommand plus the variables *STATUS\_* and *CARD\_* are appended to the active dataset.
- REPLACE should be used when the current active dataset has nothing to do with the plan file to be built. The active dataset will be replaced with one that has variables *STATUS\_*, *CARD\_*, and any other variables that are specified on the FACTORS subcommand.
- If REPLACE is specified, the FACTORS subcommand is required.

## **OUTFILE Subcommand**

OUTFILE saves the orthogonal design to an external IBM® SPSS® Statistics data file or previously declared dataset. The only specification is a name for the output file. This specification can be a filename or a previously declared dataset name. Filenames should be enclosed in quotation marks and are stored in the working directory unless a path is included as part of the file specification. Datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files.

- By default, a new data file is not created. Any new variables that are specified on a FACTORS subcommand plus the variables *STATUS\_* and *CARD\_* are appended to the active dataset.
- The output data file contains variables *STATUS\_*, *CARD\_*, and any other variables that are specified on the FACTORS subcommand.
- The file that is created by OUTFILE can be used by other syntax commands, such as PLANCARDS and CONJOINT.
- If both OUTFILE and REPLACE are specified, REPLACE is ignored.

## **MINIMUM Subcommand**

MINIMUM specifies a minimum number of cases for the plan.

- By default, the minimum number of cases necessary for the orthogonal plan is generated.



- `MINIMUM` is followed by a positive integer that is less than or equal to the total number of cases that can be formed from all possible combinations of the factor levels.
- If `ORTHOPLAN` cannot generate at least the number of cases requested on `MINIMUM`, it will generate the largest number it can that fits the specified factors and levels.

## ***HOLDOUT Subcommand***

`HOLDOUT` creates holdout cases in addition to the regular plan cases. Holdout cases are judged by the subjects but are not used when `CONJOINT` estimates utilities.

- If `HOLDOUT` is not specified, no holdout cases are produced.
- `HOLDOUT` is followed by a positive integer that is less than or equal to the total number of cases that can be formed from all possible combinations of factor levels.
- Holdout cases are generated from another random plan, not the main-effects experimental plan. The holdout cases will not duplicate the experimental cases or each other.
- The experimental and holdout cases will be randomly mixed in the generated plan or the holdout cases will be listed after the experimental cases, depending on subcommand `MIXHOLD`. The value of `STATUS_` for holdout cases is 1. Any simulation cases will follow the experimental and holdout cases.

## ***MIXHOLD Subcommand***

`MIXHOLD` indicates whether holdout cases should be randomly mixed with the experimental cases or should appear separately after the experimental plan in the file.

- If `MIXHOLD` is not specified, the default is `NO`, meaning holdout cases will appear after the experimental cases in the file.
- `MIXHOLD` followed by keyword `YES` requests that the holdout cases be randomly mixed with the experimental cases.
- `MIXHOLD` specified without a `HOLDOUT` subcommand has no effect.

# OUTPUT ACTIVATE

```
OUTPUT ACTIVATE [NAME=]name
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 15.0

- Command introduced.

## **Example**

```
GET FILE='/examples/data/SalaryData2005.sav'.  
DESCRIPTIVES salbegin salary.  
OUTPUT NAME alleduclevels.
```

```
TEMPORARY.  
SELECT IF (educ>12).  
OUTPUT NEW NAME=over12.  
DESCRIPTIVES salbegin salary.
```

```
GET FILE='/examples/data/SalaryData2000.sav'.  
TEMPORARY.  
SELECT IF (educ>12).  
DESCRIPTIVES salbegin salary.
```

```
OUTPUT ACTIVATE alleduclevels.  
DESCRIPTIVES salbegin salary.
```

## **Overview**

The OUTPUT commands (OUTPUT NEW, OUTPUT NAME, OUTPUT ACTIVATE, OUTPUT OPEN, OUTPUT SAVE, OUTPUT CLOSE) provide the ability to programmatically manage one or many output documents. These functions allow you to:

- Save an output document through syntax.
- Programmatically partition output into separate output documents (for example, results for males in one output document and results for females in a separate one).
- Work with multiple open output documents in a given session, selectively appending new results to the appropriate document.

The OUTPUT ACTIVATE command activates an open output document, making it the designated output document. Subsequent procedure output is directed to this output document until the document is closed or another output document is created, opened, or activated.

### **Basic Specification**

The basic specification for `OUTPUT ACTIVATE` is the command name followed by the name of an open output document. This is the name assigned by a previous `OUTPUT NAME`, `OUTPUT OPEN`, `OUTPUT NEW` command, or automatically assigned when the document is created or opened in the current session; it is not the file name associated with a saved output document. In the Viewer window title bar, the name is enclosed in square brackets. The `NAME` keyword is optional, but if it is used it must be followed by an equals sign.

### **Operations**

- The window containing the activated document becomes the designated output window in the user interface.
- An error occurs, but processing continues, if the named output document does not exist. Output continues to be directed to the last designated output document.

### **Example**

```
GET FILE='/examples/data/SurveyData.sav'.

TEMPORARY.
SELECT IF (Sex='Male').
FREQUENCIES VARIABLES=ALL.
OUTPUT NAME males.

TEMPORARY.
SELECT IF (Sex='Female').
OUTPUT NEW NAME=females.
FREQUENCIES VARIABLES=ALL.

GET FILE='/examples/data/Preference.sav'.

TEMPORARY.
SELECT IF (Sex='Female').
DESCRIPTIVES VARIABLES=product1 product2 product3.

TEMPORARY.
SELECT IF (Sex='Male').
OUTPUT ACTIVATE males.
DESCRIPTIVES VARIABLES=product1 product2 product3.

OUTPUT SAVE NAME=males OUTFILE='/examples/output/Males.spv'.
OUTPUT SAVE NAME=females OUTFILE='/examples/output/Females.spv'.
```

- The first `GET` command loads survey data for males and females.
- `FREQUENCIES` output for male respondents is written to the designated output document. The `OUTPUT NAME` command is used to assign the name *males* to the designated output document.
- `FREQUENCIES` output for females is written to a new output document named *females*.
- The second `GET` command loads preferences data for males and females.
- After the second `GET` command, the output document named *females* is still the designated output document. Descriptive statistics for females are appended to this output document.

- `OUTPUT ACTIVATE males` activates the output document named *males*. Descriptive statistics for males are appended to this output document.
- The two open output documents are saved to separate files. Because the operation of saving an output document does not close it, both documents remain open. The output document named *males* remains the designated output document.

# OUTPUT CLOSE

```
OUTPUT CLOSE [NAME=]{name}
                { *   }
                {ALL }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 15.0

- Command introduced.

## **Example**

```
GET FILE='/examples/data/Males.sav'.
FREQUENCIES VARIABLES=ALL.
OUTPUT SAVE OUTFILE='/examples/output/Males.spv'.
OUTPUT CLOSE *.
```

```
GET FILE='/examples/data/Females.sav'.
FREQUENCIES VARIABLES=ALL.
```

## **Overview**

The OUTPUT commands (OUTPUT NEW, OUTPUT NAME, OUTPUT ACTIVATE, OUTPUT OPEN, OUTPUT SAVE, OUTPUT CLOSE) provide the ability to programmatically manage one or many output documents. These functions allow you to:

- Save an output document through syntax.
- Programmatically partition output into separate output documents (for example, results for males in one output document and results for females in a separate one).
- Work with multiple open output documents in a given session, selectively appending new results to the appropriate document.

The OUTPUT CLOSE command closes one or all open output documents.

## **Basic Specification**

The only specification for OUTPUT CLOSE is the command name followed by the name of an open output document, an asterisk (\*), or the keyword ALL. The NAME keyword is optional, but if it is used it must be followed by an equals sign.

## **Operations**

- If a name is provided, the specified output document is closed and the association with that name is broken.

- If an asterisk (\*) is specified, the designated output document is closed. If the designated output document has a name, the association with that name is broken.
- If ALL is specified, all open output documents are closed and all associations of names with output documents are broken.
- Output documents are not saved automatically when they are closed. Use `OUTPUT SAVE` to save the contents of an output document.
- `OUTPUT CLOSE` is ignored if you specify a nonexistent document.

**Example**

```
GET FILE='/examples/data/Males.sav'.
FREQUENCIES VARIABLES=ALL.
OUTPUT SAVE OUTFILE='/examples/output/Males.spv'.
OUTPUT CLOSE *.
```

```
GET FILE='/examples/data/Females.sav'.
FREQUENCIES VARIABLES=ALL.
```

- `FREQUENCIES` produces summary statistics for each variable. Procedure output is added to the designated output document (one is created automatically if no output document is currently open).
- `OUTPUT SAVE` writes contents of the designated output document to the file */examples/output/Males.spv*.
- `OUTPUT CLOSE` closes the designated output document.
- Output from the second `FREQUENCIES` command is written to a new output document, which was created automatically when the previous designated output document was closed. If `OUTPUT CLOSE` had not been issued, output for females would have been directed to the output document that contained summaries for males.

# ***OUTPUT DISPLAY***

OUTPUT DISPLAY

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Release History***

Release 15.0

- Command introduced.

## ***Example***

OUTPUT DISPLAY.

## ***Overview***

The OUTPUT commands (OUTPUT NEW, OUTPUT NAME, OUTPUT ACTIVATE, OUTPUT OPEN, OUTPUT SAVE, OUTPUT CLOSE) provide the ability to programmatically manage one or many output documents. These functions allow you to:

- Save an output document through syntax.
- Programmatically partition output into separate output documents (for example, results for males in one output document and results for females in a separate one).
- Work with multiple open output documents in a given session, selectively appending new results to the appropriate document.

The OUTPUT DISPLAY command displays a list of open output documents and identifies the one that is currently active, and referred to as the designated output document. The only specification is the command name OUTPUT DISPLAY.

# OUTPUT EXPORT

```
OUTPUT EXPORT [NAME = {* }]  
              {name}  
  
[/CONTENTS [EXPORT = {ALL** }]]  
            {VISIBLE}  
[LAYERS = {PRINTSETTING**}]  
          {ALL }  
          {VISIBLE }  
[MODELVIEWS = {PRINTSETTING**}]  
          {ALL }  
          {VISIBLE }  
  
[/DOC DOCUMENTFILE = 'filespec']  
[NOTESCAPTIONS = {YES**}]  
              {NO }  
[WIDETABLES = {WRAP**}]  
              {SHRINK }  
              {ASIS }  
[PAGESIZE = {INCHES(width,height)}]  
            {MM(width,height) }  
[TOPMARGIN = {INCHES(n)}]  
            {MM(n) }  
[BOTTOMMARGIN = {INCHES(n)}]  
            {MM(n) }  
[LEFTMARGIN = {INCHES(n)}]  
            {MM(n) }  
[RIGHTMARGIN = {INCHES(n)}]  
            {MM(n) }  
  
[/HTML DOCUMENTFILE = 'filespec']  
[NOTESCAPTIONS = {YES**}]  
              {NO }  
[IMAGEFORMAT = {JPG**}]  
              {BMP }  
              {EMF }  
              {EPS }  
              {PNG }  
              {TIF }  
[/PDF DOCUMENTFILE = 'filespec']  
[EMBEDBOOKMARKS = {YES**}]  
              {NO }  
[EMBEDFONTS = {YES**}]  
              {NO }  
  
[/PPT DOCUMENTFILE = 'filespec']  
[NOTESCAPTIONS = {YES**}]  
              {NO }  
[SLIDETITLES = {YES**}]  
              {NO }  
[WIDETABLES = {WRAP**}]  
              {SHRINK }  
              {ASIS }  
[SLIDESIZE = {INCHES(width,height)}]  
            {MM(width,height) }  
[TOPMARGIN = {INCHES(n)}]  
            {MM(n) }  
[BOTTOMMARGIN = {INCHES(n)}]  
            {MM(n) }  
[LEFTMARGIN = {INCHES(n)}]  
            {MM(n) }  
[RIGHTMARGIN = {INCHES(n)}]  
            {MM(n) }  
  
[/TEXT DOCUMENTFILE = 'filespec']  
[ENCODING = {LOCALE**}]  
          {UTF8 }  
          {UTF16 }
```



```

[NOTESCAPTIONS = {YES**}]
    {NO }
[COLUMNWIDTH = {AUTOFIT**}]
    {width }
[SEPARATOR = {SPACE**}]
    {TAB }
[ROWBORDER = {'-'** }]]
    {'character' }
[COLUMNBORDER = {'|'** }]]
    {'character' }
[IMAGEFORMAT = {JPG**}]
    {BMP }
    {EMF }
    {EPS }
    {PNG }
    {TIF }

[/XLS DOCUMENTFILE = 'filespec']
[OPERATION = {CREATEFILE**}]
    {CREATESHEET }
    {MODIFYSHEET }
[SHEET = {'sheetname' }]
[LOCATION = {LASTCOLUMN** }]]
    {LASTROW }
    {STARTCELL('cellref')}
[NOTESCAPTIONS = {YES**}]
    {NO }

[/BMP IMAGEROOT = 'rootname']
[PERCENTSIZE = {100**}]
    {n }
[COMPRESS = {NO**}]
    {YES }

[/EMF IMAGEROOT = 'rootname']
[PERCENTSIZE = {100**}]
    {n }

[/EPS IMAGEROOT = 'rootname']
[WIDTH = {PERCENT**({100**})}]
    {n }
    {CENTIMETERS(n) }
    {INCHES(n) }
    {POINTS(n) }
[PREVIEW = {YES**}]
    {NO }
[FONTHANDLING = {REFERENCES**}]
    {CURVES }
[/JPG IMAGEROOT = 'rootname']
[PERCENTSIZE = {100**}]
    {n }
[GRAYSCALE = {NO**}]
    {YES }

[/PNG IMAGEROOT = 'rootname']
[PERCENTSIZE = {100**}]
    {n }
[DEPTH = {SCREEN** }]]
    {BLACKWHITE}
    {GRAYS256 }
    {COLORS16 }
    {COLORS256 }
    {TRUE24BIT }
    {TRUE32BIT }

[/TIF IMAGEROOT = 'rootname']
[PERCENTSIZE = {100**}]
    {n }

```

\*\* Default if the subcommand or keyword is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 17.0

- Command introduced.

### **Example**

```
OUTPUT EXPORT  
  /DOC DOCUMENTFILE=' /mydocs/myresults.doc' .
```

## **Overview**

The `OUTPUT EXPORT` command exports output from an open Viewer document to an external file format. By default, the contents of the designated Viewer document are exported, but a different Viewer document can be specified by name.

### **Options**

Options include:

- Export output in Word, Excel, HTML, PDF, PowerPoint, and text format.
- Export all items or only visible items.
- Export graphics only.

### **Basic Specification**

The basic specification is the `OUTPUT EXPORT` command, followed by one of the document type subcommands with the `DOCUMENTFILE` keyword or by one of the image type subcommands with the `IMAGEROOT` keyword.

### **Syntax Rules**

- The command name, subcommand names, and all keywords must be spelled in full. (Abbreviation is not allowed.)
- Each subcommand must be preceded by a forward slash.
- Equals signs shown in the syntax diagram are required elements.
- Subcommands can be specified in any order.
- Each subcommand can be specified only once.
- Each subcommand must be followed by at least one keyword specification.
- Within a subcommand, each keyword can be specified only once.

### **Operations**

By default, the contents of the **designated** Viewer document are exported. The designated Viewer document is the one most recently opened (by `OUTPUT NEW` or `OUTPUT OPEN`, or from the menus and dialogs), or explicitly designated (by `OUTPUT ACTIVATE` or by selecting Designate Window from the Utilities menu).

To export an open Viewer document other than the designated one, use the `NAME` keyword. [For more information, see the topic NAME Keyword on p. 1470.](#)

### **Limitations**

- Export of selected Viewer items is not supported in `OUTPUT EXPORT`. You can only export all items or all visible items.
- Exporting to PowerPoint (\*.ppt) format is supported only on Windows operating systems that have PowerPoint installed.
- EMF graphic format is available only on Windows operating systems.
- `OUTPUT EXPORT` is not supported in the IBM® SPSS® Statistics Batch Facility (available with SPSS Statistics Server). To export output created in the Batch Facility, use [OMS](#).

## **Examples**

### **Export to Word with Minimum Specifications**

```
OUTPUT EXPORT  
  /DOC DOCUMENTFILE='/mydocs/myresults.doc'.
```

All output in the designated Viewer document is exported to the specified file in Word (RTF) format. If the file already exists, the contents are overwritten.

### **Export Named Viewer Document to New Sheet in Existing Excel Document**

```
OUTPUT EXPORT NAME=Region1Results  
  /XLS DOCUMENTFILE='/mydocs/ResultsByRegion.xls'  
  OPERATION=CREATESHEET SHEET='Region 1'.
```

- Contents of the Viewer document named *Region1Results* are exported to the specified Excel file.
- If the Excel file already exists, a new worksheet named *Region 1* is added to the file. If a worksheet with that name already exists in the file, the contents of that worksheet are overwritten.
- If the Excel file does not exist, a file with the specified name is created, and a worksheet named *Region 1* is added to the new file.

## **NAME Keyword**

The optional `NAME` keyword specifies the Viewer document from which to export output. This is useful when more than one Viewer document is open and you want to ensure that the command operates on the correct document.

**NAME=\***. *Export the contents of the designated output document.* This is the default.

**NAME=name**. *Export the contents of the named output document.* This is the name assigned by a previous `OUTPUT NAME`, `OUTPUT OPEN`, `OUTPUT NEW` command, or automatically assigned when the document is created or opened in the current session; it is not the file name associated with a saved output document. In the Viewer window title bar, the name is enclosed in square brackets.

### **Example**

```
OUTPUT EXPORT NAME=Region1Results  
/XLS DOCUMENTFILE=' /mydocs/ResultsByRegion.xls'.
```

## **CONTENTS Subcommand**

The optional `CONTENTS` subcommand Viewer controls which contents are exported.

### **EXPORT Keyword**

The `EXPORT` keyword specifies whether to export all Viewer items or only those that are visible.

**ALL**. *Export all Viewer items.* All Viewer items are exported regardless of whether they are visible or hidden. This is the default.

**VISIBLE**. *Export visible items only.* Hidden items are not included.

### **LAYERS Keyword**

By default, inclusion or exclusion of pivot table layers is controlled by the table properties for each pivot table. You can override this setting and include all layers or exclude all but the currently visible layer.

**PRINTSETTING**. *Export layers as specified for printing in Table Properties.* This is the default.

**ALL**. *Export all layers.* Each layer is included as a separate table in the exported file.

**VISIBLE**. *Export visible layer only.* Only the layer currently displayed in the Viewer window is included in the exported file.

### **MODELVIEWS Keyword**

By default, inclusion or exclusion of model views is controlled by the model properties for each model. You can override this setting and include all views or exclude all but the currently visible view in the Model Viewer.

**PRINTSETTING.** *Export model views as specified for printing in Model Properties.* This is the default.

**ALL.** *Export all model views.* Each item is included as a separate image.

**VISIBLE.** *Export visible view only.* Only the model view currently displayed in the Viewer window is included.

Note: All model views are exported as images. This includes any tables.

### **Example**

```
OUTPUT EXPORT
  /CONTENTS EXPORT=VISIBLE LAYERS=ALL MODELVIEWS=VISIBLE
  /DOC DOCUMENTFILE=' /mydocs/myresults.doc' .
```

## **DOC Subcommand**

The DOC subcommand exports Viewer documents in Word (RTF) format. Pivot tables are exported as Word tables with all formatting attributes intact—for example, cell borders, font styles, and background colors. Text output is exported as formatted RTF. Charts, tree diagrams, and model views are included in PNG format.

### **DOCUMENTFILE Keyword**

The required DOCUMENTFILE keyword specifies the name of the Word document. The keyword is followed by an equals sign and a quoted file specification, as in:

```
OUTPUT EXPORT
  /DOC DOCUMENTFILE=' /mydocs/myresults.doc' .
```

### **PAGESIZE Keyword**

The PAGESIZE keyword controls the page dimensions of the exported document. The keyword is followed by an equals sign and one of the following alternatives:

**INCHES(width, height).** *Height and width specified in inches.* Both values are required and both must be positive numbers.

**MM(width, height).** *Height and width specified in millimeters.* Both values are required and must be positive numbers.

- The default page dimensions are the default settings for the default printer.
- The orientation of the page is implied by the dimensions. Specifying a width that is less than the height is equivalent to choosing portrait orientation for the page. Specifying a width that is greater than the height is equivalent to choosing landscape orientation.

Note: the page size width setting is ignored if WIDETABLES=ASIS and one or more pivot tables is wider than the specified width. The page width automatically expands to accommodate the widest table.

**TOPMARGIN, BOTTOMMARGIN, LEFTMARGIN, and RIGHTMARGIN Keywords**

These keywords control the margins for the exported document. Each keyword is followed by an equals sign and one of the following alternatives:

**INCHES(n).** *Margin specified in inches.* The value must be a positive number.

**MM(n).** *Margin specified in millimeters.* The value must be a positive number.

**NOTESCAPTIONS Keyword**

The NOTESCAPTIONS keyword controls the inclusion of pivot table footnotes and captions in the exported document. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include footnotes and captions.* This is the default.

**NO.** *Do not include footnotes and captions.*

**WIDETABLES Keyword**

The WIDETABLES keyword controls the treatment of pivot tables that are too wide for the document width (the specified page width minus the left and right margins). The keyword is followed by an equals sign and one of the following alternatives:

**WRAP.** *Wrap wide tables.* The table is divided into sections that will fit within the defined document width. Row labels are repeated for each section of the table. This is the default. If the row labels are too wide for the defined document width, the table is exported without wrapping and will appear truncated in the document.

**SHRINK.** *Shrink wide tables to fit within the document width.* Font size and column width are reduced so that the table fits within the document width.

**ASIS.** *Do not shrink or wrap wide tables.* The document width expands to accommodate the widest table.

**Example**

```
OUTPUT EXPORT
  /DOC DOCUMENTFILE=' /mydocs/myresults.doc '
  PAGESIZE=INCHES(11, 8.5)
  TOPMARGIN=INCHES(1)
  BOTTOMMARGIN=INCHES(1)
  LEFTMARGIN=INCHES(1)
  RIGHTMARGIN=INCHES(1)
  NOTESCAPTIONS=NO
  WIDETABLES=SHRINK.
```

**HTML Subcommand**

The HTML subcommand exports Viewer documents in HTML format. Output objects that would be pivot tables in the Viewer are converted to simple HTML tables. No TableLook attributes (font characteristics, border styles, colors, etc.) are supported. Text output objects are tagged <PRE> in

the HTML. Charts, tree diagrams, and model views are exported as separate files in the selected graphics format and are embedded by reference. Image file names use the HTML file name as the root name, followed by a sequential integer, starting with 0.

### **DOCUMENTFILE Keyword**

The required DOCUMENTFILE keyword specifies the name of the HTML document. The keyword is followed by an equals sign and a quoted file specification, as in:

```
OUTPUT EXPORT
  /HTML DOCUMENTFILE=' /mydocs/myresults.htm' .
```

### **NOTESCAPTIONS Keyword**

The NOTESCAPTIONS keyword controls the inclusion of pivot table footnotes and captions in the exported document. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include footnotes and captions.* This is the default.

**NO.** *Do not include footnotes and captions.*

### **IMAGEFORMAT Keyword**

The IMAGEFORMAT keyword controls the format of the exported graphics. The keyword is followed by an equals sign and one of the following alternatives:

**BMP.** *Bitmap format.*

**EMF.** *Enhanced metafile format.* A vector graphic format that is available only on Windows operating systems.

**EPS.** *Encapsulated PostScript format.* This format is often used when the quality of printed images is important.

**JPG.** *Joint Photographic Experts Group format.* This is the default.

**PNG.** *Portable Network Graphics format.*

**TIF.** *Tagged image file format.*

Note: Graphics includes charts, tree diagrams, and any Model Viewer items, including Model Viewer tables.

### **Example**

```
OUTPUT EXPORT
  /HTML DOCUMENTFILE=' /mydocs/myresults.htm'
  NOTESCAPTIONS=NO IMAGEFORMAT=PNG.
```

## **PDF Subcommand**

The PDF subcommand exports output in PDF format. Output is exported as it appears in Print Preview, with all formatting attributes intact.

### **DOCUMENTFILE Keyword**

The required DOCUMENTFILE keyword specifies the name of the PDF document. The keyword is followed by an equals sign and a quoted file specification, as in:

```
OUTPUT EXPORT  
  /PDF DOCUMENTFILE=' /mydocs/myresults.pdf '.
```

### **EMBEDBOOKMARKS Keyword**

The EMBEDBOOKMARKS keyword specifies whether to include bookmarks in the PDF document that correspond to the Viewer outline entries. Like the Viewer outline pane, bookmarks can make it much easier to navigate documents with a large number of output objects. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include bookmarks.* This is the default.

**NO.** *Do not include bookmarks.*

### **EMBEDFONTS Keyword**

The EMBEDFONTS keyword specifies whether to embed fonts in the exported document. Embedding fonts ensures that the PDF document will look the same on all computers. Otherwise, if some fonts used in the document are not available on the computer being used to view (or print) the PDF document, font substitution may yield suboptimal results. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Embed fonts.* This is the default.

**NO.** *Do not embed fonts.*

### **Other Settings That Affect PDF Output**

**Page Setup/Page Attributes.** Page size, orientation, margins, content and display of page headers and footers, and printed chart size in PDF documents are controlled by page setup and page attribute options.

**Table Properties/TableLooks.** Scaling of wide and/or long tables and printing of table layers are controlled by table properties for each table. These properties can also be saved in TableLooks.

**Default/Current Printer.** The resolution (DPI) of the PDF document is the current resolution setting for the default or currently selected printer (which can be changed using Page Setup). The maximum resolution is 1200 DPI. If the printer setting is higher, the PDF document resolution will be 1200 DPI.

*Note:* High-resolution documents may yield poor results when printed on lower-resolution printers.



**Example**

```
OUTPUT EXPORT
 /PDF DOCUMENTFILE=' /mydocs/myresults.pdf '
 EMBEDBOOKMARKS=NO EMBEDFONTS=NO.
```

**PPT Subcommand**

The PPT subcommand exports output in PowerPoint format. This is available only on Windows operating systems and requires an installed version of PowerPoint 97 or later.

- Pivot tables are exported as Word tables, and each table is a separate slide in the PowerPoint file. All formatting attributes of the pivot table are retained—for example, cell borders, font styles, and background colors.
- Graphics are exported in TIFF format, and each graphic is separate slide in the PowerPoint file.
- Text output is not included.

**DOCUMENTFILE Keyword**

The required DOCUMENTFILE keyword specifies the name of the PowerPoint document. The keyword is followed by an equals sign and a quoted file specification, as in:

```
OUTPUT EXPORT
 /PPT DOCUMENTFILE=' /mydocs/myresults.ppt ' .
```

**SLIDESIZE Keyword**

The SLIDESIZE keyword controls the slide dimensions of the exported file. The keyword is followed by an equals sign and one of the following alternatives:

**INCHES(width, height).** *Height and width specified in inches.* Both values are required and both must be positive numbers.

**MM(width, height).** *Height and width specified in millimeters.* Both values are required and must be positive numbers.

**TOPMARGIN, BOTTOMMARGIN, LEFTMARGIN, and RIGHTMARGIN Keywords**

These keywords control the margins for the exported document. Each keyword is followed by an equals sign and one of the following alternatives:

**INCHES(n).** *Margin specified in inches.* The value must be a positive number.

**MM(n).** *Margin specified in millimeters.* The value must be a positive number.

**NOTESCAPTIONS Keyword**

The NOTESCAPTIONS keyword controls the inclusion of pivot table footnotes and captions in the exported document. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include footnotes and captions.* This is the default.

**NO.** *Do not include footnotes and captions.*

**SLIDETITLES Keyword**

The SLIDETITLES keyword controls the inclusion slide titles of the exported file. The title is formed from the outline entry for the item in the outline pane of the Viewer. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include slide titles.* This is the default.

**NO.** *Do not include slide titles.*

**WIDETABLES Keyword**

The WIDETABLES keyword controls the treatment of pivot tables that are too wide for the document width (the specified page width minus the left and right margins). The keyword is followed by an equals sign and one of the following alternatives:

**WRAP.** *Wrap wide tables.* The table is divided into sections that will fit within the defined document width. Row labels are repeated for each section of the table. This is the default. If the row labels are too wide for the defined document width, the table is exported without wrapping and will appear truncated in the document.

**SHRINK.** *Shrink wide tables to fit within the document width.* Font size and column width are reduced so that the table fits within the document width.

**ASIS.** *Do not shrink or wrap wide tables.* The document width expands to accommodate the widest table.

**Example**

```
OUTPUT EXPORT
  /PPT DOCUMENTFILE='/mydocs/myresults.ppt'
  SLIDESIZE=INCHES(10, 7.5)
  SLIDETITLES=NO
  WIDETABLES=ASIS.
```

**TEXT Subcommand**

The TEXT subcommand exports output in text format.

- Text output formats include local encoding, UTF-8, and UTF-16.
- Pivot tables can be exported in tab-separated or space-separated format.
- All text output is exported in space-separated format.

- Graphics are exported as separate files in the selected graphics format. A line is inserted in the text file indicating the file name of each exported image.
- Graphics are exported to the same folder as the text document. Image file names use the text file name as the root name, followed by a sequential integer, starting with 0.

### **DOCUMENTFILE Keyword**

The required DOCUMENTFILE keyword specifies the name of the text document. The keyword is followed by an equals sign and a quoted file specification, as in:

```
OUTPUT EXPORT  
  /TEXT DOCUMENTFILE=' /mydocs/myresults.txt '.
```

### **ENCODING Keyword**

The ENCODING keyword specifies the type of text file to create. The keyword is followed by an equals sign and one of the following alternatives:

**LOCALE.** *Export in the character set (code page) of the current locale.* This is the default. For more information, see [SET Command](#), [LOCALE subcommand](#).

**UTF8.** *Export as 8-bit Unicode text.*

**UTF16.** *Export as 16-bit Unicode text.*

### **NOTESCAPTIONS Keyword**

The NOTESCAPTIONS keyword controls the inclusion of pivot table footnotes and captions in the exported document. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include footnotes and captions.* This is the default.

**NO.** *Do not include footnotes and captions.*

### **SEPARATOR Keyword**

Pivot table output can be exported as space-separated or tab-separated text. The SEPARATOR keyword is followed by an equals sign and one of the following alternatives:

**SPACE.** *Use spaces to separate columns.* This is the default.

**TAB.** *Use tabs to separate columns.*

### **COLUMNWIDTH Keyword**

For space-separated pivot table output, the COLUMNWIDTH keyword controls column width. The available alternatives are:

**COLUMNWIDTH=AUTOFIT.** *Make each column as wide as the widest label or value in that column.* Column contents will not be wrapped. This is the default.

**COLUMNWIDTH=n.** *Wrap labels or values longer than n characters wide.* The value of *n* must be a positive integer.

Note: This setting affects only pivot tables exported in space-separated format (SEPARATOR=SPACE).

### **ROWBORDER Keyword**

For space-separated pivot table output, the ROWBORDER keyword specifies the character used to create row borders. The keyword is followed by an equals sign and a single quoted character. The default is a hyphen (-).

Note: This setting affects only pivot tables exported in space-separated format (SEPARATOR=SPACE).

### **COLUMNBORDER Keyword**

For space-separated pivot table output, the COLUMNBORDER keyword specifies the character used to create column borders. The keyword is followed by an equals sign and a single quoted character. The default is the vertical bar/pipe symbol (|).

Note: This setting affects only pivot tables exported in space-separated format (SEPARATOR=SPACE).

### **IMAGEFORMAT Keyword**

The IMAGEFORMAT keyword controls the format of the exported graphics. The keyword is followed by an equals sign and one of the following alternatives:

**BMP.** *Bitmap format.*

**EMF.** *Enhanced metafile format.* A vector graphic format that is available only on Windows operating systems.

**EPS.** *Encapsulated PostScript format.* This format is often used when the quality of printed images is important.

**JPG.** *Joint Photographic Experts Group format.* This is the default.

**PNG.** *Portable Network Graphics format.*

**TIF.** *Tagged image file format.*

Note: Graphics includes charts, tree diagrams, and any Model Viewer items, including Model Viewer tables.

### **Example**

```
OUTPUT EXPORT
  /TEXT DOCUMENTFILE= '/mydocs/myresults.txt'
  ENCODING=UTF8
  SEPARATOR=TAB
  IMAGEFORMAT=PNG.
```

## ***XLS Subcommand***

The XLS subcommand exports output in Excel 97-2003 format. Pivot table rows, columns, and cells are exported as Excel rows, columns, and cells, with all formatting attributes intact — for example, cell borders, font styles, and background colors. Text output is exported with all font attributes intact. Each line in the text output is a row in the Excel file, with the entire contents of the line contained in a single cell. Charts, tree diagrams, and model views are included in PNG format.

### ***DOCUMENTFILE Keyword***

The required DOCUMENTFILE keyword specifies the name of the Excel document. The keyword is followed by an equals sign and a quoted file specification, as in:

```
OUTPUT EXPORT
  /XLS DOCUMENTFILE=' /mydocs/myresults.xls ' .
```

### ***OPERATION Keyword***

The OPERATION keyword specifies the type of operation to be performed by the export. The keyword is followed by an equals sign and one of the following alternatives:

**CREATEFILE.** *Create an Excel file.* If a file with the specified name already exists, it is overwritten. This is the default.

**CREATESHEET.** *Create a new worksheet within an Excel file.* If a worksheet with the specified name (specified on the SHEET keyword) already exists in the file, that worksheet is overwritten. If the file doesn't already exist, a new file is created with a worksheet with the specified name.

**MODIFYSHEET.** *Modify the contents of the specified worksheet.* Content is added to the worksheet based on the specification on the LOCATION keyword. You must specify a sheet name with the SHEET keyword. If you modify an existing worksheet, charts, model views, and tree diagrams are not included in the exported output.

### ***SHEET Keyword***

The SHEET keyword specifies the name of the worksheet to which items will be exported. The keyword is followed by an equals sign and a sheet name enclosed in quotes.

- Sheet names cannot exceed 31 characters and cannot contain forward or back slashes, square brackets, question marks, or asterisks.
- If the specified worksheet doesn't exist in the Excel file, a new worksheet with that name will be created, regardless of operation method.
- The SHEET keyword is required with OPERATION=MODIFYSHEET. It is optional with the other operation methods.

**LOCATION Keyword**

The `LOCATION` keyword specifies the location within the worksheet for the exported items. The keyword is followed by an equals sign and one of the following alternatives:

**LASTCOLUMN.** *Add exported items after the last column.* Exported items will be added after the last column, starting in the first row, without modifying any existing contents. This is a good choice for adding new columns to an existing worksheet. This is the default.

**LASTROW.** *Add exported items after the last row.* Exported items will be added after the last row, starting in the first column, without modifying any existing contents. This is a good choice for adding new rows to an existing worksheet.

**STARTCELL('cell').** *Add exported items beginning at a specific cell location.* Exported items will be added to the worksheet starting at the specified cell location. Any existing content in the area where the exported items are added will be overwritten. The cell reference is required and must be enclosed in quotes.

**NOTESCAPTIONS Keyword**

The `NOTESCAPTIONS` keyword controls the inclusion of pivot table footnotes and captions in the exported document. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include footnotes and captions.* This is the default.

**NO.** *Do not include footnotes and captions.*

**Example**

```
OUTPUT EXPORT
  /XLS DOCUMENTFILE=' /mydocs/myresults.xls '
  OPERATION=MODIFYSHEET
  SHEET='Region 1'
  LOCATION=STARTCELL('B2').
```

**BMP Subcommand**

For graphics exported as separate files in BMP format, the `BMP` subcommand controls settings for those graphic files.

**IMAGEROOT Keyword**

- The `IMAGEROOT` keyword is required if you are exporting graphics only, without any document type.
- The keyword is followed by an equals sign and a root name enclosed in quotes. The quoted root name can also include a directory path.
- File names are generated by adding a sequential integer to the root name, starting with 0.
- If any subcommand with a `DOCUMENTFILE` keyword is specified, `IMAGEROOT` is ignored.

**PERCENTSIZE Keyword**

The PERCENTSIZE keyword specifies the size of graphics as a percentage of the original size. The keyword is followed by an equals sign and a positive number that does not exceed 200. The default value is 100.

**COMPRESS Keyword**

The COMPRESSION keyword controls image compression to reduce file size. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Compress images.* A lossless technique is used to reduce file size without affecting image quality.

**NO.** *Do not compress images.* This is the default.

**Example**

```
OUTPUT EXPORT
  /BMP IMAGEROOT=' /mydocs/mygraphic '
  PERCENTSIZE=50 COMPRESS=YES
```

**EMF Subcommand**

For graphics exported as separate files in enhanced metafile format, the EMF subcommand controls settings for those graphic files. EMF format is available only on Windows operating systems.

**IMAGEROOT Keyword**

- The IMAGEROOT keyword is required if you are exporting graphics only, without any document type.
- The keyword is followed by an equals sign and a root name enclosed in quotes. The quoted root name can also include a directory path.
- File names are generated by adding a sequential integer to the root name, starting with 0.
- If any subcommand with a DOCUMENTFILE keyword is specified, IMAGEROOT is ignored.

**PERCENTSIZE Keyword**

The PERCENTSIZE keyword specifies the size of graphics as a percentage of the original size. The keyword is followed by an equals sign and a positive number that does not exceed 200. The default value is 100.

**Example**

```
OUTPUT EXPORT
  /EMF IMAGEROOT=' /mydocs/mygraphic '
  PERCENT=50.
```

## ***EPS Subcommand***

For graphics exported as separate files in encapsulated PostScript format, the `EPS` subcommand controls settings for those graphic files.

### ***IMAGEROOT Keyword***

- The `IMAGEROOT` keyword is required if you are exporting graphics only, without any document type.
- The keyword is followed by an equals sign and a root name enclosed in quotes. The quoted root name can also include a directory path.
- File names are generated by adding a sequential integer to the root name, starting with 0.
- If any subcommand with a `DOCUMENTFILE` keyword is specified, `IMAGEROOT` is ignored.

### ***WIDTH Keyword***

The `WIDTH` keyword controls the size of the exported image by specifying a value for the width of the image. The original aspect ratio is always maintained. The keyword is followed by an equals sign and one of the following alternatives:

**PERCENT(n).** *Width specified as percentage of the original width.* The value enclosed in parentheses must be a positive number that does not exceed 200. The default value is 100.

**INCHES(n).** *Width specified in inches.* The value enclosed in parentheses must be a positive number.

**CM(n).** *Width specified in centimeters.* The value enclosed in parentheses must be a positive number.

**POINTS(n).** *Width specified in points.* The value enclosed in parentheses must be a positive number.

### ***PREVIEW Keyword***

The `PREVIEW` keyword controls the inclusion of a TIFF preview image along with the EPS image. The preview image will be displayed if the application you are using is unable to display EPS images on the screen. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Include a preview imaged.* This is the default.

**NO.** *Do not include a preview image.*

### ***FONTHANDLING Keyword***

The `FONTHANDLING` keyword controls the treatment of fonts in EPS images. The keyword is followed by an equals sign and one of the following alternatives:

**REFERENCE.** *Use font references.* If the fonts that are used in the graphic are available on the output device, those fonts are used. Otherwise, the output device uses alternate fonts. This is the default.



**CURVES.** *Convert fonts to curves.* This option turns fonts into PostScript curve data. The text itself is no longer editable as text in applications that can edit EPS graphics. This option is useful if the fonts that are used in the chart are not available on the output device.

### **Example**

```
OUTPUT EXPORT
/EPS IMAGEROOT='/mydocs/mygraphic'
WIDTH=PERCENT(50) PREVIEW=NO FONTHANDLING=CURVES.
```

## **JPG Subcommand**

For graphics exported as separate files in JPEG format, the `JPG` subcommand controls settings for those graphic files.

### **IMAGEROOT Keyword**

- The `IMAGEROOT` keyword is required if you are exporting graphics only, without any document type.
- The keyword is followed by an equals sign and a root name enclosed in quotes. The quoted root name can also include a directory path.
- File names are generated by adding a sequential integer to the root name, starting with 0.
- If any subcommand with a `DOCUMENTFILE` keyword is specified, `IMAGEROOT` is ignored.

### **PERCENTSIZE Keyword**

The `PERCENTSIZE` keyword specifies the size of graphics as a percentage of the original size. The keyword is followed by an equals sign and a positive number that does not exceed 200. The default value is 100.

### **GRAYSCALE Keyword**

The `GRAYSCALE` keyword controls conversion of color images to grayscale images. The keyword is followed by an equals sign and one of the following alternatives:

**YES.** *Convert images to grayscale.* Color are converted to shades of gray.

**NO.** *Do not convert images to grayscale.* This is the default.

### **Example**

```
OUTPUT EXPORT
/JPG IMAGEROOT='/mydocs/mygraphic'
PERCENT=50 GRAYSCALE=YES.
```

## **PNG Subcommand**

For graphics exported as separate files in PNG format, the `PNG` subcommand controls settings for those graphic files.

**IMAGEROOT Keyword**

- The `IMAGEROOT` keyword is required if you are exporting graphics only, without any document type.
- The keyword is followed by an equals sign and a root name enclosed in quotes. The quoted root name can also include a directory path.
- File names are generated by adding a sequential integer to the root name, starting with 0.
- If any subcommand with a `DOCUMENTFILE` keyword is specified, `IMAGEROOT` is ignored.

**PERCENTSIZE Keyword**

The `PERCENTSIZE` keyword specifies the size of graphics as a percentage of the original size. The keyword is followed by an equals sign and a positive number that does not exceed 200. The default value is 100.

**DEPTH Keyword**

The `DEPTH` keyword controls the number of colors in the exported graphics. If the number of colors in the graphic exceeds the number of colors for that depth, the colors will be dithered to replicate the colors in the graphic. The keyword is followed by an equals sign and one of the following alternatives:

**SCREEN.** *Current screen depth.* The number of colors currently displayed on your computer monitor. This is the default.

**BLACKWHITE.** *Black and white.* Each color is converted to either black or white.

**GRAYS256.** *Grayscale.* Colors are converted to shades of gray.

**COLORS16.** *16 colors.*

**COLORS256.** *256 colors.*

**TRUE24BIT.** *24-bit true color.*

**TRUE32BIT.** *32-bit true color.*

**Example**

```
OUTPUT EXPORT
/PNG IMAGEROOT=' /mydocs/mygraphic '
PERCENT=50
DEPTH=COLORS256.
```

## TIF Subcommand

For graphics exported as separate files in TIFF, the `TIF` subcommand controls settings for those graphic files.

**IMAGEROOT Keyword**

- The IMAGEROOT keyword is required if you are exporting graphics only, without any document type.
- The keyword is followed by an equals sign and a root name enclosed in quotes. The quoted root name can also include a directory path.
- File names are generated by adding a sequential integer to the root name, starting with 0.
- If any subcommand with a DOCUMENTFILE keyword is specified, IMAGEROOT is ignored.

**PERCENTSIZE Keyword**

The PERCENTSIZE keyword specifies the size of graphics as a percentage of the original size. The keyword is followed by an equals sign and a positive number that does not exceed 200. The default value is 100.

**Example**

```
OUTPUT EXPORT
/TIF IMAGEROOT='/mydocs/mygraphic'
PERCENT=50.
```

# OUTPUT NAME

OUTPUT NAME [NAME]=name

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 15.0

- Command introduced.

## Example

```
GET FILE='/examples/data/SalaryData2005.sav'.
DESCRIPTIVES salbegin salary.
OUTPUT NAME alleduclevels.
```

```
TEMPORARY.
SELECT IF (educ>12).
OUTPUT NEW NAME=over12.
DESCRIPTIVES salbegin salary.
```

```
GET FILE='/examples/data/SalaryData2000.sav'.
TEMPORARY.
SELECT IF (educ>12).
DESCRIPTIVES salbegin salary.
```

```
OUTPUT ACTIVATE alleduclevels.
DESCRIPTIVES salbegin salary.
```

## Overview

The OUTPUT commands (OUTPUT NEW, OUTPUT NAME, OUTPUT ACTIVATE, OUTPUT OPEN, OUTPUT SAVE, OUTPUT CLOSE) provide the ability to programmatically manage one or many output documents. These functions allow you to:

- Save an output document through syntax.
- Programmatically partition output into separate output documents (for example, results for males in one output document and results for females in a separate one).
- Work with multiple open output documents in a given session, selectively appending new results to the appropriate document.

The OUTPUT NAME command assigns a name to the designated output document. The designated output document is the one most recently opened (by OUTPUT NEW or OUTPUT OPEN) or activated (by OUTPUT ACTIVATE). The document name is used to reference the document in any subsequent OUTPUT ACTIVATE, OUTPUT SAVE, and OUTPUT CLOSE commands.

### Basic Specification

The basic specification for OUTPUT NAME is the command name followed by a name that conforms to variable naming rules. For more information, see the topic [Variable Names](#) on p. 49. The NAME keyword is optional, but if it is used it must be followed by an equals sign.

### Operations

- The association with the existing name is broken, and the new name is assigned to the document.
- If the specified name is associated with another document, that association is broken and the name is associated with the designated output document. The document previously associated with the specified name is assigned a new unique name.

### Example

```
GET FILE='/examples/data/SurveyData.sav'.

TEMPORARY.
SELECT IF (Sex='Male').
FREQUENCIES VARIABLES=ALL.
OUTPUT NAME males.

TEMPORARY.
SELECT IF (Sex='Female').
OUTPUT NEW NAME=females.
FREQUENCIES VARIABLES=ALL.

GET FILE='/examples/data/Preference.sav'.

TEMPORARY.
SELECT IF (Sex='Female').
DESCRIPTIVES VARIABLES=product1 product2 product3.

TEMPORARY.
SELECT IF (Sex='Male').
OUTPUT ACTIVATE males.
DESCRIPTIVES VARIABLES=product1 product2 product3.

OUTPUT SAVE NAME=males OUTFILE='/examples/output/Males.spv'.
OUTPUT SAVE NAME=females OUTFILE='/examples/output/Females.spv'.
```

- The first GET command loads survey data for males and females.
- FREQUENCIES output for male respondents is written to the designated output document. The OUTPUT NAME command is used to assign the name *males* to the designated output document.
- FREQUENCIES output for female respondents is written to a new output document named *females*.
- The second GET command loads preferences data for males and females.
- Descriptive statistics for females are appended to the output document named *females* and those for males are appended to the output document named *males*. Each output document now contains both survey and preferences results.
- The two open output documents are saved to separate files. Because the operation of saving an output document does not close it, both documents remain open. The output document named *males* remains the designated output document.

# OUTPUT NEW

```
OUTPUT NEW [NAME=name]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 15.0

- Command introduced.

Release 16.0

- TYPE keyword is obsolete and is ignored.

## **Example**

```
GET FILE='/examples/data/Males.sav'.  
FREQUENCIES VARIABLES=ALL.  
OUTPUT SAVE OUTFILE='/examples/output/Males.spv'.  
OUTPUT NEW.
```

```
GET FILE='/examples/data/Females.sav'.  
FREQUENCIES VARIABLES=ALL.  
OUTPUT SAVE OUTFILE='/examples/output/Females.spv'.
```

## **Overview**

The OUTPUT commands (OUTPUT NEW, OUTPUT NAME, OUTPUT ACTIVATE, OUTPUT OPEN, OUTPUT SAVE, OUTPUT CLOSE) provide the ability to programmatically manage one or many output documents. These functions allow you to:

- Save an output document through syntax.
- Programmatically partition output into separate output documents (for example, results for males in one output document and results for females in a separate one).
- Work with multiple open output documents in a given session, selectively appending new results to the appropriate document.

The OUTPUT NEW command creates a new output document, which becomes the designated output document. Subsequent procedure output is directed to the new output document until the document is closed or another output document is created, opened, or activated.

## **Basic Specification**

The basic specification for OUTPUT NEW is simply the command name.

### **TYPE Keyword**

This keyword is obsolete and is ignored. The only valid output type is Viewer. Draft Viewer format is no longer supported. To produce text output equivalent to Draft Viewer output use OMS. For more information, see the topic OMS on p. 1403.

### **NAME Keyword**

By default, the newly created output document is provided with a unique name. You can optionally specify a custom name for the output document, overriding the default name. The document name is used to reference the document in any subsequent OUTPUT ACTIVATE, OUTPUT SAVE, and OUTPUT CLOSE commands.

- The specified name must conform to variable naming rules. For more information, see the topic Variable Names on p. 49.
- If the specified name is associated with another document, that association is broken and the name is associated with the new document. The document previously associated with the specified name is assigned a new unique name.

### **Syntax Rules**

- An error occurs if a keyword is specified more than once.
- Keywords must be spelled in full.
- Equals signs (=) used in the syntax chart are required elements.

### **Operations**

The new output document is opened in a window in the user interface and becomes the designated output window.

### **Limitations**

Because each window requires a minimum amount of memory, there is a limit to the number of windows, IBM® SPSS® Statistics or otherwise, that can be concurrently open on a given system. The particular number depends on the specifications of your system and may be independent of total memory due to OS constraints.

### **Example**

```
GET FILE='/examples/data/Males.sav'.  
FREQUENCIES VARIABLES=ALL.  
OUTPUT SAVE OUTFILE='/examples/output/Males.spv'.  
OUTPUT NEW.
```

```
GET FILE='/examples/data/Females.sav'.  
FREQUENCIES VARIABLES=ALL.  
OUTPUT SAVE OUTFILE='/examples/output/Females.spv'.
```

- FREQUENCIES produces summary statistics for each variable in */examples/data/Males.sav*. The output from FREQUENCIES is added to the designated output document (one is created automatically if no output document is currently open).

---

*OUTPUT NEW*

- `OUTPUT SAVE` writes the contents of the designated output document to */examples/output/Males.spv*.
- `OUTPUT NEW` creates a new Viewer document, which becomes the designated output document.
- The subsequent `FREQUENCIES` command produces output for females using the data in */examples/data/Females.sav*. `OUTPUT SAVE` writes this output to */examples/output/Females.spv*.

As shown in this example, `OUTPUT NEW` allows you to direct results to a new output document. If `OUTPUT NEW` were not specified, */examples/output/Females.spv* would contain frequencies for both males and females.



# OUTPUT OPEN

```
OUTPUT OPEN FILE='file specification' [NAME=name]
      [LOCK={YES**}]
      {NO }
```

\*\* Default if the subcommand or keyword is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 15.0

- Command introduced.

Release 17.0

- LOCK keyword introduced.

## **Example**

```
OUTPUT OPEN FILE='/examples/output/Q1Output.spv'.
GET FILE='/examples/data/March.sav'.
FREQUENCIES VARIABLES=ALL.
OUTPUT SAVE OUTFILE='/examples/output/Q1Output.spv'.
```

## **Overview**

The OUTPUT commands (OUTPUT NEW, OUTPUT NAME, OUTPUT ACTIVATE, OUTPUT OPEN, OUTPUT SAVE, OUTPUT CLOSE) provide the ability to programmatically manage one or many output documents. These functions allow you to:

- Save an output document through syntax.
- Programmatically partition output into separate output documents (for example, results for males in one output document and results for females in a separate one).
- Work with multiple open output documents in a given session, selectively appending new results to the appropriate document.

The OUTPUT OPEN command opens a Viewer document, which becomes the designated output document. You can use OUTPUT OPEN to append output to an existing output document. Once opened, subsequent procedure output is directed to the document until it is closed or until another output document is created, opened, or activated.

## **Basic Specification**

The basic specification for OUTPUT OPEN is the command name followed by a file specification for the file to open.

**NAME Keyword**

By default, the newly opened output document is provided with a unique name. You can optionally specify a custom name for the output document, overriding the default name. The document name is used to reference the document in any subsequent `OUTPUT ACTIVATE`, `OUTPUT SAVE`, and `OUTPUT CLOSE` commands.

- The specified name must conform to variable naming rules. [For more information, see the topic Variable Names on p. 49.](#)
- If the specified name is associated with another document, that association is broken and the name is associated with the newly opened document. The document previously associated with the specified name is assigned a new unique name.

**LOCK Keyword**

The `LOCK` keyword only applies when opening a file from a IBM® SPSS® Collaboration and Deployment Services Repository. It specifies whether the file is locked (in the repository) when opened. Locking a file prevents other users from modifying it while you are working with it. Other users can open locked files but they cannot save them back to the repository. The lock is released when the file is closed or when the current session ends.

- By default, files are opened as locked.
- The `LOCK` keyword is ignored if the specified file does not reside in a IBM SPSS Collaboration and Deployment Services Repository.

[For more information, see the topic File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects on p. 2230.](#)

**Syntax Rules**

- An error occurs if a keyword is specified more than once.
- Keywords must be spelled in full.
- Equals signs (=) used in the syntax chart are required elements.

**Operations**

- The output document is opened in a window in the user interface and becomes the designated output window.
- An error occurs, but processing continues, if the specified file is not found. Output continues to be directed to the last designated output document.
- An error occurs, but processing continues, if the specified file is not a Viewer document. Output continues to be directed to the last designated output document.
- `OUTPUT OPEN` honors file handles and changes to the working directory made with the `CD` command.

**Limitations**

- Because each window requires a minimum amount of memory, there is a limit to the number of windows, IBM® SPSS® Statistics or otherwise, that can be concurrently open on a given system. The particular number depends on the specifications of your system and may be independent of total memory due to OS constraints.
- OUTPUT OPEN is not supported in the SPSS Statistics Batch Facility (available with SPSS Statistics Server).

**Example**

```
OUTPUT OPEN FILE='/examples/output/Q1Output.spv'.  
GET FILE='/examples/data/March.sav'.  
FREQUENCIES VARIABLES=ALL.  
OUTPUT SAVE OUTFILE='/examples/output/Q1Output.spv'.
```

- OUTPUT OPEN opens the Viewer document */examples/output/Q1Output.spv*. The document contains summaries for the months of January and February.
- The GET command opens a file containing data for the month of March.
- The FREQUENCIES command produces summaries for March data, which are appended to the designated output document.
- OUTPUT SAVE saves the designated output document to */examples/output/Q1Output.spv*. The saved document contains results for each of the three months in the first quarter.

# OUTPUT SAVE

```
OUTPUT SAVE [NAME={*   }]
                {name}
      OUTFILE='file specification'
      [TYPE={SPV**}]
            (SPW  )
      [LOCK={NO**}]
            {YES  }
```

\*\* Default if the keyword is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 15.0

- Command introduced.

Release 16.0

- TYPE keyword introduced.

Release 17.0

- LOCK keyword introduced.

## **Example**

```
OUTPUT OPEN FILE='/examples/output/Q1Output.spv'.
GET FILE='/examples/data/March.sav'.
FREQUENCIES VARIABLES=ALL.
OUTPUT SAVE OUTFILE='/examples/output/Q1Output.spv'.
```

## **Overview**

The OUTPUT commands (OUTPUT NEW, OUTPUT NAME, OUTPUT ACTIVATE, OUTPUT OPEN, OUTPUT SAVE, OUTPUT CLOSE) provide the ability to programmatically manage one or many output documents. These functions allow you to:

- Save an output document through syntax.
- Programmatically partition output into separate output documents (for example, results for males in one output document and results for females in a separate one).
- Work with multiple open output documents in a given session, selectively appending new results to the appropriate document.

The OUTPUT SAVE command saves the contents of an open output document to a file.

**Basic Specification**

The basic specification for `OUTPUT SAVE` is the command name followed by a file specification for the destination file.

**NAME Keyword**

Use the `NAME` keyword to save an output document other than the designated one. Provide the name associated with the document.

**TYPE Keyword**

Use the `TYPE` keyword to specify the format of the output file—`SPV` for standard output files and `SPW` for the IBM® SPSS® Statistics Web Reports format. Files in the `SPW` format that are stored in a IBM® SPSS® Collaboration and Deployment Services Repository can be viewed and manipulated over the Web, in real time, using a standard browser with release 3.5 or higher of IBM® SPSS® Collaboration and Deployment Services.

- `spw` files created from `OUTPUT SAVE` contain all visible objects from the associated Viewer window, and pivot tables are saved as interactive, meaning they can be manipulated when viewed over the Web. If you need greater control over items saved to an `spw` file, use the `OMS` command.

**LOCK Keyword**

Use `LOCK=YES` to prevent editing of Viewer documents in IBM® SPSS® Smartreader (a separate product for working with Viewer documents). If a Viewer document is locked, you can manipulate pivot tables (swap rows and columns, change the displayed layer, etc.) but you cannot edit any output or save any changes to the Viewer document in SPSS Smartreader. This setting has no effect on Viewer documents opened in SPSS Statistics. The default is `LOCK=NO`.

**Syntax Rules**

- An error occurs if a keyword is specified more than once.
- Keywords must be spelled in full.
- Equals signs (=) used in the syntax chart are required elements.

**Operations**

- By default, the designated output document is saved. The designated output document is the one most recently opened (by `OUTPUT NEW` or `OUTPUT OPEN`) or activated (by `OUTPUT ACTIVATE`).
- If the specified file already exists, `OUTPUT SAVE` overwrites it without warning.
- An error occurs if you specify a nonexistent output document.
- An error occurs if the file specification is invalid.
- `OUTPUT SAVE` saves the document but does not close it. Use `OUTPUT CLOSE` to close the document.

---

*OUTPUT SAVE*

- `OUTPUT SAVE` honors file handles and changes to the working directory made with the `CD` command.

**Limitations**

- `OUTPUT SAVE` is not supported in the SPSS Statistics Batch Facility (available with SPSS Statistics Server).

**Example**

```
OUTPUT OPEN FILE='/examples/output/Q1Output.spv'.
GET FILE='/examples/data/March.sav'.
FREQUENCIES VARIABLES=ALL.
OUTPUT SAVE OUTFILE='/examples/output/Q1Output.spv'.
```

- `OUTPUT OPEN` opens the Viewer document */examples/output/Q1Output.spv*. The document contains summaries for the months of January and February.
- `GET` opens a file containing new data for March.
- `FREQUENCIES` produces frequencies for March data, which are appended to the designated output document.
- `OUTPUT SAVE` saves the contents of the designated output document to */examples/output/Q1Output.spv*, which now contains results for the entire first quarter.

# OVERALS

OVERALS is available in the Categories option.

```
OVERALS VARIABLES=varlist (max)

/ANALYSIS=varlist[({ORDI**})]
                {SNOM  }
                {MNOM  }
                {NUME  }

/SETS= n (# of vars in set 1, ..., # of vars in set n)

[/NOBSERVATIONS=value]

[/DIMENSION={2**  }]
                {value}

[/INITIAL={NUMERICAL**}]
                {RANDOM  }

[/MAXITER={100**}]
                {value}

[/CONVERGENCE={.00001**}]
                {value  }

[/PRINT=[DEFAULT] [FREQ**] [QUANT] [CENTROID**]
        [HISTORY] [WEIGHTS**]
        [OBJECT] [FIT] [NONE]]

[/PLOT=[NDIM=({1    ,2    }**)]
        {value,value}
        {ALL  ,MAX  }
        [DEFAULT[(n)]] [OBJECT**[(varlist)][(n)]]
        [QUANT[(varlist)][(n)]] [LOADINGS**[(n)]]
        [TRANS[(varlist)]] [CENTROID[(varlist)][(n)]]
        [NONE]]

[/SAVE=[rootname] [(value)]]

[/MATRIX=OUT({*          })]
                {'savfile'|'dataset'}
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
OVERALS VARIABLES=PRETEST1 PRETEST2 POSTEST1 POSTEST2(20)
SES(5) SCHOOL(3)
/ANALYSIS=PRETEST1 TO POSTEST2 (NUME) SES (ORDI) SCHOOL (SNOM)
/SETS=3(2,2,2).
```

## Overview

OVERALS performs nonlinear canonical correlation analysis on two or more sets of variables. Variables can have different optimal scaling levels, and no assumptions are made about the distribution of the variables or the linearity of the relationships.

### ***Options***

**Optimal Scaling Levels.** You can specify the level of optimal scaling at which you want to analyze each variable.

**Number of Dimensions.** You can specify how many dimensions OVERALS should compute.

**Iterations and Convergence.** You can specify the maximum number of iterations and the value of a convergence criterion.

**Display Output.** The output can include all available statistics, only the default statistics, or only the specific statistics that you request. You can also control whether some of these statistics are plotted.

**Saving Scores.** You can save object scores in the active dataset.

**Writing Matrices.** You can write a matrix data file containing quantification scores, centroids, weights, and loadings for use in further analyses.

### ***Basic Specification***

- The basic specification is command OVERALS, the VARIABLES subcommand, the ANALYSIS subcommand, and the SETS subcommand. By default, OVERALS estimates a two-dimensional solution and displays a table listing optimal scaling levels of each variable by set, eigenvalues and loss values by set, marginal frequencies, centroids and weights for all variables, and plots of the object scores and component loadings.

### ***Subcommand Order***

- The VARIABLES subcommand, ANALYSIS subcommand, and SETS subcommand must appear in that order before all other subcommands.
- Other subcommands can appear in any order.

### ***Operations***

- If the ANALYSIS subcommand is specified more than once, OVERALS is not executed. For all other subcommands, if a subcommand is specified more than once, only the last occurrence is executed.
- OVERALS treats every value in the range 1 to the maximum value that is specified on VARIABLES as a valid category. To avoid unnecessary output, use the AUTORECODE or RECODE command to recode a categorical variable that has nonsequential values or that has a large number of categories. For variables that are treated as numeric, recoding is not recommended because the characteristic of equal intervals in the data will not be maintained (see AUTORECODE and RECODE for more information).

### ***Limitations***

- String variables are not allowed; use AUTORECODE to recode nominal string variables.
- The data must be positive integers. Zeros and negative values are treated as system-missing, which means that they are excluded from the analysis. Fractional values are truncated after the decimal and are included in the analysis. If one of the levels of a categorical variable has been coded 0 or some negative value, and you want to treat it as a valid category, use the AUTORECODE or RECODE command to recode the values of that variable.



- OVERALS ignores user-missing value specifications. Positive user-missing values that are less than the maximum value that is specified on the VARIABLES subcommand are treated as valid category values and are included in the analysis. If you do not want the category to be included, use COMPUTE or RECODE to change the value to a value outside of the valid range. Values outside of the range (less than 1 or greater than the maximum value) are treated as system-missing and are excluded from the analysis.
- If one variable in a set has missing data, all variables in that set are missing for that object (case).
- Each set must have at least three valid (non-missing, non-empty) cases.

## Examples

```
OVERALS VARIABLES=PRETEST1 PRETEST2 POSTEST1 POSTEST2 (20)
SES(5) SCHOOL(3)
/ANALYSIS=PRETEST1 TO POSTEST2 (NUME) SES (ORDI) SCHOOL (SNOM)
/SETS=3 (2,2,2)
/PRINT=OBJECT FIT
/PLOT=QUANT(PRETEST1 TO SCHOOL) .
```

- VARIABLES defines the variables and their maximum values.
- ANALYSIS specifies that all variables from *PRETEST1* to *POSTEST2* are to be analyzed at the numeric level of optimal scaling, *SES* is to be analyzed at the ordinal level, and *SCHOOL* is to be analyzed as a single nominal. These variables are all of the variables that will be used in the analysis.
- SETS specifies that there are three sets of variables to be analyzed and two variables in each set.
- PRINT lists the object and fit scores.
- PLOT plots the single-category and multiple-category coordinates of all variables in the analysis.

## VARIABLES Subcommand

VARIABLES specifies all variables in the current OVERALS procedure.

- The VARIABLES subcommand is required and precedes all other subcommands. The actual word VARIABLES can be omitted.
- Each variable or variable list is followed by the maximum value in parentheses.

## ANALYSIS Subcommand

ANALYSIS specifies the variables to be used in the analysis and the optimal scaling level at which each variable is to be analyzed.

- The ANALYSIS subcommand is required and follows the VARIABLES subcommand.
- The specification on ANALYSIS is a variable list and an optional keyword in parentheses, indicating the level of optimal scaling.

- The variables on `ANALYSIS` must also be specified on the `VARIABLES` subcommand.
- Only active variables are listed on the `ANALYSIS` subcommand. **Active variables** are those variables that are used in the computation of the solution. **Passive variables**, those variables that are listed on the `VARIABLES` subcommand but not on the `ANALYSIS` subcommand, are ignored in the `OVERALS` solution. Object score plots can still be labeled by passive variables.

The following keywords can be specified to indicate the optimal scaling level:

<b>MNOM</b>	<i>Multiple nominal.</i> The quantifications can be different for each dimension. When all variables are multiple nominal, and there is only one variable in each set, <code>OVERALS</code> gives the same results as <code>HOMALS</code> .
<b>SNOM</b>	<i>Single nominal.</i> <code>OVERALS</code> gives only one quantification for each category. Objects in the same category (cases with the same value on a variable) obtain the same quantification. When all variables are <code>SNOM</code> , <code>ORDI</code> , or <code>NUME</code> , and there is only one variable per set, <code>OVERALS</code> gives the same results as <code>PRINCALS</code> .
<b>ORDI</b>	<i>Ordinal.</i> This setting is the default for variables that are listed without optimal scaling levels. The order of the categories of the observed variable is preserved in the quantified variable.
<b>NUME</b>	<i>Numerical.</i> Interval or ratio scaling level. <code>OVERALS</code> assumes that the observed variable already has numerical values for its categories. When all variables are quantified at the numerical level, and there is only one variable per set, the <code>OVERALS</code> analysis is analogous to classical principal components analysis.

These keywords can apply to a variable list as well as to a single variable. Thus, the default `ORDI` is not applied to a variable without a keyword if a subsequent variable on the list has a keyword.

## ***SETS Subcommand***

`SETS` specifies how many sets of variables exist and how many variables are in each set.

- `SETS` is required and must follow the `ANALYSIS` subcommand.
- `SETS` is followed by an integer to indicate the number of variable sets. Following this integer is a list of values in parentheses, indicating the number of variables in each set.
- There must be at least two sets.
- The sum of the values in parentheses must equal the number of variables specified on the `ANALYSIS` subcommand. The variables in each set are read consecutively from the `ANALYSIS` subcommand.

An example is as follows:

```
/SETS=2 (2, 3)
```

This specification indicates that there are two sets. The first two variables that are named on `ANALYSIS` are the first set, and the last three variables that are named on `ANALYSIS` are the second set.

## ***NOBSERVATIONS Subcommand***

`NOBSERVATIONS` specifies how many cases are used in the analysis.

- If `NOOBSERVATIONS` is not specified, all available observations in the active dataset are used.
- `NOOBSERVATIONS` is followed by an integer, indicating that the first  $n$  cases are to be used.

## ***DIMENSION Subcommand***

`DIMENSION` specifies the number of dimensions that you want `OVERALS` to compute.

- If you do not specify the `DIMENSION` subcommand, `OVERALS` computes two dimensions.
- `DIMENSION` is followed by an integer indicating the number of dimensions.
- If all variables are `SNOM` (single nominal), `ORDI` (ordinal), or `NUME` (numerical), the maximum number of dimensions that you can specify is the total number of variables on the `ANALYSIS` subcommand.
- If some or all variables are `MNOM` (multiple nominal), the maximum number of dimensions that you can specify is the number of `MNOM` variable levels (categories) plus the number of non-`MNOM` variables, minus the number of `MNOM` variables.
- The maximum number of dimensions must be less than the number of observations minus 1.
- If the number of sets is 2, and all variables are `SNOM`, `ORDI`, or `NUME`, the number of dimensions should not be more than the number of variables in the smaller set.
- If the specified value is too large, `OVERALS` tries to adjust the number of dimensions to the allowable maximum. `OVERALS` might not be able to adjust if there are `MNOM` variables with missing data.

## ***INITIAL Subcommand***

The `INITIAL` subcommand specifies the method that is used to compute the initial configuration.

- The specification on `INITIAL` is keyword `NUMERICAL` or `RANDOM`. If the `INITIAL` subcommand is not specified, `NUMERICAL` is the default.

**NUMERICAL**      *Treat all variables except multiple nominal as numerical.* This specification is best to use when there are no `SNOM` variables.

**RANDOM**        *Compute a random initial configuration.* This specification should be used only when some or all variables are `SNOM`.

## ***MAXITER Subcommand***

`MAXITER` specifies the maximum number of iterations that `OVERALS` can go through in its computations.

- If `MAXITER` is not specified, `OVERALS` will iterate up to 100 times.
- The specification on `MAXITER` is an integer indicating the maximum number of iterations.

## ***CONVERGENCE Subcommand***

`CONVERGENCE` specifies a convergence criterion value. `OVERALS` stops iterating if the difference in fit between the last two iterations is less than the `CONVERGENCE` value.

- The default CONVERGENCE value is 0.00001.
- The specification on CONVERGENCE is any value that is greater than 0.000001. (Values that are less than this value might seriously affect performance.)

## ***PRINT Subcommand***

PRINT controls which statistics are included in your display output. The default output includes a table that lists optimal scaling levels of each variable by set; eigenvalues and loss values by set by dimension; and the output that is produced by keywords `FREQ`, `CENTROID`, and `WEIGHTS`.

The following keywords are available:

<b>FREQ</b>	<i>Marginal frequencies for the variables in the analysis.</i>
<b>HISTORY</b>	<i>History of the iterations.</i>
<b>FIT</b>	<i>Multiple fit, single fit, and single loss per variable.</i>
<b>CENTROID</b>	<i>Category quantification scores, the projected centroids, and the centroids.</i>
<b>OBJECT</b>	<i>Object scores.</i>
<b>QUANT</b>	<i>Category quantifications and the single and multiple coordinates.</i>
<b>WEIGHTS</b>	<i>Weights and component loadings.</i>
<b>DEFAULT</b>	<i>FREQ, CENTROID, and WEIGHTS.</i>
<b>NONE</b>	<i>Summary loss statistics.</i>

## ***PLOT Subcommand***

PLOT can be used to produce plots of transformations, object scores, coordinates, centroids, and component loadings.

- If PLOT is not specified, plots of the object scores and component loadings are produced.

The following keywords can be specified on PLOT:

<b>LOADINGS</b>	<i>Plot of the component loadings.</i>
<b>OBJECT</b>	<i>Plot of the object scores.</i>
<b>TRANS</b>	<i>Plot of category quantifications.</i>
<b>QUANT</b>	<i>Plot of all category coordinates.</i>
<b>CENTROID</b>	<i>Plot of all category centroids.</i>
<b>DEFAULT</b>	<i>OBJECT and LOADINGS.</i>
<b>NONE</b>	<i>No plots.</i>

- Keywords `OBJECT`, `QUANT`, and `CENTROID` can each be followed by a variable list in parentheses to indicate that plots should be labeled with these variables. For `QUANT` and `CENTROID`, the variables must be specified on both the `VARIABLES` and `ANALYSIS` subcommands. For `OBJECT`, the variables must be specified on `VARIABLES` but need not appear on `ANALYSIS`, meaning that variables that are not used in the computations can still be used to label `OBJECT` plots. If the variable list is omitted, the default plots are produced.

- Object score plots use category labels corresponding to all categories within the defined range. Objects in a category that is outside the defined range are labeled with the label corresponding to the category immediately following the defined maximum category.
- If TRANS is followed by a variable list, only plots for those variables are produced. If a variable list is not specified, plots are produced for each variable.
- All keywords except NONE can be followed by an integer in parentheses to indicate how many characters of the variable or value label are to be used on the plot. (If you specified a variable list after OBJECT, CENTROID, TRANS, or QUANT, you can specify the value in parentheses after the list.) The value can range from 1 to 20. If the value is omitted, 12 characters are used. Spaces between words count as characters.
- If a variable label is missing, the variable name is used for that variable. If a value label is missing, the actual value is used.
- Make sure that your variable and value labels are unique by at least one letter in order to distinguish them on the plots.
- When points overlap, the points are described in a summary following the plot.

In addition to the plot keywords, the following keyword can be specified:

**NDIM**      *Dimension pairs to be plotted.* NDIM is followed by a pair of values in parentheses. If NDIM is not specified, plots are produced for dimension 1 versus dimension 2.

- The first value indicates the dimension that is plotted against all higher dimensions. This value can be any integer from 1 to the number of dimensions minus 1.
- The second value indicates the highest dimension to be used in plotting the dimension pairs. This value can be any integer from 2 to the number of dimensions.
- Keyword ALL can be used instead of the first value to indicate that all dimensions are paired with higher dimensions.
- Keyword MAX can be used instead of the second value to indicate that plots should be produced up to and including the highest dimension fit by the procedure.

### **Example**

```
OVERALS COLA1 COLA2 JUICE1 JUICE2 (4)
/ANALYSIS=COLA1 COLA2 JUICE1 JUICE2 (SNOM)
/SETS=2(2,2)
/PLOT NDIM(1,3) QUANT(5) .
```

- The NDIM(1,3) specification indicates that plots should be produced for two dimension pairs—dimension 1 versus dimension 2 and dimension 1 versus dimension 3.
- QUANT requests plots of the category quantifications. The (5) specification indicates that the first five characters of the value labels are to be used on the plots.

### **Example**

```
OVERALS COLA1 COLA2 JUICE1 JUICE2 (4)
/ANALYSIS=COLA1 COLA2 JUICE1 JUICE2 (SNOM)
/SETS=2(2,2)
```

```
/PLOT NDIM(ALL,3) QUANT(5) .
```

- This plot is the same as above except for the `ALL` specification following `NDIM`, which indicates that all possible pairs up to the second value should be plotted. `QUANT` plots will be produced for dimension 1 versus dimension 2, dimension 2 versus dimension 3, and dimension 1 versus dimension 3.

## **SAVE Subcommand**

`SAVE` lets you add variables containing the object scores that are computed by `OVERALS` to the active dataset.

- If `SAVE` is not specified, object scores are not added to the active dataset.
- A variable rootname can be specified on the `SAVE` subcommand, to which `OVERALS` adds the number of the dimension. Only one rootname can be specified, and it can contain up to six characters.
- If a rootname is not specified, unique variable names are automatically generated. The variable names are `OVEn_m`, where  $n$  is a dimension number and  $m$  is a set number. If three dimensions are saved, the first set of names are `OVE1_1`, `OVE2_1`, and `OVE3_1`. If another `OVERALS` is then run, the variable names for the second set are `OVE1_2`, `OVE2_2`, `OVE3_2`, and so on.
- Following the name, the number of dimensions for which you want object scores saved can be listed in parentheses. The number cannot exceed the value of the `DIMENSION` subcommand.
- The prefix should be unique for each `OVERALS` command in the same session. Otherwise,, `OVERALS` replaces the prefix with `DIM`, `OBJ`, or `OBSAVE`. If all of these prefixes already exist, `SAVE` is not executed.
- If the number of dimensions is not specified, the `SAVE` subcommand saves object scores for all dimensions.
- If you replace the active dataset by specifying an asterisk (\*) on a `MATRIX` subcommand, the `SAVE` subcommand is not executed.

### **Example**

```
OVERALS CAR1 CAR2 CAR3(5) PRICE(10)
/SET=2(3,1)
/ANALYSIS=CAR1 TO CAR3(SNOM) PRICE(NUM)
/DIMENSIONS=3
/SAVE=DIM(2) .
```

- Analyzed items include three single nominal variables, `CAR1`, `CAR2`, and `CAR3` (each with 5 categories) and one numeric level variable (with 10 categories).
- The `DIMENSIONS` subcommand requests results for three dimensions.
- `SAVE` adds the object scores from the first two dimensions to the active dataset. The names of these new variables will be `DIM00001` and `DIM00002`, respectively.

## **MATRIX Subcommand**

The `MATRIX` subcommand is used to write category quantifications, coordinates, centroids, weights, and component loadings to a matrix data file.

- The specification on `MATRIX` is keyword `OUT` and a quoted file specification or previously declared dataset name (`DATASET DECLARE` command), enclosed in parentheses.
- You can specify an asterisk (\*) instead of a file to replace the active dataset.
- All values are written to the same file.
- The matrix data file has one case for each value of each original variable.

The variables of the matrix data file and their values are as follows:

<b>ROWTYPE_</b>	<i>String variable containing value QUANT for the category quantifications, SCOOR_ for the single-category coordinates, MCOOR_ for multiple-category coordinates, CENTRO_ for centroids, PCENTRO_ for projected centroids, WEIGHT_ for weights, and LOADING_ for the component scores.</i>
<b>LEVEL</b>	<i>String variable containing the values (or value labels, if present) of each original variable for category quantifications. For cases with ROWTYPE_=LOADING_ or WEIGHT_, the value of LEVEL is blank.</i>
<b>VARNAME_</b>	<i>String variable containing the original variable names.</i>
<b>VARTYPE_</b>	<i>String variable containing values MULTIPLE, SINGLE N, ORDINAL, or NUMERICAL, depending on the level of optimal scaling that is specified for the variable.</i>
<b>SET_</b>	<i>The set number of the original variable.</i>
<b>DIM1...DIMn</b>	<i>Numeric variables containing the category quantifications, the single-category coordinates, multiple-category coordinates, weights, centroids, projected centroids, and component loadings for each dimension. Each variable is labeled DIMn, where n represents the dimension number. Any values that cannot be computed are assigned 0 in the file.</i>

# PACF

```
PACF VARIABLES= series names

[/DIFF={1}]
      {n}

[/SDIFF={1}]
      {n}

[/PERIOD=n]

[/{NOLOG**}]
      {LN      }

[/SEASONAL]

[/MXAUTO={16**}]
      {n      }

[/APPLY [= 'model name']]
```

**\*\*Default if the subcommand is omitted and there is no corresponding specification on the TSET command.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
PACF VARIABLES = TICKETS.
```

## **Overview**

PACF displays and plots the sample partial autocorrelation function of one or more time series. You can also display and plot the partial autocorrelations of transformed series by requesting natural log and differencing transformations from within the procedure.

## **Options**

**Modification of the Series.** You can use the LN subcommand to request a natural log transformation of the series, and you can use the SDIFF and DIFF subcommand to request seasonal and nonseasonal differencing to any degree. With seasonal differencing, you can specify the periodicity on the PERIOD subcommand.

**Statistical Output.** With the MXAUTO subcommand, you can specify the number of lags for which you want values to be displayed and plotted, overriding the maximum value that is specified on TSET. You can also use the SEASONAL subcommand to display and plot values only at periodic lags.



### **Basic Specification**

The basic specification is one or more series names. For each specified series, `PACF` automatically displays the partial autocorrelation value and standard error value for each lag. `PACF` also plots the partial autocorrelations and marks the bounds of two standard errors on the plot. By default, `PACF` displays and plots partial autocorrelations for up to 16 lags (or the number of lags that are specified on `TSET`).

### **Subcommand Order**

- Subcommands can be specified in any order.

### **Syntax Rules**

- `VARIABLES` can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each subcommand is executed.

### **Operations**

- Subcommand specifications apply to all series that are named on the `PACF` command.
- If the `LN` subcommand is specified, any differencing that is requested on that `PACF` command is done on log-transformed series.
- Confidence limits are displayed in the plot, marking the bounds of two standard errors at each lag.

### **Limitations**

- A maximum of one `VARIABLES` subcommand is allowed. There is no limit on the number of series that are named on the list.

## **Example**

```
PACF VARIABLES = TICKETS
  /LN
  /DIFF=1
  /SDIFF=1
  /PERIOD=12
  /MXAUTO=25.
```

- This example produces a plot of the partial autocorrelation function for the series `TICKETS` after a natural log transformation, differencing, and seasonal differencing have been applied to the series. Along with the plot, the partial autocorrelation value and standard error are displayed for each lag.
- `LN` transforms the data by using the natural logarithm (base  $e$ ) of the series.
- `DIFF` differences the series once.
- `SDIFF` and `PERIOD` apply one degree of seasonal differencing with a period of 12.
- `MXAUTO` specifies 25 for the maximum number of lags for which output is to be produced.

## **VARIABLES Subcommand**

VARIABLES specifies the series names and is the only required subcommand.

## **DIFF Subcommand**

DIFF specifies the degree of differencing that is used to convert a nonstationary series to a stationary series with a constant mean and variance before the partial autocorrelations are computed.

- You can specify 0 or any positive integer on DIFF.
- If DIFF is specified without a value, the default is 1.
- The number of values that are used in the calculations decreases by 1 for each degree of differencing.

### **Example**

```
PACF VARIABLES = SALES  
/DIFF=1.
```

- In this example, the series *SALES* will be differenced once before the partial autocorrelations are computed and plotted.

## **SDIFF Subcommand**

If the series exhibits a seasonal or periodic pattern, you can use the SDIFF subcommand to seasonally difference the series before obtaining partial autocorrelations. SDIFF indicates the degree of seasonal differencing.

- The specification on SDIFF can be 0 or any positive integer.
- If SDIFF is specified without a value, the default is 1.
- The number of seasons that are used in the calculations decreases by 1 for each degree of seasonal differencing.
- The length of the period that is used by SDIFF is specified on the PERIOD subcommand. If the PERIOD subcommand is not specified, the periodicity that was established on the TSET or DATE command is used (see the PERIOD subcommand).

## **PERIOD Subcommand**

PERIOD indicates the length of the period to be used by the SDIFF or SEASONAL subcommand. PERIOD indicates how many observations are in one period or season.

- The specification on PERIOD can be any positive integer that is greater than 1.
- PERIOD is ignored if it is used without the SDIFF or SEASONAL subcommand.

- If `PERIOD` is not specified, the periodicity that was established on `TSET PERIOD` is in effect. If `TSET PERIOD` is not specified, the periodicity that was established on the `DATE` command is used. If periodicity was not established anywhere, the `SDIFF` and `SEASONAL` subcommands are not executed.

### Example

```
PACF VARIABLES = SALES
  /SDIFF=1
  /PERIOD=12 .
```

- This `PACF` command applies one degree of seasonal differencing with a periodicity of 12 to the series `SALES` before partial autocorrelations are computed and plotted.

## LN and NOLOG Subcommands

`LN` transforms the data by using the natural logarithm (base  $e$ ) of the series and is used to remove varying amplitude over time. `NOLOG` indicates that the data should not be log transformed. `NOLOG` is the default.

- If you specify `LN` on a `PACF` command, any differencing that is requested on that command is performed on the log-transformed series.
- There are no additional specifications on `LN` or `NOLOG`.
- Only the last `LN` or `NOLOG` subcommand on a `PACF` command is executed.
- If a natural log transformation is requested when there are values in the series that are less than or equal to 0, `PACF` will not be produced for that series because nonpositive values cannot be log-transformed.
- `NOLOG` is generally used with an `APPLY` subcommand to turn off a previous `LN` specification.

### Example

```
PACF VARIABLES = SALES
  /LN .
```

- This command transforms the series `SALES` by using the natural log transformation and then computes and plots partial autocorrelations.

## SEASONAL Subcommand

Use `SEASONAL` to focus attention on the seasonal component by displaying and plotting autocorrelations only at periodic lags.

- There are no additional specifications on `SEASONAL`.
- If `SEASONAL` is specified, values are displayed and plotted at the periodic lags that are indicated on the `PERIOD` subcommand. If `PERIOD` is not specified, the periodicity that was established on the `TSET` or `DATE` command is used (see the `PERIOD` subcommand).
- If `SEASONAL` is not specified, partial autocorrelations for all lags (up to the maximum) are displayed and plotted.

**Example**

```
PACF VARIABLES = SALES
  /SEASONAL
  /PERIOD=12 .
```

- In this example, partial autocorrelations are displayed and plotted at every 12th lag.

**MXAUTO Subcommand**

MXAUTO specifies the maximum number of lags for a series.

- The specification on MXAUTO must be a positive integer.
- If MXAUTO is not specified, the default number of lags is the value that was set on TSET MXAUTO. If TSET MXAUTO is not specified, the default is 16.
- The value on MXAUTO overrides the value that was set on TSET MXAUTO.

**Example**

```
PACF VARIABLES = SALES
  /MXAUTO=14 .
```

- This command specifies 14 for the maximum number of partial autocorrelations that can be displayed and plotted for series *SALES*.

**APPLY Subcommand**

APPLY allows you to use a previously defined PACF model without having to repeat the specifications.

- The only specification on APPLY is the name of a previous model enclosed in quotes. If a model name is not specified, the model that was specified on the previous PACF command is used.
- To change one or more model specifications, specify the subcommands of only those portions that you want to change, placing the specifications after the APPLY subcommand.
- If no series are specified on the PACF command, the series that were originally specified with the model that is being reapplied are used.
- To change the series that are used with the model, enter new series names before or after the APPLY subcommand.

**Example**

```
PACF VARIABLES = TICKETS
  /LN
  /DIFF=1
  /SDIFF=1
  /PER=12
  /MXAUTO=25 .
PACF VARIABLES = ROUNTRP
  /APPLY .
```

- The first command specifies a maximum of 25 partial autocorrelations for the series *TICKETS* after it has been log-transformed, differenced once, and had one degree of seasonal differencing with a periodicity of 12 applied to it. This model is assigned the default name *MOD\_1*.
- The second command displays and plots partial autocorrelations for series *ROUNDTRP* by using the same model that was specified for series *TICKETS*.

## ***References***

Box, G. E. P., and G. M. Jenkins. 1976. *Time series analysis: Forecasting and control*, Rev. ed. San Francisco: Holden-Day.

# PARTIAL CORR

PARTIAL CORR is available in the Statistics Base option.

```
PARTIAL CORR VARIABLES= varlist [WITH varlist]
BY varlist [(levels)]

[/SIGNIFICANCE={TWOTAIL**}]
{ONETAIL }

[/STATISTICS={NONE**} [CORR] [DESCRIPTIVES] [BADCORR] [ALL]]

[/FORMAT={MATRIX** } ]
{SERIAL }
{CONDENSED}

[/MISSING={LISTWISE**} [{EXCLUDE**}]]
{ANALYSIS } {INCLUDE }

[/MATRIX= [IN({* }))] [OUT({* }))]
{'savfile'|'dataset'} {'savfile'|'dataset'}
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 16.0

- Added support for SET THREADS and SET MCACHE.

## Example

```
PARTIAL CORR VARIABLES=PUBTRANS MECHANIC BY NETPURSE(1).
```

## Overview

PARTIAL CORR produces partial correlation coefficients that describe the relationship between two variables while adjusting for the effects of one or more additional variables. PARTIAL CORR calculates a matrix of Pearson product-moment correlations. PARTIAL CORR can also read the zero-order correlation matrix as input. Other procedures that produce zero-order correlation matrices that can be read by PARTIAL CORR include CORRELATIONS, REGRESSION, DISCRIMINANT, and FACTOR.

## Options

**Significance Levels.** By default, the significance level for each partial correlation coefficient is based on a two-tailed test. Optionally, you can request a one-tailed test using the SIGNIFICANCE subcommand.

**Statistics.** In addition to the partial correlation coefficient, degrees of freedom, and significance level, you can use the `STATISTICS` subcommand to obtain the mean, standard deviation, and number of nonmissing cases for each variable, as well as zero-order correlation coefficients for each pair of variables.

**Format.** You can specify condensed format, which suppresses the degrees of freedom and significance level for each coefficient, and you can print only nonredundant coefficients in serial string format by using the `FORMAT` subcommand.

**Matrix Input and Output.** You can read and write zero-order correlation matrices by using the `MATRIX` subcommand.

### **Basic Specification**

The basic specification is the `VARIABLES` subcommand, which specifies a list of variables to be correlated, and one or more control variables following keyword `BY`. `PARTIAL CORR` calculates the partial correlation of each variable with every other variable that was specified on the correlation variable list.

### **Subcommand Order**

Subcommands can be specified in any order.

### **Operations**

- `PARTIAL CORR` produces one matrix of partial correlation coefficients for each of up to five order values. For each coefficient, `PARTIAL CORR` prints the degrees of freedom and the significance level.
- This procedure uses the multithreaded options specified by `SET THREADS` and `SET MCACHE`.

### **Limitations**

- A maximum of 25 variable lists on a single `PARTIAL CORR` command is allowed. Each variable list contains a correlation list, a control list, and order values.
- A maximum of 400 variables total can be named or implied per `PARTIAL CORR` command.
- A maximum of 100 control variables is allowed.
- A maximum of 5 different order values per single list is allowed. The largest order value that can be specified is 100.

## **Example**

```
PARTIAL CORR
  /VARIABLES= funding disease BY visits
  /SIGNIFICANCE=TWOTAIL
  /STATISTICS=CORR
  /MISSING=LISTWISE .
```

- The procedure produces partial correlations *funding* with *disease* controlling for *visits*.

- `STATISTICS` requests zero-order correlations in addition to partial correlations.
- All other options are set to their default values.

## VARIABLES Subcommand

`VARIABLES` requires a *correlation list* of one or more pairs of variables for which partial correlations are desired and requires a *control list* of one or more variables that will be used as controls for the variables in the correlation list, followed by optional order values in parentheses.

- The correlation list specifies pairs of variables to be correlated while controlling for the variables in the control list.
- To request a square or lower-triangular matrix, do not use keyword `WITH` in the correlation list. This specification obtains the partial correlation of every variable with every other variable in the list.
- To request a rectangular matrix, specify a list of correlation variables followed by keyword `WITH` and a second list of variables. This specification obtains the partial correlation of specific variable pairs. The first variable list defines the rows of the matrix, and the second list defines the columns.
- The control list is specified after keyword `BY`.
- The correlation between a pair of variables is referred to as a *zero-order correlation*. Controlling for one variable produces a first-order partial correlation, controlling for two variables produces a second-order partial correlation, and so on.
- To indicate the exact partials that are to be computed, you can specify order values in parentheses following the control list. These values also determine the partial correlation matrix or matrices to be printed. Up to five order values can be specified. Separate each value with at least one space or comma. The default order value is the number of control variables.
- One partial is produced for every unique combination of control variables for each order value.

### Obtaining the Partial Correlation for Specific Variable Pairs

```
PARTIAL CORR
  VARIABLES = RENT FOOD PUBTRANS WITH TEACHER MANAGER BY NETSALRY(1).
```

- `PARTIAL CORR` produces a rectangular matrix. Variables *RENT*, *FOOD*, and *PUBTRANS* form the matrix rows, and variables *TEACHER* and *MANAGER* form the columns.

### Specifying Order Values

```
PARTIAL CORR
  VARIABLES = RENT WITH TEACHER BY NETSALRY, NETPRICE (1).
PARTIAL CORR
  VARIABLES = RENT WITH TEACHER BY NETSALRY, NETPRICE (2).
PARTIAL CORR
  VARIABLES = RENT WITH TEACHER BY NETSALRY, NETPRICE (1,2).
PARTIAL CORR
  VARIABLES = RENT FOOD PUBTRANS BY NETSALRY NETPURSE NETPRICE (1,3).
```

- The first `PARTIAL CORR` produces two first-order partials: *RENT* with *TEACHER* controlling for *NETSALRY*, and *RENT* with *TEACHER* controlling for *NETPRICE*.



- The second PARTIAL CORR produces one second-order partial of *RENT* with *TEACHER* controlling simultaneously for *NETSALRY* and *NETPRICE*.
- The third PARTIAL CORR specifies both sets of partials that were specified by the previous two commands.
- The fourth PARTIAL CORR produces three first-order partials (controlling for *NETSALRY*, *NETPURSE*, and *NETPRICE* individually) and one third-order partial (controlling for all three control variables simultaneously).

## **SIGNIFICANCE Subcommand**

SIGNIFICANCE determines whether the significance level is based on a one-tailed or two-tailed test.

- By default, the significance level is based on a two-tailed test. This setting is appropriate when the direction of the relationship between a pair of variables cannot be specified in advance of the analysis.
- When the direction of the relationship can be determined in advance, a one-tailed test is appropriate.

**TWOTAIL**     *Two-tailed test of significance.* This setting is the default.

**ONETAIL**     *One-tailed test of significance.*

## **STATISTICS Subcommand**

By default, the partial correlation coefficient, degrees of freedom, and significance level are displayed. Use STATISTICS to obtain additional statistics.

- If both CORR and BADCORR are requested, CORR takes precedence over BADCORR, and the zero-order correlations are displayed.

**CORR**             *Zero-order correlations with degrees of freedom and significance level.*

**DESCRIPTIVES**     *Mean, standard deviation, and number of nonmissing cases.* Descriptive statistics are not available with matrix input.

**BADCORR**         *Zero-order correlation coefficients only if any zero-order correlations cannot be computed.* Noncomputable coefficients are displayed as a period.

**NONE**             *No additional statistics.* This setting is the default.

**ALL**                *All additional statistics that are available with PARTIAL CORR.*

## **FORMAT Subcommand**

FORMAT determines page format.

- If both CONDENSED and SERIAL are specified, only SERIAL is in effect.

<b>MATRIX</b>	<i>Display degrees of freedom and significance level in matrix format.</i> This format requires four lines per matrix row and displays the degrees of freedom and the significance level. The output includes redundant coefficients. This setting is the default.
<b>CONDENSED</b>	<i>Suppress the degrees of freedom and significance level.</i> This format requires only one line per matrix row and suppresses the degrees of freedom and significance. A single asterisk (*) following a coefficient indicates a significance level of 0.05 or less. Two asterisks (**) following a coefficient indicate a significance level of 0.01 or less.
<b>SERIAL</b>	<i>Display only the nonredundant coefficients in serial string format.</i> The coefficients, degrees of freedom, and significance levels from the first row of the matrix are displayed first, followed by all unique coefficients from the second row and so on for all rows of the matrix.

## **MISSING Subcommand**

MISSING controls the treatment of cases with missing values.

- When pairwise deletion is in effect (keyword ANALYSIS), the degrees of freedom for a particular partial coefficient are based on the smallest number of cases that are used in the calculation of any of the simple correlations.
- LISTWISE and ANALYSIS are alternatives. However, each command can be used with either INCLUDE or EXCLUDE. The default is LISTWISE and EXCLUDE.

<b>LISTWISE</b>	<i>Exclude cases with missing values listwise.</i> Cases with missing values for any of the variables that are listed for an analysis—including control variables—are not used in the calculation of the zero-order correlation coefficient. This setting is the default.
<b>ANALYSIS</b>	<i>Exclude cases with missing values on a pair-by-pair basis.</i> Cases with missing values for one or both of a pair of variables are not used in the calculation of zero-order correlation coefficients.
<b>EXCLUDE</b>	<i>Exclude user-missing values.</i> User-missing values are treated as missing. This setting is the default.
<b>INCLUDE</b>	<i>Include user-missing values.</i> User-missing values are treated as valid values.

## **MATRIX Subcommand**

MATRIX reads and writes matrix data files.

- Either `IN` or `OUT` and a matrix file in parentheses is required. When both `IN` and `OUT` are used on the same `PARTIAL CORR` procedure, they can be specified on separate `MATRIX` subcommands or they can both be specified on the same subcommand.

**OUT ('savfile']'dataset')** *Write a matrix data file or dataset.* Specify either a filename, a previously declared dataset name, or an asterisk, enclosed in parentheses. Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. If you specify an asterisk (\*), the matrix data file replaces the active dataset. If you specify an asterisk or a dataset name, the file is not stored on disk unless you use `SAVE` or `XSAVE`.

**IN ('savfile']'dataset')** *Read a matrix data file or dataset.* Specify either a filename, dataset name created during the current session, or an asterisk enclosed in parentheses. An asterisk reads the matrix data from the active dataset. Filenames should be enclosed in quotes and are read from the working directory unless a path is included as part of the file specification.

### Matrix Output

- The matrix materials that `PARTIAL CORR` writes can be used by subsequent `PARTIAL CORR` procedures or by other procedures that read correlation-type matrices.
- In addition to the partial correlation coefficients, the matrix materials that `PARTIAL CORR` writes include the mean, standard deviation, and number of cases that are used to compute each coefficient (see *Format of the Matrix Data File* on p. 1517 for a description of the file). If `PARTIAL CORR` reads matrix data and then writes matrix materials based on those data, the matrix data file that it writes will not include means and standard deviations.
- `PARTIAL CORR` writes a full square matrix for the analysis that is specified on the `VARIABLES` subcommand (or the analysis list if keyword `VARIABLES` is omitted). No matrix is written for subsequent variable lists.
- Any documents that are contained in the active dataset are not transferred to the matrix file.

### Matrix Input

- When matrix materials are read from a file other than the active dataset, both the active dataset and the matrix data file that is specified on `IN` must contain all variables that are specified on the `VARIABLES` subcommand.
- `MATRIX=IN` cannot be specified unless a active dataset has already been defined. To read an existing matrix data file at the beginning of a session, use `GET` to retrieve the matrix file and then specify `IN (*)` on `MATRIX`.
- `PARTIAL CORR` can read correlation-type matrices written by other procedures.
- The program reads variable names, variable and value labels, and print and write formats from the dictionary of the matrix data file.

### Format of the Matrix Data File

- The matrix data file includes two special variables that are created by the program: `ROWTYPE_` and `VARNAME_`.



- PARTIAL CORR writes the same matrix as in the example above. However, the matrix data file replaces the active dataset. The LIST command is executed on the matrix file, not on the CITY file.

### Using a Matrix Data File as Input

```
GET FILE='/data/personnel.sav'.
FREQUENCIES VARIABLES=AGE.
PARTIAL CORR VARIABLES=BUSDRVER MECHANIC ENGINEER TEACHER COOK
BY NETSALRY(1) /MATRIX=IN('/data/corr_matrix.sav').
```

- This example performs a frequencies analysis on file *personnel.sav* and then uses a different file for PARTIAL CORR. The file is an existing matrix data file.
- MATRIX=IN specifies the matrix data file. Both the active dataset and the *corr\_matrix.sav* file must contain all variables that are specified on the VARIABLES subcommand on PARTIAL CORR.
- The *corr\_matrix.sav* file does not replace *personnel.sav* as the active dataset.

### Using an Active Dataset That Contains Matrix Data

```
GET FILE='/data/corr_matrix.sav'.
PARTIAL CORR VARIABLES=BUSDRVER MECHANIC ENGINEER TEACHER COOK
BY NETSALRY(1)
/MATRIX=IN(*).
```

- The GET command retrieves the matrix data file *corr\_matrix.sav*.
- MATRIX=IN specifies an asterisk because the active dataset is the matrix file *CORMTX*. If MATRIX=IN('/data/corr\_matrix.sav') is specified, the program issues an error message.
- If the GET command is omitted, the program issues an error message.

```
GET FILE='/data/city.sav'.
REGRESSION MATRIX=OUT(*)
/VARIABLES=NETPURSE PUBTRANS MECHANIC BUSDRVER
/DEPENDENT=NETPURSE /ENTER.
PARTIAL CORR
VARIABLES = PUBTRANS MECHANIC BUSDRVER BY NETPURSE(1) /MATRIX=IN(*).
```

- GET retrieves the IBM® SPSS® Statistics data file *city.sav*.
- REGRESSION computes correlations among the specified variables. MATRIX=OUT(\*) writes a matrix data file that replaces the active dataset.

# PERMISSIONS

```
PERMISSIONS FILE='filespec'  
  
  /PERMISSIONS {READONLY }  
                {WRITEABLE}
```

## Example

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

```
PERMISSIONS FILE='/mydir/mydata.sav'  
  /PERMISSIONS READONLY.
```

## Overview

PERMISSIONS changes the read/write permissions for the specified file, using the operating system facilities for changing permissions.

### Syntax Rules

- A FILE specification and a PERMISSIONS subcommand are both required.
- The file specification should be enclosed in single quotation marks or double quotation marks.

## PERMISSIONS Subcommand

<b>READONLY</b>	<i>File permissions are set to read-only for all users.</i> The file cannot be saved by using the same file name with subsequent changes unless the read/write permissions are changed in the operating system or a subsequent PERMISSIONS command specifies PERMISSIONS=WRITEABLE.
<b>WRITEABLE</b>	<i>File permissions are set to allow writing for the file owner.</i> If file permissions were set to read-only for other users, the file remains read-only for them.

Your ability to change the read/write permissions may be restricted by the operating system.

# PLANCARDS

PLANCARDS is available in the Conjoint option.

```
PLANCARDS
  [/FACTORS=varlist]

  [/FORMAT={LIST}]
           {CARD}
           {BOTH}

  [/TITLE='string']

  [/FOOTER='string']

  [/OUTFILE=file]
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- PAGINATE subcommand is obsolete and no longer supported.

## **Example**

```
PLANCARDS /TITLE='Car for Sale'.
```

## **Overview**

PLANCARDS produces profiles, or cards, from a plan file for a conjoint analysis study. The plan file can be generated by ORTHOPLAN or entered by the user. The printed profiles can be used as the experimental stimuli that subjects judge in terms of preference.

## **Options**

**Format.** You can produce profiles in the format of a single list or formatted so that each profile is displayed separately.

**Titles and Footers.** You can specify title and footer labels that appear at the top and bottom of the output (for single list format) or at the top and bottom of each profile (when the profiles are displayed separately).

## **Basic Specification**

- The basic specification is PLANCARDS, which produces a listing of profiles, using all variables in the active dataset except *STATUS\_* and *CARD\_* as factors.

**Subcommand Order**

- Subcommands can be named in any order.

**Operations**

- PLANCARDS assumes that the active dataset represents a plan for a full-profile (full-concept) conjoint study. Each “case” in such a file is one profile in the conjoint experimental plan.
- Factor and factor-level labels in the active dataset—generated by ORTHOPLAN or by the VARIABLE and VALUE LABELS commands—are used in the output.
- The command SPLIT FILE is ignored for single-profile format. In listing format, each subfile represents a different plan, and a new listing begins for each subfile.
- The WEIGHT command is ignored by PLANCARDS.

**Limitations**

- Missing values are not recognized as missing and are treated like other values.

**Examples**

```

ORTHOPLAN FACTORS=SPEED 'Highest possible speed'
  (70 '70 mph' 100 '100 mph' 130 '130 mph')
  WARRANTY 'Length of warranty' ('1 year' '3 year' '5 year')
  SEATS 'Number of seats' (2, 4) /MINIMUM=9 /HOLDOUT=6.
PLANCARDS FORMAT=BOTH
  /TITLE='Car for Sale'.

```

- ORTHOPLAN generates a set of profiles (cases) for a full-profile conjoint analysis in the active dataset.
- PLANCARDS displays the profiles, along with the title *Car for Sale*.

**Example: User-entered Plan**

```

DATA LIST FREE/ COST NEWNESS EXPER NAME REP
              GUARAN TRIAL TRUST.
VARIABLE LABELS
  COST 'Product cost'
  NEWNESS 'Product newness'
  EXPER 'Brand experience'
  NAME "Manufacturer's Name"
  REP "Distributor's reputation"
  GUARAN 'Money-back Guarantee'
  TRIAL 'Free sample/trial'
  TRUST 'Endorsed by a trusted person'.
VALUE LABELS
  COST 1 'LOW' 2 'HIGH'/
  NEWNESS 1 'NEW' 2 'OLD'/
  EXPER 1 'SOME' 2 'NONE'/
  NAME 1 'ESTABLISHED' 2 'UNKNOWN'/
  REP 1 'GOOD' 2 'UNKNOWN'/
  GUARAN 1 'YES' 2 'NO'/
  TRIAL 1 'YES' 2 'NO'/
  TRUST 1 'YES' 2 'NO'.
BEGIN DATA
  1 2 2 1 2 2 2 1
  2 2 2 1 1 1 2 1
  2 2 1 2 2 1 1 1

```



```

2 1 2 1 2 2 1 2
2 1 1 2 2 2 2 1
2 1 2 2 1 1 2 2
1 1 2 2 1 2 1 1
1 1 1 1 2 1 2 2
1 2 1 2 1 2 2 2
1 1 1 1 1 1 1 1
2 2 1 1 1 2 1 2
1 2 2 2 2 1 1 2
END DATA.
PLANCARDS.

```

- In this example, the plan is entered and defined by the user rather than by ORTHOPLAN.
- PLANCARDS uses the information in the active dataset to produce a set of profiles. Because no format is specified, the default format (single list) is used. The variables and values in this example were taken from Akaah & Korgaonkar (Akaah and Korgaonkar, 1988).

## ***FACTORS Subcommand***

FACTORS identifies the variables to be used as factors and the order in which their labels are to appear in the output. String variables are permitted.

- Keyword FACTORS is followed by a variable list.
- By default, if FACTORS is not specified, all variables in the active dataset (except those variables that are named *STATUS\_* or *CARD\_*) are used as factors in the order in which they appear in the file. (See the ORTHOPLAN command for information about variables *STATUS\_* and *CARD\_*.)

## ***FORMAT Subcommand***

FORMAT specifies how the profiles should be displayed. The choices are **listing** format (LIST keyword) and **single-profile** format (CARD keyword). Listing format displays the profiles in the form of a single list. For single-profile format, output is displayed so that each profile is presented separately.

- The keyword FORMAT is followed by LIST, CARD, or BOTH. (ALL is an alias for BOTH.)
- The default format is LIST.
- With LIST format, holdout profiles are differentiated from experimental profiles, and simulation profiles are listed separately following the experimental and holdout profiles. With CARD format, holdout profiles are not differentiated, and simulation profiles are not produced.
- If FORMAT=LIST is specified along with the OUTFILE subcommand, the OUTFILE subcommand is ignored (OUTFILE only applies to CARD format). Specifying OUTFILE with FORMAT=BOTH is equivalent to OUTFILE with FORMAT=CARD.

### ***Example***

```

PLANCARDS FORMAT=CARD /OUTFILE='DESIGN.FRM'
/TITLE=' ' 'Profile #)CARD' /FOOTER='RANK:'.

```

- FORMAT=CARD specifies that the output will be in single-profile format.

- The profiles are written to the file *DESIGN.FRM*.
- Each profile in *DESIGN.FRM* will have the title Profile #*n* at the top and the label RANK: at the bottom, where *n* is a profile identification number.

The output for the first two profiles is shown below.

Figure 178-1

*Single-profile format*

```

Profile #1

Product cost  LOW
Product newness  OLD
Brand experience  NONE
Manufacturer's Name  ESTABLISHED
Distributor's reputation  UNKNOWN
Money-back Guarantee  NO
Free sample/trial  NO
Endorsed by a trusted person  YES

RANK:

Profile #2

Product cost  HIGH
Product newness  OLD
Brand experience  NONE
Manufacturer's Name  ESTABLISHED
Distributor's reputation  GOOD
Money-back Guarantee  YES
Free sample/trial  NO
Endorsed by a trusted person  YES

RANK:

. . .

```

## ***OUTFILE Subcommand***

OUTFILE names an external file where profiles in single-profile format are to be written. Profiles in listing format are not written to an external file.

- By default, no external file is written.
- The OUTFILE keyword is followed by the name of an external file. The file is specified in the usual manner for your system.
- If the OUTFILE subcommand is specified along with FORMAT=LIST, the OUTFILE subcommand is ignored (OUTFILE only applies to FORMAT=CARD).

## ***TITLE Subcommand***

TITLE specifies a string to be used at the top of the output (in listing format) or at the top of each new profile (in single-profile format).

- Default titles are provided, except for output that is directed to an external file with the OUTFILE subcommand.
- The keyword TITLE is followed by a string enclosed in apostrophes.

- Quotation marks can be used to enclose the string instead of apostrophes when you want to use an apostrophe in the title.
- Multiple strings per TITLE subcommand can be specified; each string will appear on a separate line.
- Use an empty string ( ' ' ) to cause a blank line.
- Multiple TITLE subcommands can be specified; each subcommand will appear on a separate line.
- If the special character sequence )CARD is specified anywhere in the title, PLANCARDS will replace it with the sequential profile number in single-profile-formatted output. This character sequence is not translated in listing format.

## **FOOTER Subcommand**

FOOTER specifies a string to be used at the bottom of the output (in listing format) or at the bottom of each profile (in single-profile format).

- If FOOTER is not used, nothing appears after the last attribute.
- FOOTER is followed by a string enclosed in apostrophes.
- Quotation marks can be used to enclose the string instead of apostrophes when you want to use an apostrophe in the footer.
- Multiple strings per FOOTER subcommand can be specified; each string will appear on a separate line.
- Use an empty string ( ' ' ) to cause a blank line.
- Multiple FOOTER subcommands can be specified; each subcommand will appear on a separate line.
- If the special character sequence )CARD is specified anywhere in the footer, PLANCARDS will replace it with the sequential profile number in single-profile-formatted output. This character sequence is not translated in listing format.

### **Example**

```
PLANCARDS
  TITLE='Profile # )CARD' ' '
    'Circle the number in the scale at the bottom that
    'indicates how likely you are to purchase this item.' ' '
/FOOTER= '0  1  2  3  4  5  6  7  8  9 10'
          'Not at all      May or may      Certainly'
          'likely to      not          would'
          'purchase      purchase    purchase'
          '-----'
/FORMAT=CARD
/OUTFILE='DESIGN.FRM' .
```

The above example would produce the following output, in *DESIGN.FRM*, for the first profile:

**Figure 178-2**

*Footer with multiple strings*

Profile # 1

Circle the number in the scale at the bottom that indicates how likely you are to purchase this item.

Product cost   LOW  
Product newness   OLD  
Brand experience   NONE  
Manufacturer's Name   ESTABLISHED  
Distributor's reputation   UNKNOWN  
Money-back Guarantee   NO  
Free sample/trial   NO  
Endorsed by a trusted person   YES

0	1	2	3	4	5	6	7	8	9	10
Not at all				May or may				Certainly		
likely to				not				would		
purchase				purchase				purchase		

-----

# PLS

PLS is available in the Statistics Base option.

PLS is an extension command that requires the Python Extension Module to be installed on the system where you plan to run PLS. The PLS Extension Module must be installed separately and the installer can be downloaded from <http://www.spss.com/devcentral>.

*Note:* The PLS Extension Module is dependent upon Python software. SPSS Inc. is not the owner or licensor of the Python software. Any user of Python must agree to the terms of the Python license agreement located on the Python Web site. SPSS Inc. is not making any statement about the quality of the Python program. SPSS Inc. fully disclaims all liability associated with your use of the Python program.

```
PLS dependent variable [MLEVEL={N}] [REFERENCE={FIRST }]
                        {O}           {LAST**}
                        {S}           {value }
    [dependent variable...]
    [BY factor list] [WITH covariate list]

[/ID VARIABLE = variable]

[/MODEL effect [...effect]]

[/OUTDATASET [CASES=IBM SPSS Statistics dataset]
             [LATENTFACTORS=IBM SPSS Statistics dataset]
             [PREDICTORS=IBM SPSS Statistics dataset]]

[/CRITERIA  LATENTFACTORS={5**   }]
             {integer}
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- Command introduced.

## **Example**

```
PLS Response WITH Price.
```

## **Overview**

The PLS procedure estimates partial least squares regression models. Partial least squares is a predictive technique that is an alternative to ordinary least squares (OLS) regression, canonical correlation, or structural equation modeling for analysis of systems of independent and response

variables. It is particularly useful when predictor variables are highly correlated or when the number of predictors exceeds the number of cases.

PLS combines features of principal components analysis and multiple regression. It first extracts a set of latent factors that explains as much of the covariance as possible between the independent and dependent variables. Then a regression step predicts values of the dependent variables using the decomposition of the independent variables.

Partial least squares regression is also known as “Projection to Latent Structure”.

### **Options**

**Response Variables.** PLS estimates univariate and multivariate models. If you specify one or more categorical dependent variables, a classification model is estimated. If you specify one or more scale dependent variables, a regression model is estimated. Mixed regression and classification models are supported.

**Predictors.** Predictors can be categorical or continuous variables. Both main effects and interaction terms can be estimated.

**Method.** You can specify the maximum number of latent factors to extract. By default, five latent factors are extracted.

**Export.** You can save casewise, factorwise, and predictorwise model results to IBM® SPSS® Statistics datasets.

### **Basic Specification**

- PLS is an extension command and will not be recognized by the system until you use the `EXTENSION` command to add PLS to the command table. The syntax diagram for PLS is defined in `plscommand.xml`, which is installed in the `\extensions` subdirectory of the main installation directory. [For more information, see the topic EXTENSION on p. 702.](#)
- The minimum specification is one or more dependent variables and one or more predictors.
- The procedure displays the following tables: proportion of variance explained (by latent factor), latent factor weights, latent factor loadings, independent variable importance in projection (VIP), and regression parameter estimates (by dependent variable).

### **Operations**

- All model variables are centered and standardized, including indicator variables representing categorical variables.
- If a `WEIGHT` variable is specified, its values are used as frequency weights. Weight values are rounded to the nearest whole number before use. Cases with missing weights or weights less than 0.5 are not used in the analyses.
- User- and system-missing values are treated as invalid.
- Memory allocated via `SET WORKSPACE` is unavailable to extension commands; when running PLS on large datasets, you may actually *lower* the size of your workspace.

**Syntax Rules**

- The PLS command is required. All subcommands are optional.
- Only a single instance of each subcommand is allowed.
- An error occurs if an attribute or keyword is specified more than once within a subcommand.
- Equals signs and parentheses shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.
- Empty subcommands are not allowed.

**Examples**

```
PLS Response WITH Price.
```

- PLS estimates a univariate prediction model that regresses *Response* on *Price*.
- Measurement level is not explicitly specified for dependent variable *Response*. Thus, if *Response* is categorical (nominal or ordinal), a classification model is estimated; otherwise, if *Response* is scale, a regression model is estimated.
- Variable *Price* is treated as a continuous predictor (covariate).

**Classification Model (Explicitly Specified)**

```
PLS Response MLEVEL=N REFERENCE=5 BY Region WITH Price.
```

- PLS estimates a univariate prediction model.
- Since dependent variable *Response* is specified as nominal (N), a classification model is estimated. Value 5 is used as the reference category.
- Variable *Region* is treated as a categorical predictor and *Price* is treated as a continuous predictor (covariate). By default, a main effects model is estimated.

**Multivariate Model**

```
PLS Q1 MLEVEL=S Q2 MLEVEL=S BY Region Edlevel
  /MODEL Region Edlevel Region*Edlevel.
```

- PLS estimates a multivariate regression model.
- Both dependent variables, *Q1* and *Q2*, are treated as scale (continuous) variables.
- Variables *Region* and *Edlevel* are treated as categorical predictors.
- MODEL specifies a fully factorial ANOVA model that includes main effects for *Region* and *Edlevel* as well as the interaction between the two predictors.

**Requesting optional plots**

```
PLS insales MLEVEL=S BY type WITH price engine_s horsepow wheelbas width
  length curb_wgt fuel_cap mpg
  /CRITERIA LATENTFACTORS=5
  /OUTDATASET CASES=indvCases LATENTFACTORS=latentFactors
  PREDICTORS=indepVars.
```

- The procedure fits a partial least squares regression model to the scale variable *lnsales*, using *type* as a factor and *price* through *mpg* as covariates.
- The `OUTDATASET` subcommand requests that output variables related to individual cases, latent factors, and predictors be saved to new datasets *indvCases*, *latentFactors*, and *indepVars*, respectively. These selections will also produce plots of latent factor scores, latent factor weights, and variable importance to projection (VIP) by latent factor.
- All other options are set to their default values.

## Variable Lists

The command line variable lists specify the dependent variables, any categorical predictors (factors), and any scale predictors (covariates). `ALL` and `TO` keywords are supported in all variable lists.

### Dependent Variables

A list of one or more dependent variables must be the first specification on the `PLS` command.

- Dependent variables can be numeric or string.
- If a dependent variable is specified more than once, only the last specification is honored for the variable.

### MLEVEL Keyword

By default, the measurement level recorded in the data dictionary is honored for dependent variables.

- The `MLEVEL` keyword can be specified after the name of any dependent variable to override its dictionary level of measurement. Specify `N` (nominal), `O` (ordinal), or `S` (scale). Note that the `PLS` procedure does not distinguish between ordinal and nominal dependent variables; it treats both as categorical.
- If `MLEVEL` is not specified and no setting is recorded in the data dictionary, then a numeric variable is treated as scale and a string variable is treated as categorical.
- A string dependent variable may be defined as ordinal or nominal only.
- `MLEVEL` applies only to the variable that immediately precedes it in the dependent variable list. For example, `Age Income [S]` and `Age TO Income [S]` assign the scale level of measurement to *Income* only. The keyword `ALL` followed by a measurement level specification applies that measurement level to all dependent variables.

### REFERENCE Keyword

The `REFERENCE` keyword specifies the value of the dependent variable to use as the reference category for parameter estimation. `REFERENCE` can be specified after the name of any dependent variables but is honored only for categorical dependent variables. It is silently ignored for scale variables.



- Specify `FIRST`, `LAST`, or a value. If you specify `FIRST`, the lowest value is treated as the reference category. `LAST` (the default) treats highest value of the dependent variable as the reference category.
- If you specify a value it must exist in the data and be consistent with the type and format of the dependent variable. String and formatted (for example, date or time) values must be quoted.
- `REFERENCE` applies only to the variable that immediately precedes it in the dependent variable list. The keyword `ALL` followed by a `REFERENCE` specification applies that reference category to all categorical dependent variables.

### ***Predictor Variables***

- At least one predictor must be specified.
- The names of the categorical predictors, if any, must be preceded by the keyword `BY`. If keyword `BY` is specified with no factors, then a warning is issued and `BY` is ignored.
- The names of the covariates, if any, must be preceded by the keyword `WITH`. If keyword `WITH` is specified with no covariates, then a warning is issued and `WITH` is ignored.
- Categorical predictor variables can be numeric or string.
- Covariates must be numeric.
- If the dependent variable is specified within a categorical predictor list or a covariate list, then it is ignored in the list.
- All variables specified within a categorical predictor or covariate list must be unique. If duplicate variables are specified within a list, then the duplicates are ignored.
- If duplicate variables are specified across the predictor lists, then an error is issued.

## ***ID Subcommand***

The `ID` subcommand specifies a variable that is used as a unique identifier in casewise output and saved datasets (see the [OUTDATASET Subcommand](#)). If this option is not specified, case number is used as the identifier.

Specify a string or numeric variable. An error is issued if the identifier variable is specified as a dependent or predictor variable.

## ***MODEL Subcommand***

The `MODEL` subcommand is used to specify model effects.

- Specify a list of terms to be included in the model.
- If the `MODEL` subcommand is not specified, the default model includes main effects for any categorical predictors and covariates.
- To include a term for the main effect of a categorical predictor or covariate, enter the name of the variable.

- To include a term for an interaction among predictors, use the keyword `BY` or the asterisk (\*) to join the variables involved in the interaction. For example, `A*B` means a two-way interaction effect of A and B. `A*A` is not allowed because variables inside an interaction effect must be distinct.
- Interactions among covariates are valid, as are interactions between categorical predictors and covariates.
- Nested terms are not supported in the `PLS` procedure.

## ***OUTDATASET Subcommand***

The `OUTDATASET` subcommand requests output datasets and plots for model estimates. You can obtain casewise, factorwise, and predictorwise results. None are produced by default.

### ***CASES Keyword***

`CASES` saves the following casewise model estimates: predicted values, residuals, distance to latent factor model, and latent factor scores. It also plots latent factor scores.

Specify the name of the IBM® SPSS® Statistics dataset. The dataset name must be unique within an invocation of the `PLS` procedure. If you specify the name of an existing dataset, its contents are replaced; otherwise, a new dataset is created. An error is generated if an external file is specified.

### ***LATENTFACTORS Keyword***

`LATENTFACTORS` saves latent factor loadings and latent factor weights. It also plots latent factor weights.

Specify the name of the SPSS Statistics dataset. The dataset name must be unique within an invocation of the `PLS` procedure. If you specify the name of an existing dataset, its contents are replaced; otherwise, a new dataset is created. An error is generated if an external file is specified.

### ***PREDICTORS Keyword***

`PREDICTORS` saves regression parameter estimates and variable importance to projection (VIP). It also plots VIP by latent factor.

Specify the name of the SPSS Statistics dataset. The dataset name must be unique within an invocation of the `PLS` procedure. If you specify the name of an existing dataset, its contents are replaced; otherwise, a new dataset is created. An error is generated if an external file is specified.

## ***CRITERIA Subcommand***

The `CRITERIA` subcommand specifies model estimation criteria.

***LATENTFACTORS Keyword***

The `LATENTFACTORS` keyword specifies an upper limit on the number of latent factors that are extracted. By default, a maximum of five factors are extracted. The value must be a positive integer.

The number of latent factors that is actually extracted may be fewer than the number requested. For example, it is limited by the number of cases and predictors (whichever is smaller).

# PLUM

PLUM is available in the Statistics Base option.

```
PLUM dependent variable [BY factor varlist] [WITH covariate varlist]

[/CRITERIA = [CIN({95** })] [DELTA({0** })] [MXITER({100**})] [MXSTEP({5**})]
             {value}           {value } {n }           {n }
             [LCONVERGE({0** })] [PCONVERGE({1.0E-6**})] [SINGULAR({1.0E-8**})]
             {value}           {value }           {value }           {value }
             [BIAS] ]

[/LINK = {CAUCHIT}]
         {CLOGLOG}
         {LOGIT**}
         {NLOGLOG}
         {PROBIT }

[/LOCATION = [effect effect ...] ]

[/MISSING = {EXCLUDE**}]
           {INCLUDE }

[/PRINT = [CELLINFO] [CORB] [COVB] [FIT] [HISTORY({1})] [KERNEL]
          {n}
          [TPARALLEL] [PARAMETER] [SUMMARY]]

[/SAVE = [ESTPROB [(rootname [{25**})] ] [PREDCAT [(newname)] ] [PCPROB [(newname)] ]
         {n } [ACPROB [(newname)] ] ]

[/SCALE = [effect effect ...] ]

[/TEST [(valuelist)] = ['label'] effect valuelist [effect valuelist] ...;
       [effect valuelist [effect valuelist] ...;] ... ]

[/TEST [(valuelist)] = ['label'] ALL list; [ALL list;] ... ].
```

\*\* Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
PLUM chist.
```

## Overview

This procedure makes use of a general class of models to allow you to analyze the relationship between a polytomous ordinal dependent variable and a set of predictors. These models utilize the ordinal nature of the dependent variable and eliminate the need for rescaling.

## Options

**Link Functions.** Five link functions are available for specifying the model with the LINK subcommand.

**Tuning the Algorithm.** You can control the values of algorithm-tuning parameters with the `CRITERIA` subcommand.

**Optional Output.** You can request additional output through the `PRINT` subcommand.

### **Basic Specification**

The basic specification is one dependent variable.

### **Syntax Rules**

- A minimum of one dependent variable must be specified.
- The variable specification must come first and can be specified only once.
- Subcommands can be specified in any order.
- When subcommands (except the `TEST` subcommand) are repeated, previous specifications are discarded and the last subcommand is in effect.
- Empty subcommands (except the `LOCATION` and the `SCALE` subcommands) are ignored. An empty `LOCATION` or `SCALE` subcommand indicates a simple additive model.
- The words `BY`, `WITH`, and `WITHIN` are reserved keywords in this procedure.

## **Example**

```
PLUM
  chist BY numcred othnstal housng WITH age duration
  /CRITERIA = CIN(95) DELTA(0) LCONVERGE(0) MXITER(100) MXSTEP(5)
  PCONVERGE(1.0E-6) SINGULAR(1.0E-8)
  /LINK = CLOGLOG
  /PRINT = FIT PARAMETER SUMMARY TPARALLEL
  /SAVE = PREDCAT .
```

- *chist* is the dependent variable, *numcred*, *othnstal*, and *housing* are factors, and *age* and *duration* are covariates.
- `LOCATION` is not specified, so by default the location model is based on the main effects of all factors and covariates.
- `CRITERIA` specifies that the confidence level to use is 95, no delta value should be added to cells with observed zero frequency, and neither the log-likelihood nor parameter estimates convergence criteria should be used. This means that the procedure will stop when either 100 iterations or 5 step-halving operations have been performed.
- `LINK` specifies that the complementary log-log function should be used.
- `PRINT` specifies that the goodness-of-fit statistics, parameter statistics, model summary, and test of parallel lines should be displayed.
- `SAVE` specifies that the predicted category for each case should be saved to the active dataset.

## **Variable List**

The variable list specifies the dependent variable, factors, and covariates in the model.

- The dependent variable must be the first specification on the command line.

- The dependent variable is assumed to be an ordinal variable and can be of any type (numeric versus string). The order is determined by sorting the level of the dependent variable in ascending order. The lowest value defines the first category.
- Factor variables can be of any type (numeric versus string). Factor levels are sorted in ascending order. The lowest value defines the first category.
- Covariate variables must be numeric.
- Names of the factors follow the dependent variable separated by the keyword `BY`.
- Enter the covariates, if any, following the factors. Use the keyword `WITH` to separate covariates from factors (if any) and the dependent variable.

## **Weight Variable**

- If a `WEIGHT` variable is specified, this procedure will take the non-missing weight values, rounded to the nearest integer, as frequencies.
- Cases with negative frequencies are always excluded.

## **CRITERIA Subcommand**

The `CRITERIA` subcommand offers controls on the iterative algorithm used for estimation, specifies numerical tolerance for checking singularity, and offers options to customize your output.

<b>BIAS</b>	<i>Bias value added to all observed cell frequencies.</i> Specify a non-negative value less than 1. The default value is 0.0.
<b>CIN</b>	<i>Confidence interval level.</i> Specify a value greater than or equal to 0 and less than 100. The default value is 95.
<b>DELTA</b>	<i>Delta value added to observed zero frequency.</i> Specify a non-negative value less than 1. The default value is 0.0.
<b>LCONVERGE</b>	<i>Log-likelihood function convergence criterion.</i> Convergence is assumed if the absolute change or relative change in the log-likelihood function is less than this value. The criterion is not used if the value is 0. Specify a non-negative value. The default value is 0.
<b>MXITER</b>	<i>Maximum number of iterations.</i> Specify a non-negative integer. The default value is 100. Specifying 0 gives the initial estimates.
<b>MXSTEP</b>	<i>Maximum step-halving allowed.</i> Specify a positive integer. The default value is 5.
<b>PCONVERGE</b>	<i>Parameter estimates convergence criterion.</i> Convergence is assumed if the maximum absolute change in each of the parameter estimates is less than this value. The criterion is not used if the value is 0. Specify a non-negative value. The default value is $10^{-6}$ .
<b>SINGULAR</b>	<i>Value used as tolerance in checking singularity.</i> Specify a positive value. The default value is $10^{-8}$ .

## **LINK Subcommand**

The `LINK` subcommand offers five link functions to specify the model.

- If `LINK` is not specified, `LOGIT` is the default.
- The five keywords are mutually exclusive. Only one of them can be specified and only once.

<b>CAUCHIT</b>	<i>Cauchit function.</i> $f(x) = \tan(\pi(x - 0.5))$ .
<b>CLOGLOG</b>	<i>Complementary log-log function.</i> $f(x) = \log(-\log(1 - x))$ .
<b>LOGIT</b>	<i>Logit function.</i> $f(x) = \log(x / (1 - x))$ . This is the default link function.
<b>NLOGLOG</b>	<i>Negative log-log function.</i> $f(x) = -\log(-\log(x))$ .
<b>PROBIT</b>	<i>Probit function.</i> $f(x) = \Phi^{-1}(x)$ , where $\Phi^{-1}$ is the inverse standard normal cumulative distribution function.

## LOCATION Subcommand

The `LOCATION` subcommand specifies the location model.

- Specify a list of terms to be included in the location model, separated by commas or spaces.
- The default location model is generated if the subcommand is not specified or empty. The default model contains the intercept, all of the covariates (if specified) in the order in which they are specified, and all of the main factorial effects in the order in which they are specified on the variable list.
- To include the intercept term explicitly, enter the keyword `INTERCEPT` on the subcommand.
- To include a main effect term, enter the name of the factor on the subcommand.
- To include an interaction effect term among factors, use the keyword `BY` or the asterisk (\*) to join factors involved in the interaction. For example, `A*B*C` means a three-way interaction effect of A, B, and C, where A, B, and C are factors. The expression `A BY B BY C` is equivalent to `A*B*C`. Factors inside an interaction effect must be distinct. Expressions such as `A*C*A` and `A*A` are invalid. The keyword `INTERCEPT` cannot be used to construct an interaction term.
- To include a nested effect term, use the keyword `WITHIN` or a pair of parentheses on the subcommand. For example, `A(B)` means that A is nested within B, where A and B are factors. The expression `A WITHIN B` is equivalent to `A(B)`. Factors inside a nested effect must be distinct. Expressions such as `A(A)` and `A(B*A)` are invalid.
- Multiple level nesting is supported. For example, `A(B(C))` means that B is nested within C, and A is nested within B(C). When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.
- Nesting within an interaction effect is valid. For example, `A(B*C)` means that A is nested within B\*C.
- Interactions among nested effects are allowed. The correct syntax is the interaction followed by the common nested effect inside the parentheses. For example, interaction between A and B within levels of C should be specified as `A*B(C)` instead of `A(C)*B(C)`.
- To include a covariate term in the model, enter the name of the covariate on the subcommand.
- Covariates can be connected, but not nested, using the keyword `BY` or the asterisk (\*) operator. For example, `X*X` is the product of X and itself. This is equivalent to a covariate whose values are the square of those of X. On the contrary, `X(Y)` is invalid.

- Factor and covariate effects can be connected in many ways. No effects can be nested within a covariate effect. Suppose A and B are factors and X and Y are covariates. Examples of valid combination of factor and covariate effects are A\*X, A\*B\*X, X(A), X(A\*B), X\*A(B), X\*Y(A\*B), and A\*B\*X\*Y.

### Example

```
PLUM
  chist BY numcred othnstal
  /CRITERIA = CIN(95) DELTA(0) LCONVERGE(0) MXITER(100) MXSTEP(5)
             PCONVERGE(0)
  /LOCATION = numcred othnstal numcred*othnstal.
```

- LOCATION specifies that the location model consists of *numcred*, *othnstal*, and their interaction effect.

## MISSING Subcommand

By default, cases with missing values for any of the variables on the variable list are excluded from the analysis. The MISSING subcommand allows you to include cases with user-missing values.

- If MISSING is not specified, the default is EXCLUDE.
- Listwise deletion is always used in this procedure.
- Keywords EXCLUDE and INCLUDE are mutually exclusive. Only one of them can be specified and only once.

**EXCLUDE**            *Exclude both user-missing and system-missing values.* This is the default.  
**INCLUDE**            *User-missing values are treated as valid.* System-missing values cannot be included in the analysis.

## PRINT Subcommand

The PRINT subcommand controls the display of optional output. If no PRINT subcommand is specified, default output includes a case-processing summary table.

**CELLINFO**            *Cell information.* Observed and expected frequencies by category and cumulative, Pearson residual for cumulative and category frequencies, and observed and expected probabilities of each response category separately and cumulatively by covariate pattern combination.

**CORB**                *Asymptotic correlation matrix of the parameter estimates.*

**COVB**                *Asymptotic covariance matrix of the parameter estimates.*

**FIT**                  *Goodness-of-fit statistics.* The Pearson chi-square and the likelihood-ratio chi-square statistics. The statistics are computed based on the classification specified on the variable list.

**HISTORY**            *Iteration history.* The table contains log-likelihood function value and parameter estimates every *n* iterations. The default value is *n* = 1. The first and the last iterations are always printed if HISTORY is specified and regardless of the value of *n*.

**KERNEL**              *Use the kernel of the log-likelihood function for display instead of the complete log-likelihood function.*



<b>TPARALLEL</b>	<i>Test of parallel lines assumption.</i> Produce a chi-squared score test of the parallel lines assumption.
<b>PARAMETER</b>	<i>Parameter statistics.</i> The parameter estimates, the standard errors, the significances, and the confidence interval.
<b>SUMMARY</b>	<i>Model summary.</i> The Cox & Snell's $R^2$ , the Nagelkerke's $R^2$ , and the McFadden's $R^2$ statistics.

## SAVE Subcommand

The SAVE subcommand puts casewise post-estimation statistics back into the active file.

- The new variables must have valid variable names that are not in use in the working file.
- The rootname must be a valid variable name.
- The new variables are saved to the working file in the order the keywords are specified on the subcommand.

<b>ESTPROB</b>	<i>Estimated probabilities of classifying a factor/covariate pattern into the response categories.</i> The predicted probabilities of the first $n$ categories are saved. The default number of categories is 25. To specify a number of categories without a rootname, put a colon before the number.
<b>PREDCAT</b>	<i>The response category that has the maximum expected probability for a factor/covariate pattern.</i>
<b>PCPROB</b>	<i>Estimated probability of classifying a factor/covariate pattern into the predicted category.</i> This probability is the maximum of the estimated probabilities of the factor/covariate pattern.
<b>ACPROB</b>	<i>Estimated probability of classifying a factor/covariate pattern into the actual category.</i>

### Example

```
PLUM
  chist BY numcred othnstal
  /CRITERIA = CIN(95) DELTA(0) LCONVERGE(0) MXITER(100) MXSTEP(5)
  PCONVERGE(0)
  /SAVE = ACPROB(correct) PRPROB.
```

- SAVE specifies that the estimated probabilities of correctly classifying each case should be saved to the variable *correct*. The estimated probabilities of classifying each case into the predicted category are saved to the default variable *pcp\_k*, where  $k$  is the smallest integer for which *pcp\_k* does not already exist.

## SCALE Subcommand

The SCALE subcommand specifies the scale component in the model.

- Specify a list of terms to be included in the model, separated by commas or spaces.
- The model will have no scale component if the subcommand is omitted.
- No scale component is generated if the subcommand is not specified or empty.
- To include a main effect term, enter the name of the factor on the subcommand.
- The keyword INTERCEPT is not allowed on the subcommand.

- To include an interaction effect term among factors, use the keyword `BY` or the asterisk (\*) to join factors involved in the interaction. For example, `A*B*C` means a three-way interaction effect of A, B, and C, where A, B, and C are factors. The expression `A BY B BY C` is equivalent to `A*B*C`. Factors inside an interaction effect must be distinct. Expressions such as `A*C*A` and `A*A` are invalid.
- To include a nested effect term, use the keyword `WITHIN` or a pair of parentheses on the subcommand. For example, `A(B)` means that A is nested within B, where A and B are factors. The expression `A WITHIN B` is equivalent to `A(B)`. Factors inside a nested effect must be distinct. Expressions such as `A(A)` and `A(B*A)` are invalid.
- Multiple level nesting is supported. For example, `A(B(C))` means that B is nested within C, and A is nested within B(C). When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus, `A(B)(C)` is not valid.
- Nesting within an interaction effect is valid. For example, `A(B*C)` means that A is nested within B\*C.
- Interactions among nested effects are allowed. The correct syntax is the interaction followed by the common nested effect inside the parentheses. For example, interaction between A and B within levels of C should be specified as `A*B(C)` instead of `A(C)*B(C)`.
- To include a covariate term in the model, enter the name of the covariate on the subcommand.
- Covariates can be connected, but not nested, using the keyword `BY` or the asterisk (\*) operator. For example, `X*X` is the product of X and itself. This is equivalent to a covariate whose values are the square of those of X. On the contrary, `X(Y)` is invalid.
- Factor and covariate effects can be connected in many ways. No effects can be nested within a covariate effect. Suppose A and B are factors, and X and Y are covariates. Examples of valid combination of factor and covariate effects are `A*X`, `A*B*X`, `X(A)`, `X(A*B)`, `X*A(B)`, `X*Y(A*B)`, and `A*B*X*Y`.

## ***TEST Subcommand***

The `TEST` subcommand allows you to customize your hypothesis tests by directly specifying null hypotheses as linear combinations of parameters.

- `TEST` is offered only through syntax.
- Multiple `TEST` subcommands are allowed. Each is handled independently.
- The basic format of the `TEST` subcommand is an optional list of values enclosed in a pair of parentheses, an optional label in quotes, an effect name or the keyword `ALL`, and a list of values.
- To specify the coefficient for the intercept, use the keyword `INTERCEPT`. The number of values after `INTERCEPT` must be equal to the number of response categories minus 1.
- When multiple linear combinations are specified within the same `TEST` subcommand, a semicolon terminates each linear combination, except the last one.
- The linear combinations are separately tested for each category of the dependent variable and then simultaneously tested for all the categories.

- If specified, the value list that immediately follows the subcommand name is the constant that the linear combinations are equated to under the null hypotheses. If this value list is omitted, the constants are assumed to be all zeros.
- The optional label is a string with a maximum length of 255 bytes. Only one label per TEST subcommand can be specified.
- Only valid effects appearing or implied on the LOCATION or the SCALE subcommands can be specified in a linear combination. If an effect appears in both subcommands, then enter the effect only once on the TEST subcommand.
- To specify coefficient for the intercept, use the keyword INTERCEPT. Only one value is expected to follow INTERCEPT.
- The number of values following an effect name must equal the number of parameters (including the redundant ones) corresponding to that effect. For example, if the effect A\*B takes up six parameters, then exactly six values must follow A\*B.
- A number can be specified as a fraction with a positive denominator. For example, 1/3 or -1/3 are valid, but 1/-3 is invalid.
- When ALL is specified, only a list of values can follow. The number of values must equal the combined number of LOCATION and SCALE parameters (including the redundant ones).
- Effects appearing or implied on the LOCATION or the SCALE subcommands but not specified on the TEST are assumed to take the value 0 for all their parameters.
- Effect names and the ALL keywords are mutually exclusive within a single TEST subcommand.
- If ALL is specified for the first row in a TEST matrix, then all subsequent rows should begin with the ALL keyword.
- If effects are specified for the first row in a TEST matrix, then all subsequent rows should use effect name (thus ALL is not allowed).

### Example

```

PLUM
  chist BY housng
  /CRITERIA = CIN(95) DELTA(0) LCONVERGE(0) MXITER(100) MXSTEP(5)
             PCONVERGE(1.0E-6) SINGULAR(1.0E-8)
  /LINK = CLOGLOG
  /PRINT = CELLINFO CORB COVB FIT HISTORY(1) PARAMETER
             SUMMARY TPARALLEL
  /TEST(0 0) = ALL 1 -1 0 0 0 0 0;
               ALL 0 0 1 -1 0 0 0.

```

- There are a total of seven parameter coefficients in the model; four for the thresholds, and three for the factor *housng*. TEST specifies two separate tests: one in which the first and second thresholds are tested for equality, and one in which the third and fourth thresholds are tested for equality.

# POINT

```
POINT KEY=varname  
  [FILE='file'] [ENCODING='encoding specification']
```

## **Release History**

Release 16.0

- ENCODING subcommand added for Unicode support.

## **Example**

```
POINT FILE=DRIVERS /KEY=#FRSTAGE.
```

## **Overview**

POINT establishes the location at which sequential access begins (or resumes) in a keyed file. A keyed file is a file that provides access to information by a record key. An example of a keyed file is a file containing a social security number and other information about a firm's employees. The social security number can be used to identify the records in the file. For additional information about keyed files, see `KEYED DATA LIST`.

POINT prepares for reading the key-sequenced dataset sequentially from a point that the key value controls. Data selection commands can then be used to limit the file to the portion that you want to analyze. A `DATA LIST` command is used to read the data. To read keyed files (and also direct-access files), see `KEYED DATA LIST`.

## **Basic Specification**

The basic specification is the `KEY` subcommand and a string variable. The value of the string variable is used as the file key for determining where sequential retrieval (via `DATA LIST`) begins or resumes.

## **Subcommand Order**

- Subcommands can be named in any order.
- Each `POINT` command must precede its corresponding `DATA LIST` command.

## **Syntax Rules**

- `POINT` can be used more than once to change the order of retrieval during processing.
- `POINT` must be specified in an input program and therefore cannot be used to add cases to an existing file.

**Operations**

- The next `DATA LIST` command that is executed after the `POINT` command (for the same file) will read a record whose key value is at least as large as the value of the specified key. To prevent an infinite loop in which the same record is read again and again, the value of the variable that is specified on `KEY` must change from case to case, or the `POINT` command must be set up to execute only once.
- If the file contains a record whose key exactly matches the value of the `KEY` variable, the next execution of `DATA LIST` will read that record, the second execution of `DATA LIST` will read the next record, and so on.
- If an exact match between key and variable is not found, the results depend on the operating system. On IBM implementations, reading begins or resumes at the record that has the next higher key. If the value of the key is shorter than the file key, the value of the key variable is logically extended with the lowest character in the collating sequence. For example, if the value of the key variable is the single letter M, retrieval begins or resumes at the first record that has a key (regardless of length) beginning with the letter M or a character that is higher in the collating sequence.
- `POINT` does not report on whether the file contains a record that exactly matches the specified key. To check for missing records, use `LIST` to display the data that were read by the subsequent `DATA LIST` command.

**Examples****Basic Example**

```
FILE HANDLE DRIVERS/ file specifications.
POINT FILE=DRIVERS /KEY=#FRSTAGE.
```

- `FILE HANDLE` defines the handle for the data file to be read by `POINT`. The handle is specified on the `FILE` subcommand on `POINT`.
- `KEY` on `POINT` specifies the key variable. The key variable must be a string, and it must already exist as the result of a prior `DATA LIST`, `KEYED DATA LIST`, or transformation command.

**Selecting a Subset of Records from a Keyed File**

```
FILE HANDLE      DRIVERS/ file specifications.
INPUT PROGRAM.
STRING          #FRSTAGE(A2).
DO IF          #FRSTAGE = ' '.          /* First case check
+ COMPUTE      #FRSTAGE = '26'.        /* Initial key
+ POINT        FILE=DRIVERS /KEY=#FRSTAGE.
END IF.
DATA LIST       FILE=DRIVERS NOTABLE/
AGE 19-20(A) SEX 21(A) TICKETS 12-13.
DO IF          AGE > '30'.
+ END FILE.
END IF.
END INPUT PROGRAM.
LIST.
```

- This example illustrates how to execute POINT for only the first case. The file contains information about traffic violations, and it uses the individual's age as the key. Ages between 26 and 30 are selected.
- FILE HANDLE specifies the file handle DRIVERS.
- The INPUT PROGRAM and END INPUT PROGRAM commands begin and end the block of commands that build cases. POINT must appear in an input program.
- STRING declares the string variable #FRSTAGE, whose value will be used as the key on the POINT command. Because #FRSTAGE is a string variable, it is initialized as blanks.
- The first DO IF-END IF structure is executed only if no records have been read (that is, when #FRSTAGE is blank). When #FRSTAGE is blank, COMPUTE resets #FRSTAGE to 26, which is the initial value. POINT is executed, and it causes the first execution of DATA LIST to read a record whose key is at least 26. Because the value of #FRSTAGE is now 26, the DO IF-END IF structure is not executed again.
- DATA LIST reads the variables AGE, SEX, and TICKETS from the file DRIVERS.
- The second DO IF-END IF structure executes an END FILE command as soon as a record is read that contains a driver's age that is greater than 30. The program does not add this last case to the working file when it ends the file (see END FILE).

## FILE Subcommand

FILE specifies a file handle for the keyed data file. The file handle must have been previously defined on a FILE HANDLE command.

- FILE is optional.
- If FILE is omitted, POINT reads from the last file that is specified on an input command, such as DATA LIST.

### Example

```
FILE HANDLE DRIVERS/ file specifications.
POINT FILE=DRIVERS /KEY=#NXTCASE.
```

- FILE HANDLE specifies DRIVERS as the file handle for the data. The FILE subcommand on POINT specifies file handle DRIVERS.

## ENCODING Subcommand

ENCODING specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is UTF8. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).

- The quoted encoding value can be: `Locale` (the current locale setting), `UTF8`, `UTF16`, `UTF16BE` (big endian), `UTF16LE` (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- In Unicode mode, the defined width of string variables is tripled for code page and UTF-16 text data files. Use `ALTER TYPE` to automatically adjust the defined width of string variables.

## ***KEY Subcommand***

`KEY` specifies the variable whose value will be used as the file key for determining where sequential retrieval by `DATA LIST` will begin or resume. This variable must be a string variable, and it must already exist as the result of a prior `DATA LIST`, `KEYED DATA LIST`, or transformation command.

- `KEY` is required. Its only specification is a single variable. The variable can be a permanent variable or a scratch variable.
- Although the keys on a file are inherently numbers, such as social security numbers, the `STRING` function can be used to convert the numeric variable to a string. [For more information, see the topic String/numeric conversion functions on p. 111.](#)

### ***Example***

```
FILE HANDLE DRIVERS/ file specifications.  
POINT FILE=DRIVERS /KEY=#NXTCASE.
```

- `KEY` indicates that the value of the existing scratch variable `#FRSTAGE` will be used as the key to reading each record.
- Variable `#FRSTAGE` must be an existing string variable.

# P PLOT

P PLOT is available in the Statistics Base option.

```
P PLOT VARIABLES= varlist

[/DISTRIBUTION={NORMAL(a,b)** } ]
                {EXPONENTIAL(a)}
                {WEIBUL(a,b)}
                {PARETO(a,b)}
                {LNORMAL(a,b)}
                {BETA(a,b)}
                {GAMMA(a,b)}
                {LOGISTIC(a,b)}
                {LAPLACE(a,b)}
                {UNIFORM(a,b)}
                {HNORMAL(a)}
                {CHI(df)}
                {STUDENT(df)}

[/FRACTION={BLOM**}
           {RANKIT}
           {TUKEY}
           {VW}
           ]

[/TIES={MEAN**}
      {LOW}
      {HIGH}
      {BREAK}
      ]

[/{NOSTANDARDIZE**}
  {STANDARDIZE}
  ]

[/TYPE={Q-Q**}
      {P-P}
      ]

[/PLOT={BOTH**}
      {NORMAL}
      {DETRENDED}
      ]

[/DIFF={1}
      {n}
      ]

[/SDIFF={1}
      {n}
      ]

[/PERIOD=n]

[/{NOLOG**}
  {LN}
  ]

[/APPLY [= 'model name']]

[/TEMPLATE= 'filename']
```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 19.0

- `TEMPLATE` subcommand introduced.



### Example

```
P PLOT VARIABLES = VARX  
  /FRACTION=TUKEY  
  /DIFF=2.
```

## Overview

P PLOT (alias NP PLOT) produces probability plots of one or more sequence or time series variables. The variables can be standardized, differenced, and/or transformed before plotting. Expected normal values or deviations from expected normal values can be plotted.

### Options

**Variable Modification.** You can use the LN subcommand to request a natural log transformation of the sequence or time series variables, and you can use the SDIFF and DIFF subcommands to request seasonal and nonseasonal differencing to any degree. With seasonal differencing, you can specify the periodicity on the PERIOD subcommand. You can also plot standardized series by using the STANDARDIZE subcommand.

**Plot Type.** You can request p-p (proportion-proportion) or q-q (quantile-quantile) plots on the TYPE subcommand. With the PLOT subcommand, you can display normal plots, detrended plots, or both.

**Distribution Type.** You can specify the distribution type on the DISTRIBUTION subcommand. The cumulative distribution function (CDF) and the inverse distribution function (IDF) for the specified distribution type are used to compute the expected values in the p-p and q-q plots, respectively.

**Score Calculations.** On the FRACTION subcommand, you can specify one of several fractional rank formulas to use for estimating the empirical distribution in p-p plots and computing expected quantiles in q-q plots. You can specify the treatment of tied values on the TIE subcommand.

**Templates.** You can specify a template, using the TEMPLATE subcommand, to override the default chart attribute settings on your system.

### Basic Specification

The basic specification is one or more variable names.

- For each specified variable, P PLOT produces two q-q plots of the observed values (one plot versus expected normal values and the other plot versus deviations from normal values. By default, expected normal values are calculated by using Blom's transformation.
- Observed values define the horizontal axis, and expected normal values or deviations define the vertical axis.

### Subcommand Order

- Subcommands can be specified in any order.

### Syntax Rules

- VARIABLES can be specified only once.

- Other subcommands can be specified more than once, but only the last specification of each subcommand is executed.

### **Operations**

- Subcommand specifications apply to all plots that are produced by P PLOT.
- If the LN subcommand is specified, any differencing or standardization that is requested on that P PLOT is done on the log-transformed series.
- If differencing (DIFF or SDIFF) is specified, any standardization is done on the differenced series.

### **Limitations**

- A maximum of 1 VARIABLES subcommand is allowed. There is no limit on the number of variables that are named on the list.

## **Example**

```
P PLOT VARIABLES = VARX
      /FRACTION=TUKEY
      /DIFF=2 .
```

- This command produces two normal q-q plots of VARX (one plot not detrended and the other plot detrended).
- The expected quantile values are calculated by using Tukey's transformation.
- The variable is differenced twice before plotting.

## **VARIABLES Subcommand**

VARIABLES specifies the sequence or time series variables to be plotted and is the only required subcommand.

## **DISTRIBUTION Subcommand**

DISTRIBUTION specifies the distribution type of your data. The default is NORMAL if the subcommand is not specified or is specified without a keyword. If the parameters of the distribution type are not specified, DISTRIBUTION estimates them from the sample data and displays them with the plots.

<b>NORMAL(a,b)</b>	<i>Normal distribution.</i> The location parameter $a$ can be any numeric value, while the scale parameter $b$ must be positive. If they are not specified, DISTRIBUTION estimates them from the sample mean and sample standard deviation.
<b>EXPONENTIAL(a)</b>	<i>Exponential distribution.</i> The scale parameter $a$ must be positive. If the parameter is not specified, DISTRIBUTION estimates it from the sample mean. Negative observations are not allowed.
<b>WEIBULL(a,b)</b>	<i>Weibull distribution.</i> The scale and shape parameters $a$ and $b$ must be positive. If they are not specified, DISTRIBUTION estimates them using the least square method. Negative observations are not allowed.

<b>PARETO(a,b)</b>	<i>Pareto distribution.</i> The threshold and shape parameters $a$ and $b$ must be positive. If they are not specified, DISTRIBUTION assumes $a$ equals the minimum observation and estimates $b$ by the maximum likelihood method. Negative observations are not allowed.
<b>LNORMAL(a,b)</b>	<i>Lognormal distribution.</i> The scale and shape parameters $a$ and $b$ must be positive. If they are not specified, DISTRIBUTION estimates them from the mean and standard deviation of the natural logarithm of the sample data. Negative observations are not allowed.
<b>BETA(a,b)</b>	<i>Beta distribution.</i> The shape1 and shape2 parameters $a$ and $b$ must be positive. If they are not specified, DISTRIBUTION estimates them from the sample mean and sample standard deviation. All observations must be between 0 and 1, inclusive.
<b>GAMMA(a,b)</b>	<i>Gamma distribution.</i> The shape and scale parameters $a$ and $b$ must be positive. If they are not specified, DISTRIBUTION estimates them from the sample mean and sample standard deviation. Negative observations are not allowed.
<b>LOGISTIC(a,b)</b>	<i>Logistic distribution.</i> LOGISTIC takes a location and a scale parameter ( $a$ and $b$ ). The scale parameter ( $b$ ) must be positive. If the parameters are not specified, DISTRIBUTION estimates them from the sample mean and sample standard deviation.
<b>LAPLACE(a,b)</b>	<i>Laplace or double exponential distribution.</i> LAPLACE takes a location and a scale parameter ( $a$ and $b$ ). The scale parameter ( $b$ ) must be positive. If the parameters are not specified, DISTRIBUTION estimates them from the sample mean and sample standard deviation.
<b>UNIFORM(a,b)</b>	<i>Uniform distribution.</i> UNIFORM takes a minimum and a maximum parameter ( $a$ and $b$ ). Parameter $a$ must be equal to or greater than $b$ . If the parameters are not specified, DISTRIBUTION assumes them from the sample data.
<b>HNORMAL(a)</b>	<i>Half-normal distribution.</i> Data are assumed to be location-free or centralized. (Location parameter=0.) You can specify the scale parameter $a$ or let DISTRIBUTION estimate it by using the maximum likelihood method.
<b>CHI(df)</b>	<i>Chi-square distribution.</i> You must specify the degrees of freedom ( $df$ ). Negative observations are not allowed.
<b>STUDENT(df)</b>	<i>Student's t distribution.</i> You must specify the degrees of freedom ( $df$ ).

## **FRACTION Subcommand**

FRACTION specifies the formula to be used in estimating the empirical distribution in p-p plots and calculating the expected quantile values in q-q plots.

- Only one formula can be specified. If more than one formula is specified, only the first formula is used.
- If the FRACTION subcommand is not specified, BLOM is used by default.
- These formulas produce noticeable differences for short series only.

Four formulas are available:

<b>BLOM</b>	Blom's transformation, defined by the formula $(r - (3/8)) / (n + (1/4))$ , where $n$ is the number of observations and $r$ is the rank, ranging from 1 to $n$ (Blom, 1958).
<b>RANKIT</b>	Formula $(r - (1/2)) / n$ , where $n$ is the number of observations and $r$ is the rank, ranging from 1 to $n$ (Chambers, Cleveland, Kleiner, and Tukey, 1983).

- TUKEY** Tukey's transformation, defined by the formula  $(r - (1/3)) / (n + (1/3))$ , where  $n$  is the number of observations and  $r$  is the rank, ranging from 1 to  $n$  (Tukey, 1962).
- VW** Van der Waerden's transformation, defined by the formula  $r / (n + 1)$ , where  $n$  is the number of observations and  $r$  is the rank, ranging from 1 to  $n$  (Lehmann, 1975).

### Example

```
P PLOT VARIABLES = VARX
      / FRACTION=VW.
```

- This P PLOT command uses van der Waerden's transformation to approximate the proportion estimate  $p$ , which is used in the inverse distribution function.
- By default, two q-q plots are produced.

## TIES Subcommand

TIES determines the way that tied values are handled. The default method is MEAN.

- MEAN** Mean rank of tied values is used for ties. This setting is the default.
- LOW** Lowest rank of tied values is used for ties.
- HIGH** Highest rank of tied values is used for ties.
- BREAK** Consecutive ranks with ties sharing the same value. Each distinct value of the ranked variable is assigned a consecutive rank. Ties share the same rank.

## TYPE Subcommand

TYPE specifies the type of plot to produce. The default is Q-Q. The plots show a quantile-quantile plot and a proportion-proportion plot using the same data (with a normal distribution).

- Q-Q** Quantile-quantile plots. The quantiles of the observed values are plotted against the quantiles of the specified distribution.
- P-P** Proportion-proportion plots. The observed cumulative proportion is plotted against the expected cumulative proportion if the data were a sample from a specified distribution.

Figure 182-1  
Normal q-q plot of household income

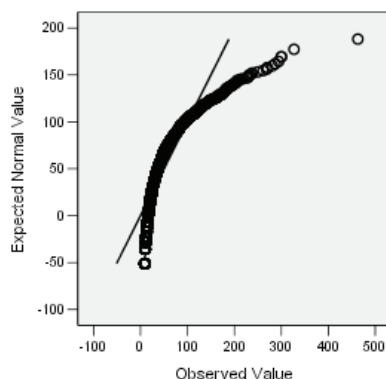
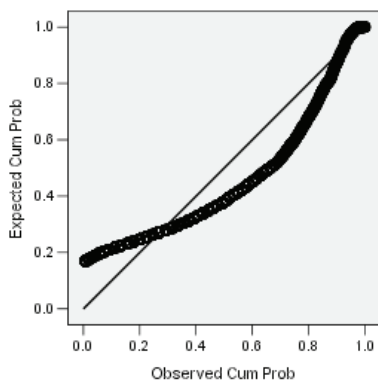


Figure 182-2  
Normal p-p plot of household income



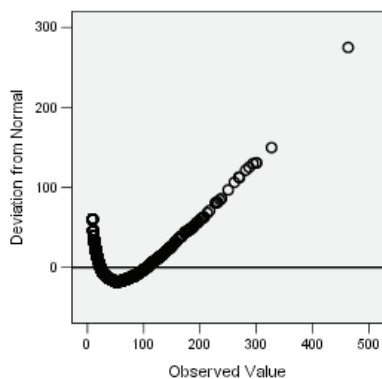
## PLOT Subcommand

PLOT specifies whether to produce a plot of observed values versus expected values, a plot of observed values versus deviations from expected values, or both. The plots shown in [TYPE Subcommand](#) are nondetrended plots. The figure below shows a detrended q-q plot.

<b>BOTH</b>	<i>Display both detrended and nondetrended normal plots.</i> This is the default.
<b>NORMAL</b>	<i>Display nondetrended normal plots.</i> The observed values are plotted against the expected values.
<b>DETRENDED</b>	<i>Display detrended plots.</i> The observed values are plotted against the deviations from the expected values.

- If you specify PLOT more than once, only the last specification is executed.
- Deviations are calculated by subtracting the expected value from the observed value.
- In low resolution, a dash is used in a detrended plot to indicate where the deviation from the expected is 0.

Figure 182-3  
Detrended normal q-q plot of household income



## ***STANDARDIZE and NOSTANDARDIZE Subcommands***

STANDARDIZE transforms the sequence or time series variables into a sample with a mean of 0 and a standard deviation of 1. NOSTANDARDIZE is the default and indicates that the series should not be standardized.

- There are no additional specifications on the STANDARDIZE or NOSTANDARDIZE subcommands.
- Only the last STANDARDIZE or NOSTANDARDIZE subcommand on the P PLOT command is executed.
- The STANDARDIZE and NOSTANDARDIZE subcommands have no effect on expected values, which are always standardized.
- NOSTANDARDIZE is generally used with an APPLY subcommand to turn off a previous STANDARDIZE specification.

### ***Example***

```
P PLOT VARIABLES = VARX  
/STANDARDIZE.
```

- This example produces two q-q normal-probability plots of *VARX* with standardized observed values.

## ***DIFF Subcommand***

DIFF specifies the degree of differencing that is used before plotting to convert a nonstationary variable into a stationary variable with a constant mean and variance.

- You can specify any positive integer on DIFF.
- If DIFF is specified without a value, the default is 1.
- The number of plotted values decreases by 1 for each degree of differencing.

### ***Example***

```
P PLOT VARIABLES = TICKETS  
/DIFF=2.
```

- In this example, *TICKETS* is differenced twice before the expected and observed values are plotted.

## ***SDIFF Subcommand***

If the variable exhibits a seasonal or periodic pattern, you can use the SDIFF subcommand to seasonally difference the variable before plotting.

- The specification on SDIFF indicates the degree of seasonal differencing and can be any positive integer.
- If SDIFF is specified without a value, the degree of seasonal differencing defaults to 1.

- The number of plotted seasons decreases by 1 for each degree of seasonal differencing.
- The length of the period that is used by `SDIFF` is specified on the `PERIOD` subcommand. If the `PERIOD` subcommand is not specified, the periodicity that is established on the `TSET` or `DATE` command is used (see [PERIOD Subcommand](#)).

## PERIOD Subcommand

`PERIOD` indicates the length of the period to be used by the `SDIFF` subcommand.

- The specification on `PERIOD` indicates how many observations are in one period or season. You can specify any positive integer on `PERIOD`.
- The `PERIOD` subcommand is ignored if it is used without the `SDIFF` subcommand.
- If `PERIOD` is not specified, the periodicity that is established on `TSET` `PERIOD` is in effect. If `TSET` `PERIOD` is not specified either, the periodicity that is established on the `DATE` command is used. If periodicity was not established anywhere, the `SDIFF` subcommand will not be executed.

### Example

```
PLOT VARIABLES = TICKETS
  /SDIFF=1
  /PERIOD=12 .
```

- This command applies 1 degree of seasonal differencing with 12 observations per season to the variable `TICKETS`.

## LN and NOLOG Subcommands

`LN` transforms the data by using the natural logarithm (base  $e$ ) to remove varying amplitude. `NOLOG` indicates that the data should not be log transformed. `NOLOG` is the default.

- There are no additional specifications on `LN` or `NOLOG`.
- Only the last `LN` or `NOLOG` subcommand on a `PLOT` command is executed.
- If a natural log transformation is requested, cases with values that are less than or equal to 0 will be set to system-missing, because nonpositive values cannot be log-transformed.
- `NOLOG` is generally used with an `APPLY` subcommand to turn off a previous `LN` specification.

### Example

```
PLOT VARIABLES = TICKETS
  /FRACTION=TUKEY
  /DIFF=1
  /LN.
PLOT VARIABLES = EARNINGS
  /APPLY
  /NOLOG.
```

- The first command requests a natural log transformation of variable `TICKETS` before plotting.
- The second command applies the previous `PLOT` specifications to variable `EARNINGS`. However, `EARNINGS` is not log-transformed before plotting.

## **APPLY Subcommand**

APPLY allows you to produce a plot by using previously defined specifications without having to repeat the P PLOT subcommands.

- The only specification on APPLY is the name of a previous model in quotation marks. If a model name is not specified, the model that is specified on the previous P PLOT command is used.
- To change any plot specifications, specify the subcommands of only those portions that you want to change. Make these entries after the APPLY subcommand.
- If no variables are specified, the variables that were specified for the original plot are used.
- To change the variables that are used with the model, enter new variable names before or after the APPLY subcommand.
- The distribution type is applied, but the parameters are not applied.

### **Example**

```
P PLOT VARIABLES = X1
  /FRACTION=TUKEY.
P PLOT VARIABLES = Z1
  /APPLY.
```

- The first command produces two q-q normal-probability plots of *X1*, using Tukey's transformation to compute the expected values.
- The second command requests the same plots for variable *Z1*.

### **Example**

```
P PLOT VARIABLES = X1 Y1 Z1
  /FRACTION=VW.
P PLOT APPLY
  /FRACTION=BLOM.
```

- The first command uses van der Waerden's transformation to calculate expected normal values of *X1*, *Y1*, and *Z1*.
- The second command uses Blom's transformation for the same three series.

### **Example**

```
P PLOT VARIABLES = VARX
  /FRACTION=RANKIT
  /DIFF
  /STANDARDIZE.
P PLOT VARIABLES = VARY
  /APPLY
  /NOSTANDARDIZE.
```

- The first command differences and standardizes series *VARX* and then produces a normal probability plot by using the RANKIT transformation.
- The second command applies the previous plot specifications to *VARY* but does not standardize the series.



## **TEMPLATE Subcommand**

TEMPLATE uses an existing file as a template and applies it to any charts generated by the current PLOT command. The template overrides the default settings that are used to create any charts, and the specifications on the current PLOT command override the template. Templates are created in the Chart Editor by saving an existing chart as a template.

### **Example**

```
PLOT VARIABLES = varx
  /FRACTION=TUKEY
  /DIFF=2
  /TEMPLATE= '/templates/mytemplate.sgt'.
```

## **References**

- Blom, G. 1958. *Statistical estimates and transformed beta variables*. New York: John Wiley and Sons.
- Chambers, J. M., W. S. Cleveland, B. Kleiner, and P. A. Tukey. 1983. *Graphical methods for data analysis*. Boston: Duxbury Press.
- Lehmann, E. L. 1975. *Nonparametrics: Statistical methods based on ranks*. San Francisco: Holden-Day.
- Tukey, J. W. 1962. The future of data analysis. *Annals of Mathematical Statistics*, 33:22, 1–67.

# PREDICT

```
PREDICT [{start date      }] [THRU [{end date      }]]  
        {start case number}      {end case number}  
                                   {END          }
```

## Example

```
PREDICT Y 61 THRU Y 65.
```

## Overview

PREDICT specifies the observations that mark the beginning and end of the forecast period. If the forecast period extends beyond the length of the series, PREDICT extends the series in the active dataset to allow room for the forecast observations.

### Basic Specification

The minimum specification on PREDICT is either the start or the end of the range, or it is keyword THRU. PREDICT sets up a forecast period beginning and ending with the specified dates or case numbers. The default starting point is the observation immediately after the end of the series or, if USE is specified, the observation immediately after the end of the use range (the historical period). The default end is the last observation in the series.

### Operations

- PREDICT is executed when the data are read for the next forecasting procedure (for example, CURVEFIT in the Core system, and 2SLS in the Regression add-on module).
- PREDICT is ignored by non-forecasting procedures.
- Case number specifications refer to the sequential numbers that are assigned to cases as they are read.
- If the forecast period extends beyond the length of the series, PREDICT extends the series in the active dataset to allow room for the forecast observations.
- New observations that are added to the end of existing series will contain non-missing date variables, forecast values (variable *FIT#n*), confidence interval limits (variables *LCL#n* and *UCL#n*), and, for ARIMA models, standard error of the predicted value (*SEP#n*). For all other variables, including the original series, the new cases will be system-missing.
- PREDICT cannot forecast beyond the end of the series for ARIMA with regressors and 2SLS. However, PREDICT can forecast values for the dependent variable if the independent variables have valid values in the predict period.
- If the use and predict periods overlap, the model is still estimated by using all observations in the use period.

- `USE` and `PREDICT` can be used together to perform forecasting validation. To do this, specify a use period that ends before the existing end of the series, and specify a predict period starting with the next observation.
- If there is a gap between the end of the use period and the start of the specified predict period, the program uses the first observation after the end of the use period as the start of the predict period. (This setting is the default.)
- The `DATE` command turns off all existing `USE` and `PREDICT` specifications.
- `PREDICT` remains in effect in a session until it is changed by another `PREDICT` command or until a new `DATE` command is issued.
- If more than one forecasting procedure is specified after `PREDICT`, the `USE` command should be specified between procedures so that the original series—without any new, system-missing cases—will be used each time. Alternatively, you can specify
 

```
TSET NEWVAR = NONE
```

 before the first forecasting procedure so that you can evaluate model statistics without creating new variables or adding new cases with missing values to the original series.

### **Limitations**

A maximum of one range (one start and/or one end) can be specified per `PREDICT` command.

## **Syntax Rules**

- You can specify a start, an end, or both.
- The start and end are specified as either date specifications or case (observation) numbers.
- Date specifications and case numbers cannot be mixed on one `PREDICT` command.
- Keyword `THRU` is required if the end of the range is specified.
- Keyword `THRU` by itself defines a `PREDICT` range starting with the first observation after the use range and ending with the end of the series. If `USE` has not been specified, `PREDICT THRU` is meaningless.

## **Date Specifications**

- A date specification consists of `DATE` keywords and values (see the `DATE` command on p. 586). These specifications must correspond to existing date variables.
- If more than one date variable exists, the highest-order variable must be included in the date specification.
- Values on keyword `YEAR` must have the same format (two or four digits) as the `YEAR` specifications on the `DATE` command.

## **Case Specifications**

The case number specification is the sequence number of the case (observation) as it is read by the program.

**Valid Range**

- The start date must precede the end date.
- The start case number must be less than the end case number.
- The start can be any observation ranging from the second observation in the historical period that is specified on `USE` to the observation immediately following the end of the historical period. If `USE` is not specified, the start can be any observation ranging from the second observation in the series to the observation immediately following the end of the series.
- For most models, the start of the predict period should not be too close to the start of the use period.
- The predict and use periods should not be exactly the same.
- The start of the predict period should not precede the start of the use period.

**Examples*****Specifying the Forecast Period as a Date Range***

```
PREDICT Y 61 THRU Y 65.
```

- This command specifies a forecast period from 1961 to 1965.
- The active dataset must include variable `YEAR_`, which, in this example, contains only the last two digits of each year.
- If variable `MONTH_` also exists, the above command is equivalent to

```
PREDICT Y 61 M 1 THRU Y 65 M 12.
```

```
PREDICT W 28 THRU W 56.
```

- This command specifies a forecast period from the 28th week to the 56th week.
- The active dataset must include variable `WEEK_`.
- If variable `DAY_` also exists, the above command is equivalent to

```
PREDICT W 28 D 1 THRU W 56 D 7.
```

***Specifying the Forecast Period as a Case Range***

```
PREDICT 61 THRU 65.
```

- This command specifies a forecast period from the 61st case (observation) to the 65th case.

***Using the Default Start Date***

```
PREDICT THRU Y 65.
```

- This command uses the default start date, which is the observation immediately following the end of the use period. If `USE` is not specified, the default start is the observation immediately following the end of the series.
- The forecast period extends from the start date through year 1965.

- The active dataset must include variable *YEAR\_*.
- Keyword `THRU` is required.

***Specifying the Forecast Period by Using Date Variables***

`PREDICT THRU CYCLE 4 OBS 17.`

- This example uses the date variables *OBS\_* and *CYCLE\_*, which must exist in the active dataset.
- *CYCLE*, the highest order, must be included on `PREDICT`.
- Keyword `THRU` is required.
- The forecast period extends from the default start to the 17th observation of cycle 4.

# PREFSCAL

PREFSCAL is available in the Categories option.

```
PREFSCAL VARIABLES = varlist

[/INPUT = [ROWS({n      })]
           {rowid}
           [SOURCES({1**      })]]
           {n      }
           {sourceid}

[/PROXIMITIES = {DISSIMILARITIES**}]
                {SIMILARITIES      }

[/WEIGHTS = varlist]

[/INITIAL = {CLASSICAL[({TRIANGLE**})]}]
            {SPEARMAN      }
            {CORRESPONDENCE      }
            {ROSSCLIFF      }
            {CENTROIDS[({1})]}      }
            {n      }
            {RANDOM[({1})]}      }
            {n      }
            {(filespec) [varlist]      }

[/CONDITION = {ROW**}]
              {MATRIX}
              {UNCONDITIONAL      }

[/TRANSFORMATION = {NONE[ (INTERCEPT) ]}
                   {LINEAR[ (INTERCEPT) ]}
                   {ORDINAL[ ({UNTIE      })]}
                   {KEEP TIES**}
                   {SMOOTH[ ({UNTIE      })]}
                   {KEEP TIES}
                   {SPLINE[ ({INTERCEPT} [ORDER={2}] [INKNOT={1}])]}
                   {n      }
                   {n      }

[/MODEL = {IDENTITY** }
          {WEIGHTED      }
          {GENERALIZED}

[/RESTRICTIONS = {ROW      ({NONE**      }) (filespec) [varlist]}}]
                 {COORDINATES}
                 {COLUMN({NONE**      }) (filespec) [varlist]}}]
                 {COORDINATES}

[/PENALTY = [LAMBDA({0.5**})]
            {value}
            [OMEGA({1.0**})]
            {value}

[/CRITERIA = [DIMENSIONS({2**      })]
             {min[,max]}]
             [MAXITER({5000**})]
             {value }
             [DIFFSTRESS({0.000001**})]
             {value }
             [MINSTRESS({0.0001**})]
             {value }

[/PRINT = [NONE] [INPUT] [MULTIPLE] [INITIAL]
          [HISTORY] [MEASURES**] [DECOMPOSITION]
          [COMMON**] [DISTANCES] [WEIGHTS**]
          [INDIVIDUAL] [TRANSFORMATION]]

[/PLOT = [NONE] [MULTIPLE] [INITIAL] [STRESS]
```

```

[COMMON**] [WEIGHTS**]
[INDIVIDUAL[(valuelist) [...]]]
[TRANSFORMATIONS[(valuelist) [...]]]
[SHEPARD[(valuelist) [...]]]
[FIT[(valuelist) [...]]]
[RESIDUALS[(valuelist) [...]]]

[/OPTIONS = [MARKERS(rowid)]
             [COLORS(rowid)]]

[/OUTFILE = [COMMON('savfile'|'dataset')]
            [WEIGHTS('savfile'|'dataset')]
            [DISTANCES('savfile'|'dataset')]
            [TRANSFORMATIONS('savfile'|'dataset')]]

```

\* Default if the keyword is omitted.

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 14.0

- Command introduced.

### **Example**

```
PREFSCAL var01 TO var02.
```

## **Overview**

PREFSCAL performs multidimensional unfolding of proximity data to find a least-squares representation of the row and column objects in a low-dimensional space. Individual differences models are allowed for multiple sources. A majorization algorithm minimizes penalized Stress and guarantees monotone convergence for optionally transformed, metric and nonmetric data under a variety of models and constraints.

### **Options**

**Data Input.** You can read one or more rectangular matrices of proximities. Additionally, you can read weights, an initial configuration, and fixed coordinates.

**Methodological Assumptions.** On the `CONDITION` subcommand, you can specify transformations for all sources (unconditional), separate transformations for each source (matrix-conditional), or separate transformations for each row (row-conditional). Using the `TRANSFORMATION` subcommand, you can treat proximities as nonmetric (ordinal or smooth ordinal), as quasi-metric (splines), or as metric (linear with or without intercept). Ordinal and smooth ordinal transformations can keep tied observations tied (discrete) or untie them (continuous). You can use the `PROXIMITIES` subcommand to specify whether your proximities are similarities or dissimilarities.

**Model Selection.** You can specify multidimensional unfolding models by selecting a combination of PREFSCAL subcommands, keywords, and criteria. The subcommand MODEL offers the Identity model and two individual differences models. You can specify other selections on the CRITERIA subcommand.

**Penalties.** You can specify two penalty parameters in order to avoid degenerate solutions. On the PENALTY subcommand, LAMBDA is specified for the strength, and OMEGA is specified for the range. Penalized Stress penalizes solutions with insufficient variation in the transformed proximities.

**Constraints.** You can specify fixed coordinates on the RESTRICTIONS subcommand to restrict some or all common space coordinates of either row objects or column objects.

**Output.** You can produce output that includes the original and transformed proximities, history of iterations, common and individual configurations, individual space weights, distances, and decomposition of the Stress. Plots can be produced of common and individual configurations (biplots), individual space weights, transformations, fit, and residuals.

### **Basic Specification**

The basic specification is PREFSCAL followed by a variable list. By default, PREFSCAL produces a two-dimensional metric Euclidean multidimensional unfolding solution (Identity model). Input is expected to contain one or more rectangular matrices with proximities that are dissimilarities. The transformation of the proximities is row-conditional. The analysis uses a classical scaling start as initial configuration. By default, output includes fit and Stress values, the coordinates of the common space, and a joint plot of the common space configuration.

### **Syntax Rules**

- If there is only one source, the model is always assumed to be Identity.
- In the case of duplicate or contradicting subcommand specification, only the later subcommand applies.
- There is no constraint with respect to the syntax order.

### **Limitations**

- PREFSCAL needs at least two objects in each set. At least two variables need to be specified in the variable list (two column objects), and the active data file needs to contain at least two cases (two row objects).
- PREFSCAL does not honor SPLIT FILE.

## **Examples**

```
PREFSCAL
VARIABLES=TP BT EMM JD CT BMM HRB TmD BTJ TMn CB DP GD CC CMB
/INPUT=SOURCES(srcid )
/INITIAL=CLASSICAL (SPEARMAN)
/CONDITION=ROW
/TRANSFORMATION=NONE
/PROXIMITIES=DISSIMILARITIES
/MODEL=WEIGHTED
/CRITERIA=DIMENSIONS(2,2) DIFFSTRESS(.000001) MINSTRESS(.0001)
MAXITER(5000)
```



```

/PENALTY=LAMBDA(0.5) OMEGA(1.0)
/PRINT=MEASURES COMMON
/PLOT=COMMON WEIGHTS INDIVIDUAL ( ALL ) .

```

- This syntax specifies an analysis on variables *tp* (*Toast pop-up*) through *cmb* (*Corn muffin and butter*). The variable *srcid* is used to identify the sources.
- The `INITIAL` subcommand specifies that the starting values be imputed using Spearman distances.
- The `MODEL` subcommand specifies a weighted Euclidean model, which allows each individual space to weight the dimensions of the common space differently.
- The `PLOT` subcommand requests plots of the common space, individual spaces, and individual space weights.
- All other parameters fall back to their default values.

## VARIABLES Subcommand

The variable list identifies the columns in the proximity matrix or matrices that PREFSCAL reads. Each variable identifies one column of the proximity matrix, with each case in the active dataset representing one row.

- Only numeric variables may be specified.
- PREFSCAL reads data row by row, and the columns are represented by the variables on the variable list.

### Example

```

DATA LIST
  /var01 var02.
BEGIN DATA
  1 6
  5 4
  4 2
END DATA.
PREFSCAL var01 TO var02.

```

- This example specifies an analysis on a  $3 \times 2$  proximity matrix (3 rows and 2 columns).

## INPUT Subcommand

The `INPUT` subcommand specifies the number of rows in one source, the number of sources, or both. Specifying a row identifier, *rowid*, or a source identifier, *sourceid*, specifically identifies either the row objects or sources and provides a variable that may contain row object or source labels. Specifying only one keyword computes the number of row objects or sources according to the following formula:  $C=R \times S$ , where  $C$  is the number of cases,  $R$  is the number of row objects,

and S is the number of sources. By specifying both keywords, PREFSCAL takes the first  $R \times S$  cases from the active file.

- ROWS**                    *Number of rows.* This specifies the number of row objects in one source. A variable in parentheses specifies a row identifier. The values must be positive nonzero integers. The values of this variable specify the identifier of the rows that contain the proximities. Within one source, the values of this identifier only need to be discriminating. Over sources, the values of the row identifier must be in the same order.
- SOURCES**                *Number of sources.* This keyword specifies the number of sources. By default, the number of sources is 1. Otherwise, the number of cases in the active file must be dividable by the number of sources. A variable in parentheses specifically specifies a source identifier. The values of this identifier specify the sources and must be positive nonzero integers. The rows within one source must be consecutive cases in the active data file.

- *rowid* and *sourceid* may not be specified on the PREFSCAL variable list.
- Duplicate cell specification is not taken care of. In this case, the final specification applies.

### Example

```
DATA LIST
  /var01 TO var15 rowid sourceid.
BEGIN DATA
  13 12 07 03 05 04 08 11 10 15 02 01 06 09 14   1  1
  15 11 06 03 10 05 14 08 09 12 07 01 04 02 13   2  1
  15 10 12 14 03 02 09 08 07 11 01 06 04 05 13   3  1
  (...)
  10 03 02 14 09 01 08 12 13 04 11 05 15 06 07   41  1
  13 03 01 14 04 10 05 15 06 02 11 07 12 08 09   42  1
  15 03 05 12 02 08 07 13 01 04 06 10 11 09 14   1  2
  15 04 03 11 07 05 14 01 02 06 08 13 09 12 10   2  2
  15 02 07 12 05 06 04 08 01 03 09 10 11 14 13   3  2
  (...)
  (...)
  09 04 07 10 11 02 08 12 13 05 14 06 15 01 03   41  6
  15 10 01 12 02 06 08 14 13 11 09 03 04 05 07   42  6
END DATA.
```

- The active data file has 252 cases, containing 6 sources with 42 row objects per source and containing 15 column objects. Additionally, 2 identifying variables, *rowid* and *sourceid*, are specified to identify the row objects and sources, respectively.

```
PREFSCAL var01 TO var15
  /INPUT = ROWS(42).
```

- PREFSCAL reads 15 columns and 42 rows per source—thus, 6 sources in total (252/42).

```
PREFSCAL var01 TO var15
  /INPUT = SOURCES(6).
```

- PREFSCAL reads 15 columns and 6 sources, with 42 rows each (252/6).

```
PREFSCAL var01 TO var15
  /INPUT = ROWS(rowid).
```

- PREFSCAL reads 15 columns and 42 rows per source. The row objects are specified by *rowid*, which ranges from 1 to 42, the number of row objects in this case (per source, thus 6 sources). When a lower value is found in the row object identifier variable, a new source is started.

```
PREFSCAL var01 TO var15
  /INPUT = SOURCES(sourceid).
```

- PREFSCAL reads 15 columns and 6 sources. The sources are specified by *sourceid*, which ranges from 1 to the number of sources (in this case, from 1 to 6). When a higher value is found in the source identifier variable, a new source is started.

```
COMPUTE rowid = 1+MOD($casenum-1,42).
COMPUTE sourceid = 1+TRUNC(($casenum-1)/42).
SORT CASES BY sourceid (A) rowid (A).
VALUE LABELS sourceid 1 'overall' 2 'bacon' 3 'cereal'
                4 'pancakes' 5 'normal' 6 'snack'.
PREFSCAL var01 TO var15
  /INPUT = ROWS(rowid) SOURCES(sourceid).
```

- First, a row object identifier and a source identifier are computed. The active data file is sorted by *rowid* and *sourceid*. The variable *sourceid* is given value labels. PREFSCAL reads 15 columns and 6 sources per row object, with 42 row objects in total. The first nine case of the active data file look like this:

```
13 12 7 3 5 4 8 11 10 15 2 1 6 9 14 1 1
15 3 5 12 2 8 7 13 1 4 6 10 11 9 14 1 2
14 10 9 3 7 6 12 15 8 11 5 1 4 2 13 1 3
14 9 7 1 11 6 12 15 13 10 3 4 2 5 8 1 4
10 12 8 4 9 6 7 15 14 13 3 1 5 2 11 1 5
14 11 8 4 9 6 7 15 10 12 3 2 5 1 13 1 6
15 11 6 3 10 5 14 8 9 12 7 1 4 2 13 2 1
15 4 3 11 7 5 14 1 2 6 8 13 9 12 10 2 2
15 8 2 3 10 1 14 5 6 9 11 12 4 13 7 2 3
(...)
```

## ***PROXIMITIES Subcommand***

The PROXIMITIES subcommand specifies the type of proximities that are used in the analysis. The term **proximity** is used for either similarity or dissimilarity data. Internally, PREFSCAL works with dissimilarities. Therefore, PREFSCAL converts similarities into dissimilarities by reflecting the data about its midpoint (depending on the conditionality chosen on the CONDITION subcommand), thus preserving the endpoints and the range.

**DISSIMILARITIES** *Dissimilarity data.* This specification is the default when PROXIMITIES is not specified. Small dissimilarities correspond to small distances, and large dissimilarities correspond to large distances.

**SIMILARITIES** *Similarity data.* Small similarities correspond to large distances, and large similarities correspond to small distances.

**Example**

```
PREFSCAL var01 TO var09
  /PROXIMITIES = SIMILARITIES.
```

- In this example, PREFSCAL expects the proximities to be similarities.

**WEIGHTS Subcommand**

The WEIGHTS subcommand specifies the variables that contain the nonnegative weights for the proximities that are included in the active dataset.

- The number and order of the variables in the variable list is important. The first variable in the WEIGHTS variable list corresponds to the first variable on the PREFSCAL variable list. This correspondence is repeated for all variables on the variable lists. Every proximity has its own weight. Therefore, the number of variables on the WEIGHTS subcommand must be equal to the number of variables on the PREFSCAL variable list.
- Negative weights are not allowed. If negative weights are specified, an error message is issued, and the procedure aborts.
- The weight variable (set with WEIGHT BY) allows for the weighting of entire rows. The weight variable must contain positive, nonzero values.

**Example**

```
DATA LIST FILE = 'breakfast.dat' FREE
  /var01 TO var15 wgt01 TO wgt15.
PREFSCAL var01 TO var15
  /WEIGHTS = wgt01 TO wgt15.
```

- In this example, the PREFSCAL variable list indicate that there are 15 column objects, of which the weights can be found in *wgt01* to *wgt15*.
- *wgt01* contains the weights for *var01*, *wgt02* contains the weights for *var02*, and so on.

**INITIAL Subcommand**

INITIAL defines the initial or starting configuration of the common space for the analysis. When a reduction in dimensionality is specified on the CRITERIA subcommand, a derivation of coordinates in the higher dimensionality is used as starting configuration in the lower dimensionality.

- You can specify one of the five keywords that are listed below.
- You can specify a variable list containing the initial configuration.

**CLASSICAL**

*Classical scaling start.* This specification is the default. The rectangular proximity matrix is used to supplement the intra-blocks (values between rows and between columns) of the complete symmetrical MDS matrix by means of the triangular inequality or Spearman distances. When the complete matrix is formed, a classical scaling solution is used as initial configuration.

**ROSSCLIFF**

*Ross-Cliff start.* The Ross-Cliff start uses the results of a singular value decomposition on the double centered and squared proximity matrix as the initial values for the row and column objects.

<b>CORRESPONDENCE</b>	<i>Correspondence start.</i> The correspondence start uses the results of a correspondence analysis on the reversed data (similarities instead of dissimilarities) with symmetric normalization of row and column scores. For more information, see the topic <b>CORRESPONDENCE</b> on p. 355.
<b>CENTROIDS(n)</b>	<i>Centroids start.</i> PREFSCAL starts by positioning the row objects in the configuration by using an eigenvalue decomposition. Then, the column objects are positioned at the centroid of first choices (or second if $n=2$ or third if $n=3$ , etc.). The number of choices ( $n$ ) must be a positive integer between 1 and the number of columns. The default is 1.
<b>RANDOM(n)</b>	<i>(Multiple) random start.</i> You can specify the number of random starts ( $n$ ), where $n$ is any positive integer. The random sequence can be controlled by the <b>SET SEED</b> procedure (thus, not by a subcommand within the <b>PREFSCAL</b> procedure). All $n$ analyses start with a different random configuration. In the output, all $n$ final Stress values are reported, as well as the initial seeds of each analysis (for reproduction purposes), followed by the full output of the analysis with the lowest penalized Stress value. The default number of random starts is 1.

### **CLASSICAL Keyword**

<b>TRIANGLE</b>	<i>Imputation using the triangle inequality.</i> If <b>TRIANGLE</b> is specified, the intra-blocks are filled by using the triangular inequality.
<b>SPEARMAN</b>	<i>Imputation with Spearman distances.</i> If <b>SPEARMAN</b> is specified, the Spearman distances between all objects are used to create a symmetrical MDS matrix.

Instead of these keywords, a filespec in parentheses can be given to specify the IBM® SPSS® Statistics data file containing the coordinates of the initial configuration. The row and column coordinates are stacked, with the column coordinates following the row coordinates. The closing parenthesis of the filespec can be followed by a variable list. If the variable list is omitted, the procedure automatically selects the first *MAXDIM* variables in the external file, where *MAXDIM* is the maximum number of dimensions that are requested for the analysis on **/CRITERIA = DIMENSIONS(min, max)**. Missing values are not allowed as initial coordinates. An error is issued whenever this situation occurs.

### **Example**

```
PREFSCAL var01 TO var15
  /INITIAL = RANDOM(100).
```

- This example performs 100 analyses (each analysis starting with a different random configuration). The results of the analysis with the lowest final Stress are displayed in the output.

## **CONDITION Subcommand**

CONDITION specifies the range of proximities that are compared within one transformation list. The TRANSFORMATION subcommand specifies the type of transformation.

<b>ROW</b>	<i>Row conditional.</i> Only the proximities within each row are compared with each other. The comparison is carried out for each row separately. This setting is the default.
<b>MATRIX</b>	<i>Matrix conditional.</i> Only the proximities within each source are compared with each other. The comparison is carried out for each source separately.
<b>UNCONDITIONAL</b>	<i>Unconditional.</i> This specification is appropriate when the proximities in all sources can be compared with each other, and it results in a single transformation of all sources simultaneously.

- Note that if there is only one source, MATRIX and UNCONDITIONAL yield the same result.

### **Example**

```
PREFSCAL var01 TO var09
  /CONDITION = UNCONDITIONAL
  /TRANSFORMATION = LINEAR (INTERCEPT) .
```

- In this example, the proximities are linearly transformed, including an intercept. The transformation is carried out over all proximities simultaneously.

## **TRANSFORMATION Subcommand**

The TRANSFORMATION subcommand offers four different options for optimal transformation of the original proximities. The resulting values are called transformed proximities. The distances between the objects in the configuration should match these transformed proximities as closely as possible. The CONDITION subcommand specifies over which proximities the transformation is computed. The default transformation is ORDINAL with ties kept tied.

<b>NONE</b>	<i>No scale transformation.</i> The INTERCEPT keyword can be specified in parentheses following the NONE keyword. If INTERCEPT is specified, an intercept is estimated in the transformation.
<b>LINEAR</b>	<i>Linear transformation.</i> With this transformation, the transformed proximities are proportional to the original proximities (that is, the transformation function estimates a slope, and the intercept is fixed at 0). The INTERCEPT keyword can be specified in parentheses following the LINEAR keyword. If INTERCEPT is specified, an intercept is estimated in the transformation, resulting in an interval transformation. Without the keyword INTERCEPT, LINEAR only estimates a slope, which coincides with a ratio transformation.
<b>ORDINAL</b>	<i>Ordinal transformation.</i> The transformed proximities have the same order as the original proximities. ORDINAL can be followed by a keyword in parentheses to indicate how to handle tied proximities.

<b>SMOOTH</b>	<i>Smooth ordinal transformation.</i> The transformed proximities have the same order as the original proximities, including a smoothness restriction. This restriction takes the differences between subsequent values into account. Restricting subsequent differences allows for a smooth ordinal transformation. <i>SMOOTH</i> can be followed by a keyword in parentheses to indicate how to handle tied proximities.
<b>SPLINE</b>	<i>Monotone spline transformation.</i> The transformed proximities are a smooth nondecreasing piecewise polynomial transformation of the original proximities of the chosen degree. The pieces are specified by the number and placement of the interior knots, of which the number can be specified with <i>INKNOT</i> .

### **ORDINAL and SMOOTH Keywords**

<b>UNTIE</b>	<i>Untie ties.</i> Allowing tied proximities to be untied during transformations (also known as the primary approach to ties).
<b>KEEPTIES</b>	<i>Keep ties tied.</i> Keeping tied proximities tied during transformations (also known as secondary approach to ties). This setting is the default.

### **SPLINE Keyword**

<b>INTERCEPT</b>	<i>Include intercept.</i> If <i>INTERCEPT</i> is specified, an intercept is estimated in the transformation. Omitting this keyword sets the lower exterior knot equal to 0.
<b>DEGREE</b>	<i>The degree of the polynomial.</i> If <i>DEGREE</i> is not specified, the degree is assumed to be 2. The integer range of <i>DEGREE</i> is 1, 2, or 3.
<b>INKNOT</b>	<i>The number of interior knots.</i> If <i>INKNOT</i> is not specified, the number of interior knots is assumed to be 1. The integer range of <i>INKNOT</i> is between 0 and the number of different proximities minus 2.

### **Example**

```
PREFSCAL var01 TO var15
  /TRANSFORMATION = ORDINAL(UNTIE) .
```

- In this example, the proximities are ordinally transformed, where tied proximities are allowed to be untied.
- The *CONDITION* subcommand is omitted, and thus, the default conditionality *ROW* is in effect, which implies that the transformation is performed for each row separately.

## **MODEL Subcommand**

*MODEL* defines the scaling model for the analysis if more than one source is present. *IDENTITY* is the default model. The other models are individual differences models.

<b>IDENTITY</b>	<i>Identity model.</i> All sources have the same individual configuration. This model is the default model, and it is not an individual differences model.
<b>WEIGHTED</b>	<i>Weighted Euclidean model.</i> This model is an individual differences model (and equivalent to the <i>INDSCAL</i> model). Each source has an individual space, in which every dimension of the common space is weighted differentially.
<b>GENERALIZED</b>	<i>Generalized Euclidean model.</i> This model is an individual differences model (and equivalent to the <i>IDIOSCAL</i> model). Each source has an individual space that is equal to a differential rotation of the common space, followed by a differential weighting of the dimensions.

- If `IDENTITY` is specified for only one source, this subcommand is silently ignored.
- If an individual differences model is specified for only one source, a warning is issued, and the model is set to `IDENTITY`.

**Example**

```
PREFSCAL var01 TO var15
  /INPUT = SOURCES(6)
  /MODEL = WEIGHTED.
```

- A weighted Euclidean model is fitted for the six specified sources. As indicated on the `INPUT` subcommand, the number of cases must be dividable by 6 in this case.

**RESTRICTIONS Subcommand**

`PREFSCAL` allows (some) coordinates to be fixed in the common space configuration. Fixing an entire set (all row objects or all column objects) corresponds to performing external unfolding.

<b>ROW</b>	<i>Row restriction.</i> <code>PREFSCAL</code> allows one row object, multiple row objects, or all row objects to be free ( <code>NONE</code> ) or fixed to given coordinates ( <code>COORDINATES</code> ).
<b>COLUMN</b>	<i>Column restriction.</i> <code>PREFSCAL</code> allows one column object, multiple column objects, or all column objects to be free ( <code>NONE</code> ) or fixed to given coordinates ( <code>COORDINATES</code> ).

**ROW or COLUMN Keywords**

<b>NONE</b>	<i>No restriction.</i> The specified set of objects ( <code>ROW</code> or <code>COLUMN</code> ) has no restriction.
<b>COORDINATES</b>	Coordinates must be followed by a filespec in parentheses to specify the external IBM® SPSS® Statistics data file that contains the fixed coordinates for the specified set of objects. Following the parenthesized filespec, a variable list can be given. If the variable list is omitted, the procedure automatically selects the first <i>MAXDIM</i> variables in the external SPSS Statistics data file, where <i>MAXDIM</i> is the maximum number of dimensions that are requested for the analysis on <code>/CRITERIA = DIMENSIONS(min, max)</code> .  The number of cases for each variable in the external SPSS Statistics data file must be equal to the number of objects of the specified set ( <code>ROW</code> or <code>COLUMN</code> ).  A missing value can be used to indicate that the coordinate on that dimension is free. The coordinates of objects with nonmissing values are kept fixed during the analysis.

**Example**

```
PREFSCAL var01 TO var15
  /RESTRICTIONS = ROW(NONE)
  /RESTRICTIONS = COLUMN(COORDINATES ("indcol.sav")) .
```

- In this example, there are 15 column objects.



- The coordinates of the row objects are not restricted. Although this specification is the default, it is explicitly stated here in the syntax.
- The column objects have restrictions on the coordinates. The fixed coordinates are specified in the SPSS Statistics data file *indcol.sav*. If *indcol.sav* contains more than two variables, only the first two variables are taken as fixed coordinates, because the maximum dimensionality is 2, and specific variables on the `RESTRICTIONS` subcommand are not given.

## ***PENALTY Subcommand***

The `PENALTY` subcommand specifies the values for the penalty parameters. The two keywords can be used to set the strength and the range of the penalty. The penalty itself is based on the coefficient of variation of the transformed proximities.

<b>LAMBDA</b>	<i>Strength parameter.</i> This parameter sets the strength of the penalty. The default value is 0.75. The range of this parameter is between 0 (exclusive) and 1 (inclusive). The smaller the values of lambda, the stronger the penalty (and vice versa).
<b>OMEGA</b>	<i>Range parameter.</i> This parameter sets the range of the penalty (that is, the moment the penalty becomes active). The parameter must have a non-negative value. If OMEGA is 0, the penalty is inactive. Increasing OMEGA provides a more active penalty. By default (OMEGA = 1.0), the range is equal to the variation coefficient of the original proximities. If OMEGA is increased, the function will search for a solution with a higher variation of the transformed proximities

### ***Example***

```
PREFSCAL var01 TO var09
  /PENALTY = LAMBDA(0.5)
             OMEGA(2.0).
```

- In this example, the variation range is increased by setting OMEGA equal to 2.0.

## ***CRITERIA Subcommand***

You can use `CRITERIA` to set the dimensionality and criteria for terminating the algorithm. You can specify one or more of the following keywords:

<b>DIMENSIONS(min,max)</b>	<i>Minimum and maximum number of dimensions.</i> By default, PREFSCAL computes a solution in two dimensions. The minimum and maximum number of dimensions can be any integer between 1 and the number of objects minus 1 inclusive, as long as the minimum is less than or equal to the maximum. PREFSCAL starts computing a solution in the largest dimensionality and reduces the dimensionality in steps of one, until the lowest dimensionality is reached. Specifying a single value represents both minimum and maximum number of dimensions; thus, <code>DIMENSIONS(4)</code> is equivalent to <code>DIMENSIONS(4,4)</code> .
<b>MAXITER(n)</b>	<i>Maximum number of iterations.</i> By default, $n=5000$ , specifying the maximum number of iterations that are performed while one of the convergence criteria below ( <code>DIFFSTRESS</code> and <code>MINSTRESS</code> ) is not reached. Decreasing this number might give less accurate results, but will take less time. The value $n$ must have a non-negative integer value.

<b>DIFFSTRESS</b>	<i>Convergence criterion.</i> PREFSCAL minimizes the goodness-of-fit index “penalized Stress.” By default, PREFSCAL stops iterating when the relative difference in consecutive penalized Stress values is less than or equal to 0.000001. To obtain a more accurate solution, you can specify a smaller value. The specified value must be nonnegative.
<b>MINSTRESS</b>	<i>Minimum Stress value.</i> By default, PREFSCAL stops iterating when the penalized Stress value itself is small (that is, less than or equal to 0.001). To obtain a more accurate solution, you can specify a smaller value. The specified value must be nonnegative.

**Example**

```
PREFSCAL var01 TO var15
  /CRITERIA = DIMENSIONS(2,4) MAXITER(10000) DIFFSTRESS(1.0E-8) .
```

- The maximum number of dimensions equals 4, and the minimum number of dimensions equals 2. PREFSCAL computes a four-dimensional, three-dimensional, and two-dimensional solution, respectively.
- The maximum number of iterations is set to 10000.
- The relative difference in penalized Stress convergence criterion is sharpened to 1.0E-8.

**PRINT Subcommand**

The PRINT subcommand controls the display of tables. By default, PREFSCAL displays the Stress and fit values for each analysis, the coordinates of the common space, and, if applicable, the individual space weights.

- Omitting the PRINT subcommand or specifying PRINT without keywords is equivalent to specifying COMMON and WEIGHTS.
- If a keyword(s) is specified, only the output for that particular keyword(s) is displayed.
- Inapplicable keywords are silently ignored. That is, a specified keyword for which no output is available—for example, the keyword INDIVIDUAL with only one source specified—will be silently ignored.

<b>NONE</b>	<i>No optional output.</i> Displays only the penalized Stress and corresponding fit values.
<b>INPUT</b>	<i>Input data.</i> Displays tables of the original proximities and, if present, the data weights, the initial configuration, and the fixed coordinates.
<b>MULTIPLE</b>	<i>Multiple random starts.</i> Displays the random number seed and penalized Stress value of each random start.
<b>INITIAL</b>	<i>Initial common space.</i> Displays the coordinates of the initial common space.
<b>HISTORY</b>	<i>History of iterations.</i> Displays the history of iterations of the main algorithm.
<b>MEASURES</b>	<i>Fit measures.</i> Displays different measures. The table contains several goodness-of-fit, badness-of-fit, Stress, and fit values. This setting is specified by default.
<b>DECOMPOSITION</b>	<i>Decomposition of Stress.</i> Displays a objects, rows, and sources decomposition of penalized Stress, including row, column, and source totals.
<b>COMMON</b>	<i>Common space.</i> Displays the coordinates of the common space. This is specified by default.

<b>DISTANCES</b>	<i>Distances.</i> Displays the distances between the objects in the configuration. This keyword must be used in combination with <code>COMMON</code> or <code>INDIVIDUAL</code> to actually produce a table with distances.
<b>WEIGHTS</b>	<i>Individual space weights.</i> Displays the individual space weights, if applicable (that is, if one of the individual differences models is specified on the <code>MODEL</code> subcommand). Depending on the model, the space weights are decomposed in rotation weights and dimension weights, which are also displayed. This setting is specified by default.
<b>INDIVIDUAL</b>	<i>Individual spaces.</i> The coordinates of the individual spaces are displayed only if one of the individual differences models is specified on the <code>MODEL</code> subcommand.
<b>TRANSFORMA-TION</b>	<i>Transformed proximities.</i> Displays the transformed proximities.

### Example

```
PREFSCAL var01 TO var15
  /INPUT = ROWS(42)
  /MODEL = WEIGHTED
  /PRINT = HISTORY COMMON MEASURES.
```

- Here, a weighted Euclidean model is specified with multiple sources.
- The optional output consists of a table with the history of iterations, the coordinates of the common space, and Stress and fit measures.

## PLOT Subcommand

The `PLOT` subcommand controls the display of plots. By default, `PREFSCAL` displays the object points of the common space and, if applicable, the individual space weights.

- Omitting the `PLOT` subcommand or specifying `PLOT` without keywords produces the default plots.
- If a keyword(s) is specified, only the plot for that particular keyword(s) is displayed.
- Inapplicable keywords (for example, `STRESS` with equal minimum and maximum number of dimensions on the `CRITERIA` subcommand) are silently ignored.
- Multiple value lists are allowed for `INDIVIDUAL`, `TRANSFORMATIONS`, `SHEPARD`, `FIT`, and `RESIDUALS`. For each value list, a separate plot will be displayed.

<b>NONE</b>	<i>No plots.</i> <code>PREFSCAL</code> does not produce any plot.
<b>MULTIPLE</b>	<i>Multiple random starts.</i> Displays a stacked histogram of penalized Stress, displaying both Stress and penalty.
<b>INITIAL</b>	<i>Initial common space.</i> Displays a scatterplot matrix of the coordinates of the initial common space.
<b>STRESS</b>	<i>Scree plot.</i> Produces a lineplot of penalized Stress versus dimensions. This plot is only produced if the maximum number of dimensions is larger than the minimum number of dimensions.
<b>COMMON</b>	<i>Common space.</i> A scatterplot matrix of coordinates of the common space is displayed. This setting is the default.

**WEIGHTS**

*Individual space weights.* A scatterplot is produced for the individual space weights. This setting is only applicable if one of the individual differences models is specified on the `MODEL` subcommand. For the weighted Euclidean model, the weights for all sources are displayed in a plot, with one dimension on each axis. For the generalized Euclidean model, one plot is produced per dimension, indicating both rotation and weighting of that dimension for each source.

**INDIVIDUAL(valuelist)**

*Individual spaces.* The coordinates of the individual spaces are displayed in scatterplot matrices. This setting is only applicable if one of the individual differences models is specified on the `MODEL` subcommand. For each source that is specified on the value list, a scatterplot matrix of coordinates of the individual space is displayed. The sources are specified by a number between 1 and the total number of sources or is specified by a value from the `sourceid`, which is specified on the `INPUT` subcommand.

**TRANSFORMATIONS(valuelist)**

*Transformation plots.* A line plot is produced of the original proximities versus the transformed proximities. On the value list, the names (identifiers) for which the plot is to be produced must be specified. Because the `CONDITION` subcommand allows for the specification of multiple transformation lists, the value lists depend on the conditionality. In case of row-conditional transformations, the names are row identifiers (either a number between 1 and the total number of rows, or a value from the `rowid`, which is specified on the `INPUT` subcommand). In the case of matrix-conditional transformations, the values indicate sources identifiers (either a number between 1 and the total number of sources, or a value from the `sourceid`, which is specified on the `INPUT` subcommand). An unconditional transformation only consists of one list and does not allow further specification.

**SHEPARD(valuelist)**

*Shepard plots.* The original proximities versus both transformed proximities and distances. The distances are indicated by points, and the transformed proximities are indicated by a line. On the value list, the names (identifiers) for which the plot is to be produced must be specified. Because the `CONDITION` subcommand allows for the specification of multiple transformation lists, the value lists depend on the conditionality. In case of row-conditional transformations, the names are row identifiers (either a number between 1 and the total number of rows, or a value from the `rowid`, which is specified on the `INPUT` subcommand). In the case of matrix-conditional transformations, the values indicate sources identifiers (either a number between 1 and the total number of sources, or a value from the `sourceid`, which is specified on the `INPUT` subcommand). An unconditional transformation only consists of one list and does not allow further specification.

<b>FIT(valuelist)</b>	<i>Scatterplots of Fit.</i> The transformed proximities versus the distances are plotted in a scatterplot. On the value list, the names (identifiers) of the sources for which the plot is to be produced must be specified. The sources are specified by a number between 1 and the total number of sources or are specified by a value from the sourceid, which is specified on the INPUT subcommand.
<b>RESIDUALS(valuelist)</b>	<i>Residuals plots.</i> The transformed proximities versus the residuals (transformed proximities minus distances) are plotted in a scatterplot. On the value list, the names (identifiers) of the sources for which the plot is to be produced must be specified. The sources are specified by a number between 1 and the total number of sources or are specified by a value from the sourceid, which is specified on the INPUT subcommand.

**Example**

```
PREFSCAL var01 TO var15
  /INPUT = SOURCE(6)
  /MODEL = WEIGHTED
  /CRITERIA = DIMENSIONS(3)
  /PLOT = COMMON
        INDIVIDUAL(2)
        TRANSFORMATIONS(1 TO 42)(1 2)
        FIT(2).
```

- Here, the syntax specifies a weighted Euclidean model with six sources in three dimensions.
- COMMON produces a scatterplot matrix defined by dimensions 1, 2, and 3.
- A scatterplot matrix with threedimensions is produced only for the source 2.
- Two transformation plots are produced, one plot with all 42 rows and one plot with only row 1 and 2. Rows are specified with the TRANSFORMATIONS keyword because the default value on CONDITION is ROW.
- A scatterplot of fit is produced for the source 2.

**OPTIONS Subcommand**

The OPTIONS subcommand specifies additional markings for the row objects in plots. For this purpose, the values of variables are used to specify markers and colors for the row objects.

<b>MARKERS(variable)</b>	<i>Row object markers.</i> The values of the variable are used to cycle through all possible markers.
<b>COLORS(variable)</b>	<i>Row object colors.</i> The values of the variable are used to cycle through all colors.

**Example**

```
DATA LIST
  /var01 TO var15 rowid gender age.
PREFSCAL var01 TO var15
  /INPUT = ROW(rowid)
```

```
/OPTIONS = MARKERS(gender) COLORS(age) .
```

- In the joint plot of the common space configuration, the row objects are labeled with the values or value labels of the variable *rowid*. Additionally, the points are marked according to the values on the variable *gender* and are colored depending on the values of the variable *age*.

## OUTFILE Subcommand

OUTFILE saves coordinates of the common space, individual space weights, distances, and transformed proximities to an external IBM® SPSS® Statistics data file or previously declared dataset (DATASET DECLARE command). The data file/dataset name must be different for each keyword.

### COMMON (‘savfile’|‘dataset’)

*Common space coordinates.* The coordinates of the common space. The columns (variables) represent the dimensions *DIM\_1*, *DIM\_2*, ..., *DIM\_n* of the common space. The number of cases in the external file equals the total number of objects (row plus column objects).

### WEIGHTS (‘savfile’|‘dataset’)

*Individual space weights.* The individual space weights. The columns represent the dimensions *DIM\_1*, *DIM\_2*, ..., *DIM\_n* of the space weights. The number of cases depends on the individual differences model specified on the MODEL subcommand. The weighted Euclidean model uses diagonal weight matrices. Only the diagonals are written to file, and the number of cases is equal to the number of sources. The generalized Euclidean model has full-rank nonsingular weight matrices, one matrix for each source. The matrices are stacked beneath each other in the external SPSS Statistics data file. The number of cases equals the number of sources times the number of dimensions.

### DISTANCES (‘savfile’|‘dataset’)

*Distances.* The matrices containing the distances between the objects for each source are stacked beneath each other in the external SPSS Statistics data file. The number of variables in the data file is equal to the total number of objects (*ROW\_1*, *ROW\_2*, ..., *ROW\_n*, *COL\_1*, *COL\_2*, ..., *COL\_m*). The number of cases in the data file is equal to the total number of objects times the number of sources.

### TRANSFORMATION (‘file’|‘dataset’)

*Transformed proximities.* The matrices containing the transformed proximities for each source are stacked beneath each other in the external SPSS Statistics data file. The number of variables in the external file is equal to the total number of objects (*ROW\_1*, *ROW\_2*, ..., *ROW\_n*, *COL\_1*, *COL\_2*, ..., *COL\_m*). The number of cases in the external file is equal to the total number of objects times the number of sources.

### Example

```
PREFSCAL var01 TO var15  
/OUTFILE = COMMON('/data/start.sav') .
```

- Here, the coordinates of the common space are written to the external SPSS Statistics data file *start.sav*.
- Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files.

# **PRESERVE**

PRESERVE

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Overview**

PRESERVE stores current SET specifications that can later be restored by the RESTORE command. PRESERVE and RESTORE are especially useful with the macro facility. PRESERVE-RESTORE sequences can be nested up to five levels.

### **Basic Specification**

The only specification is the command keyword. PRESERVE has no additional specifications.

### **Limitations**

PRESERVE does not store SET THREADS and SET MCACHE settings.

## **Example**

```
PRESERVE.  
SET TNUMBERS=BOTH.  
FREQUENCIES VARIABLES=DIVISION  
RESTORE.  
FREQUENCIES VARIABLES=DIVISION.
```

- PRESERVE stores all current SET specifications.
- SET TNUMBERS=BOTH changes the TNUMBERS setting to display both values and value labels in pivot tables.
- The pivot table results from the first FREQUENCIES command will display both category values and labels.
- RESTORE reestablishes the SET specifications that were in effect prior to when PRESERVE was specified.
- The pivot table results from the second FREQUENCIES command will use the TNUMBERS setting in effect prior to PRESERVE (the default is to display only labels).



# PRINCALS

PRINCALS is available in the Categories option.

```
PRINCALS VARIABLES=varlist(max)

[/ANALYSIS=varlist[({ORDI**})]]
                {SNOM  }
                {MNOM  }
                {NUME  }

[/NOBSERVATIONS=value]

[/DIMENSION={2**  }]
                {value}

[/MAXITER={100**}]
                {value}

[/CONVERGENCE={.00001**}]
                {value  }

[/PRINT=[DEFAULT] [FREQ**] [EIGEN**] [LOADINGS**] [QUANT]
                [HISTORY] [CORRELATION] [OBJECT] [ALL] [NONE]]

[/PLOT=[NDIM=({1    ,2    }**)]
                {value,value}
                {ALL  ,MAX  }
                [DEFAULT] [(n)]] [OBJECT**] [(varlist)] [(n)]]
                [QUANT**] [(varlist)] [(n)]] [LOADINGS] [(n)]]
                [ALL] [(n)]] [NONE]]

[/SAVE=[rootname] [(value)]]

[/MATRIX=OUT({*    })]
                {'file'|'dataset'}
```

**\*\***Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Overview

PRINCALS (*principal components analysis by means of alternating least squares*) analyzes a set of variables for major dimensions of variation. The variables can be of mixed optimal scaling levels, and the relationships among observed variables are not assumed to be linear.

### Options

**Optimal Scaling Level.** You can specify the optimal scaling level for each variable to be used in the analysis.

**Number of Cases.** You can restrict the analysis to the first *n* observations.

**Number of Dimensions.** You can specify how many dimensions PRINCALS should compute.

**Iterations and Convergence.** You can specify the maximum number of iterations and the value of a convergence criterion.

**Display Output.** The output can include all available statistics, only the default statistics, or only the specific statistics you request. You can also control whether some of these statistics are plotted.

**Saving Scores.** You can save object scores in the active dataset.

**Writing Matrices.** You can write a matrix data file containing category quantifications and loadings for use in further analyses.

### ***Basic Specification***

- The basic specification is the PRINCALS command and the VARIABLES subcommand. PRINCALS performs the analysis assuming an ordinal level of optimal scaling for all variables and uses all cases to compute a two-dimensional solution. By default, marginal frequencies, eigenvalues, and summary measures of fit and loss are displayed, and quantifications and object scores are plotted.

### ***Subcommand Order***

- The VARIABLES subcommand must precede all others.
- Other subcommands can appear in any order.

### ***Operations***

- If the ANALYSIS subcommand is specified more than once, PRINCALS is not executed. For all other subcommands, only the last occurrence of each subcommand is executed.
- PRINCALS treats every value in the range of 1 to the maximum value specified on VARIABLES as a valid category. Use the AUTORECODE or RECODE command if you want to recode a categorical variable with nonsequential values or with a large number of categories to avoid unnecessary output. For variables treated as numeric, recoding is *not* recommended because the intervals between consecutive categories will not be maintained.

### ***Limitations***

- String variables are not allowed; use AUTORECODE to recode nominal string variables into numeric ones before using PRINCALS.
- The data must be positive integers. Zeros and negative values are treated as system-missing and are excluded from the analysis. Fractional values are truncated after the decimal and are included in the analysis. If one of the levels of a categorical variable has been coded 0 or a negative value and you want to treat it as a valid category, use the AUTORECODE or RECODE command to recode the values of that variable (see AUTORECODE and RECODE for more information).
- PRINCALS ignores user-missing value specifications. Positive user-missing values less than the maximum value on the VARIABLES subcommand are treated as valid category values and are included in the analysis. If you do not want the category included, you can use COMPUTE or RECODE to change the value to something outside of the valid range. Values outside of the range (less than 1 or greater than the maximum value) are treated as system-missing.

## Example

```
PRINCALS VARIABLES=ACOLA BCOLA(2) PRICEA PRICEB(5)
/ANALYSIS=ACOLA BCOLA(SNOM) PRICEA PRICEB(NUMB)
/PRINT=QUANT OBJECT.
```

- `VARIABLES` defines the variables and their maximum number of levels.
- The `ANALYSIS` subcommand specifies that variables `ACOLA` and `BCOLA` are single nominal (`SNOM`) and that variables `PRICEA` and `PRICEB` are numeric (`NUMB`).
- The `PRINT` subcommand lists the category quantifications and object scores.
- By default, plots of the category quantifications and the object scores are produced.

## VARIABLES Subcommand

`VARIABLES` specifies all of the variables that will be used in the current `PRINCALS` procedure.

- The `VARIABLES` subcommand is required and precedes all other subcommands. The actual word `VARIABLES` can be omitted.
- Each variable or variable list is followed by the maximum number of categories (levels) in parentheses.
- The number specified in parentheses indicates the number of categories *and* the maximum category value. For example, `VARI(3)` indicates that `VARI` has three categories coded 1, 2, and 3. However, if a variable is not coded with consecutive integers, the number of categories used in the analysis will differ from the number of observed categories. For example, if a three category variable is coded {2, 4, 6}, the maximum category value is 6. The analysis treats the variable as having six categories, three of which are not observed and receive quantifications of 0.
- To avoid unnecessary output, use the `AUTORECODE` or `RECODE` command before `PRINCALS` to recode a categorical variable that was coded with nonsequential values. As noted in “Limitations,” recoding is *not* recommended with variables treated as numeric (see `AUTORECODE` and `RECODE` for more information).

### Example

```
DATA LIST FREE/V1 V2 V3.
BEGIN DATA
3 1 1
6 1 1
3 1 3
3 2 2
3 2 2
6 2 2
6 1 3
6 2 2
3 2 2
6 2 1
END DATA.
AUTORECODE V1 /INTO NEWVAR1.
PRINCALS VARIABLES=NEWVAR1 V2(2) V3(3).
```

- `DATA LIST` defines three variables, `V1`, `V2`, and `V3`.

- *V1* has two levels, coded 3 and 6, *V2* has two levels, coded 1 and 2, and *V3* has three levels, coded 1, 2, and 3.
- The `AUTORECODE` command creates *NEWVARI* containing recoded values of *V1*. Values of 3 are recoded to 1, and values of 6 are recoded to 2.
- A maximum value of 2 can then be specified on the `VARIABLES` subcommand as the maximum category value for both *NEWVARI* and *V2*. A maximum value of 3 is specified for *V3*.

## ***ANALYSIS Subcommand***

`ANALYSIS` specifies the variables to be used in the computations and the optimal scaling level used by `PRINCALS` to quantify each variable or variable list.

- If `ANALYSIS` is not specified, an ordinal level of optimal scaling is assumed for all variables.
- The specification on `ANALYSIS` is a variable list and an optional keyword in parentheses to indicate the optimal scaling level.
- The variables on the variable list must also be specified on the `VARIABLES` subcommand.
- Variables listed on the `VARIABLES` subcommand but not on the `ANALYSIS` subcommand can still be used to label object scores on the `PLOT` subcommand.

The following keywords can be specified to indicate the optimal scaling level:

<b>MNOM</b>	<i>Multiple nominal.</i> The quantifications can be different for each dimension. When all variables are multiple nominal, <code>PRINCALS</code> gives the same results as <code>HOMALS</code> .
<b>SNOM</b>	<i>Single nominal.</i> <code>PRINCALS</code> gives only one quantification for each category. Objects in the same category (cases with the same value on a variable) obtain the same quantification. When <code>DIMENSION=1</code> and all variables are <code>SNOM</code> , this solution is the same as that of the first <code>HOMALS</code> dimension.
<b>ORDI</b>	<i>Ordinal.</i> This is the default for variables listed without optimal scaling levels and for all variables if the <code>ANALYSIS</code> subcommand is not used. The order of the categories of the observed variable is preserved in the quantified variable.
<b>NUME</b>	<i>Numerical.</i> This is the interval or ratio level of optimal scaling. <code>PRINCALS</code> assumes that the observed variable already has numerical values for its categories. When all variables are at the numerical level, the <code>PRINCALS</code> analysis is analogous to classical principal components analysis.

These keywords can apply to a variable list as well as to a single variable. Thus, the default `ORDI` is not applied to a variable without a keyword if a subsequent variable on the list has a keyword.

## ***NOBSERVATIONS Subcommand***

`NOBSERVATIONS` specifies how many cases are used in the analysis.

- If `NOBSERVATIONS` is not specified, all available observations in the active dataset are used.
- `NOBSERVATIONS` is followed by an integer indicating that the first *n* cases are to be used.

## ***DIMENSION Subcommand***

`DIMENSION` specifies the number of dimensions that you want `PRINCALS` to compute.

- If you do not specify the `DIMENSION` subcommand, `PRINCALS` computes two dimensions.
- `DIMENSION` is followed by an integer indicating the number of dimensions.
- If all of the variables are `SNOM` (single nominal), `ORDI` (ordinal), or `NUME` (numerical), the maximum number of dimensions you can specify is the smaller of the number of observations minus 1 *or* the total number of variables.
- If some or all of the variables are `MNOM` (multiple nominal), the maximum number of dimensions is the smaller of the number of observations minus 1 *or* the total number of valid `MNOM` variable levels (categories) plus the number of `SNOM`, `ORDI`, and `NUME` variables, minus the number of `MNOM` variables without missing values.
- `PRINCALS` adjusts the number of dimensions to the maximum if the specified value is too large.
- The minimum number of dimensions is 1.

## ***MAXITER Subcommand***

`MAXITER` specifies the maximum number of iterations `PRINCALS` can go through in its computations.

- If `MAXITER` is not specified, `PRINCALS` will iterate up to 100 times.
- `MAXITER` is followed by an integer indicating the maximum number of iterations allowed.

## ***CONVERGENCE Subcommand***

`CONVERGENCE` specifies a convergence criterion value. `PRINCALS` stops iterating if the difference in total fit between the last two iterations is less than the `CONVERGENCE` value.

- If `CONVERGENCE` is not specified, the default value is 0.00001.
- The specification on `CONVERGENCE` is a convergence criterion value.

## ***PRINT Subcommand***

`PRINT` controls which statistics are included in your output. The default output includes frequencies, eigenvalues, loadings, and summary measures of fit and loss.

`PRINT` is followed by one or more of the following keywords:

<b>FREQ</b>	<i>Marginal frequencies for the variables in the analysis.</i>
<b>HISTORY</b>	<i>History of the iterations.</i>
<b>EIGEN</b>	<i>Eigenvalues.</i>
<b>CORRELATION</b>	<i>Correlation matrix for the transformed variables in the analysis.</i> No correlation matrix is produced if there are any missing data.
<b>OBJECT</b>	<i>Object scores.</i>
<b>QUANT</b>	<i>Category quantifications and category coordinates for <code>SNOM</code>, <code>ORDI</code>, and <code>NUME</code> variables and category quantifications in each dimension for <code>MNOM</code> variables.</i>
<b>LOADINGS</b>	<i>Component loadings for <code>SNOM</code>, <code>ORDI</code>, and <code>NUME</code> variables.</i>
<b>DEFAULT</b>	<i><code>FREQ</code>, <code>EIGEN</code>, <code>LOADINGS</code>, and <code>QUANT</code>.</i>

<b>ALL</b>	<i>All of the available statistics.</i>
<b>NONE</b>	<i>Summary measures of fit.</i>

## **PLOT Subcommand**

PLOT can be used to produce plots of category quantifications, object scores, and component loadings.

- If PLOT is not specified, plots of the object scores and the quantifications are produced.
- No plots are produced for a one-dimensional solution.

PLOT is followed by one or more of the following keywords:

<b>LOADINGS</b>	<i>Plots of the component loadings of SNOM, ORDI, and NUME variables.</i>
<b>OBJECT</b>	<i>Plots of the object scores.</i>
<b>QUANT</b>	<i>Plots of the category quantifications for MNOM variables and plots of the single-category coordinates for SNOM, ORDI, and NUME variables.</i>
<b>DEFAULT</b>	<i>QUANT and OBJECT.</i>
<b>ALL</b>	<i>All available plots.</i>
<b>NONE</b>	<i>No plots.</i>

- The keywords OBJECT and QUANT can each be followed by a variable list in parentheses to indicate that plots should be labeled with these variables. For QUANT, the variables must be specified on both the VARIABLES and ANALYSIS subcommands. For OBJECT, the variables must be specified on VARIABLES but need not appear on the ANALYSIS subcommand. This means that variables not included in the computations can still be used to label OBJECT plots. If the variable list is omitted, only the default plots are produced.
- Object scores plots labeled with variables that appear on the ANALYSIS subcommand use category labels corresponding to all categories within the defined range. Objects in a category that is outside the defined range are labeled with the label corresponding to the next category greater than the defined maximum category.
- Object scores plots labeled with variables not included on the ANALYSIS subcommand use all category labels, regardless of whether or not the category value is inside the defined range.
- All of the keywords except NONE can be followed by an integer in parentheses to indicate how many characters of the variable or value label are to be used on the plot. (If you specify a variable list after OBJECT or QUANT, you can specify the value in parentheses after the list.) The value can range from 1 to 20. If the value is omitted, 12 characters are used. Spaces between words count as characters.
- The LOADINGS plots and one of the QUANT plots use variable labels; all other plots that use labels use value labels.
- If a variable label is missing, the variable name is used for that variable. If a value label is missing, the actual value is used.
- You should make sure that your variable and value labels are unique by at least one letter in order to distinguish them on the plots.
- When points overlap, the points involved are described in a summary following the plot.

**Example**

```
PRINCALS VARIABLES COLA1 (4) COLA2 (4) COLA3 (4) COLA4 (2)
/ANALYSIS COLA1 COLA2 (SNOM) COLA3 (ORDI) COLA4 (ORDI)
/PLOT OBJECT(COLA4) .
```

- Four variables are included in the analysis.
- OBJECT requests a plot of the object scores labeled with the values of *COLA4*. Any object whose *COLA4* value is not 1 or 2 is labeled 3 (or the value label for category 3, if defined).

**Example**

```
PRINCALS VARIABLES COLA1 (4) COLA2 (4) COLA3 (4) COLA4 (2)
/ANALYSIS COLA1 COLA2 (SNOM) COLA3 (ORDI)
/PLOT OBJECT(COLA4) .
```

- Three variables are included in the analysis.
- OBJECT requests a plot of the object scores labeled with the values of *COLA4*, a variable not included in the analysis. Objects are labeled using all values of *COLA4*.

In addition to the plot keywords, the following can be specified:

**NDIM**      *Dimension pairs to be plotted.* NDIM is followed by a pair of values in parentheses. If NDIM is not specified, plots are produced for dimension 1 versus dimension 2.

- The first value indicates the dimension that is plotted against all higher dimensions. This value can be any integer from 1 to the number of dimensions minus 1.
- The second value indicates the highest dimension to be used in plotting the dimension pairs. This value can be any integer from 2 to the number of dimensions.
- The keyword ALL can be used instead of the first value to indicate that all dimensions are paired with higher dimensions.
- The keyword MAX can be used instead of the second value to indicate that plots should be produced up to, and including, the highest dimension fit by the procedure.

**Example**

```
PRINCALS COLA1 COLA2 COLA3 COLA4 (4)
/PLOT NDIM(1,3) QUANT(5) .
```

- The NDIM(1,3) specification indicates that plots should be produced for two dimension pairs—dimension 1 versus dimension 2 and dimension 1 versus dimension 3.
- QUANT requests plots of the category quantifications. The (5) specification indicates that the first five characters of the value labels are to be used on the plots.

**Example**

```
PRINCALS COLA1 COLA2 COLA3 COLA4 (4)
```

```
/PLOT NDIM(ALL,3) QUANT(5) .
```

- This plot is the same as above except for the ALL specification following NDIM. This indicates that all possible pairs up to the second value should be plotted, so QUANT plots will be produced for dimension 1 versus dimension 2, dimension 2 versus dimension 3, and dimension 1 versus dimension 3.

## **SAVE Subcommand**

SAVE lets you add variables containing the object scores computed by PRINCALS to the active dataset.

- If SAVE is not specified, object scores are not added to the active dataset.
- A variable rootname can be specified on the SAVE subcommand to which PRINCALS adds the number of the dimension. Only one rootname can be specified, and it can contain up to six characters.
- If a rootname is not specified, unique variable names are automatically generated. The variable names are *PRIn\_m*, where *n* is a dimension number and *m* is a set number. If three dimensions are saved, the first set of names is *PR11\_1*, *PR12\_1*, and *PR13\_1*. If another PRINCALS is then run, the variable names for the second set are *PR11\_2*, *PR12\_2*, *PR13\_2*, and so on.
- Following the name, the number of dimensions for which you want to save object scores can be listed in parentheses. The number cannot exceed the value of the DIMENSION subcommand.
- If the number of dimensions is not specified, the SAVE subcommand saves object scores for all dimensions.
- If you replace the active dataset by specifying an asterisk (\*) on a MATRIX subcommand, the SAVE subcommand is not executed.
- The prefix should be unique for each PRINCALS command in the same session. If it is not, PRINCALS replaces the prefix with *DIM*, *OBJ*, or *OBSAVE*. If all of these already exist, SAVE is not executed.

### **Example**

```
PRINCALS CAR1 CAR2 CAR3(5) PRICE (10)
/ANALYSIS=CAR1 TO CAR3 (SNOM) PRICE (NUM)
/DIMENSIONS=3
/SAVE=DIM(2) .
```

- Three nominal variables, *CAR1*, *CAR2*, and *CAR3*, each with five categories, and one numerical (interval level) variable, with ten categories, are analyzed in this PRINCALS example.
- The DIMENSIONS subcommand requests results for three dimensions.
- SAVE adds the object scores from the first two dimensions to the active dataset. The names of these new variables will be *DIM00001* and *DIM00002*, respectively.



## **MATRIX Subcommand**

The `MATRIX` subcommand is used to write category quantifications, single-category coordinates, and component loadings to a matrix data file.

- The specification on `MATRIX` is the keyword `OUT` and a quoted file specification of previously declared dataset name (`DATASET DECLARE` command), enclosed in parentheses.
- You can specify an asterisk (\*) instead of a file to replace the active dataset .
- The category quantifications, coordinates, and component loadings are written to the same file.
- The matrix data file has one case for each value of each original variable.

The variables of the matrix data file and their values are:

<b>ROWTYPE_</b>	<i>String variable rowtype_ containing value QUANT for the category quantifications, SCOOR_ for single-category coordinates, MCOOR_ for multiple-category coordinates, and LOADING_ for the component scores.</i>
<b>LEVEL</b>	<i>String variable containing the values (or value labels if present) of each original variable for category quantifications. For cases with ROWTYPE_=LOADING_, the value of LEVEL is blank.</i>
<b>VARNAME_</b>	<i>String variable containing the original variable names.</i>
<b>VARTYPE_</b>	<i>String variable containing values MULTIPLE, SINGLE N, ORDINAL, or NUMERICAL, depending on the optimal scaling level specified for the variable.</i>
<b>DIM1...DIMn</b>	<i>Numeric variables containing category quantifications, the single-category coordinates, and component loadings for each dimension. Each variable is labeled DIMn, where n represents the dimension number. The single-category coordinates and component loadings are written only for SNOM, ORDI, and NUME variables.</i>

# PRINT

```
PRINT [OUTFILE=file] [ENCODING='encoding specification']
  [RECORDS={1}] [{NOTABLE}]
    {n}   {TABLE }

/{1   } varlist [{col location [(format)]}] [varlist...]
 {rec #}         {(format list)   }
                  {*}              }

[/{2   }...]
 {rec #}
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- ENCODING subcommand added for Unicode support.

## **Example**

```
PRINT / MOHIRED YRHIRED DEPT SALARY NAME.
```

## **Overview**

PRINT displays the values of variables for each case in the data. PRINT is simple enough for a quick check on data definitions and transformations and flexible enough for formatting simple reports.

### **Options**

**Formats.** You can specify formats for the variables (see [Formats on p. 1590](#)).

**Strings.** You can specify string values within the variable specifications. The strings can be used to label values or to create extra space between values. Strings can also be used as column headings. (See [Strings on p. 1591](#).)

**Output File.** You can use the `OUTFILE` subcommand to direct the output to a specified file.

**Summary Table.** You can use the `TABLE` subcommand to display a table that summarizes variable formats.

### **Basic Specification**

The basic specification is a slash followed by a variable list. The output displays values for all variables that are named on the list.

### ***Subcommand Order***

Subcommands can be specified in any order. However, all subcommands must be specified before the slash that precedes the start of the variable specifications.

### ***Syntax Rules***

- A slash must precede the variable specifications. The first slash begins the definition of the first (and possibly only) line per case of the PRINT display.
- Specified variables must already exist, but they can be numeric, string, scratch, temporary, or system variables. Subscripted variable names, such as  $X(I)$  for the first element in vector  $X$ , cannot be used.
- Keyword ALL can be used to display the values of all user-defined variables in the active dataset.

### ***Operations***

- PRINT is executed once for each case that is constructed from the data file.
- PRINT is a transformation and is not executed unless it is followed by a procedure or the EXECUTE command.
- Because PRINT is a transformation command, the output might be mixed with casewise procedure output. Procedures that produce individual case listings (such as LIST) should not be used immediately after PRINT. An intervening EXECUTE or procedure command should be specified.
- Values are displayed with a blank space between them. However, if a format is specified for a variable, the blank space for that variable's values is suppressed.
- Values are displayed in the output as the data are read. The PRINT output appears before the output from the first procedure.
- If more variables are specified than can be displayed in 132 columns or within the width that is specified on SET WIDTH, the program displays an error message. You must reduce the number of variables or split the output into several records.
- User-missing values are displayed exactly like valid values. System-missing values are represented by a period.

## ***Examples***

### ***Displaying Values for a Selected List of Variables***

```
PRINT / MOHIRED YRHIRED DEPT SALARY NAME .
FREQUENCIES VARIABLES=DEPT .
```

- PRINT displays values for each variable on the variable list. The FREQUENCIES procedure reads the data and causes PRINT to be executed.
- All variables are displayed by using their dictionary formats. One blank space separates the values of each variable.

### Displaying Values for All User-Defined Variables

```
PRINT /ALL.
EXECUTE.
```

- `PRINT` displays values for all user-defined variables in the active dataset. The `EXECUTE` command executes `PRINT`.

## Formats

By default, `PRINT` uses the dictionary print formats. You can specify formats for some or all variables that are specified on `PRINT`. For a string variable, the specified format must have a width at least as large as the width of the dictionary format. String values are truncated if the specified width is smaller than the width of the dictionary format.

- Format specifications can be either column-style or FORTRAN-like (see `DATA LIST`). The column location that is specified with column-style formats or that is implied with FORTRAN-like formats refers to the column in which the variable will be displayed.
- A format specification following a list of variables applies to all variables in the list. Use an asterisk to prevent the specified format from applying to variables that precede the asterisk. The specification of column locations implies a default print format, and that format applies to all previous variables if no asterisk is used.
- Printable numeric formats are `F`, `COMMA`, `DOLLAR`, `CC`, `DOT`, `N`, `E`, `PCT`, `PIBHEX`, `RBHEX`, `Z`, and the date and time formats. Printable string formats are `A` and `AHEX`. Note that hex and binary formats use different widths. For example, the `AHEX` format must have a width that is twice the width of the corresponding `A` format. For more information about specifying formats and more information about the available formats, see `DATA LIST` and Variable Types and Formats on p. 54.
- Format specifications are in effect only for the `PRINT` command. The specifications do not change the dictionary print formats.
- When a format is specified for a variable, the automatic blank following the variable in the output is suppressed. To preserve the blank between variables, use a string (see Strings on p. 1591), specify blank columns in the format, or use an `X` or `T` format element (see `DATA LIST` for information about `X` and `T`).

### Example

```
PRINT / TENURE (F2.0) ' ' MOHIRED YRHIRED DEPT *
      SALARY85 TO SALARY88 (4(DOLLAR8,1X)) NAME.
EXECUTE.
```

- Format `F2.0` is specified for `TENURE`. A blank string is specified after `TENURE` because the automatic blank following the variable is suppressed by the format specification.
- `MOHIRED`, `YRHIRED`, and `DEPT` are displayed with default formats because the asterisk prevents them from receiving the `DOLLAR8` format that is specified for `SALARY85` to `SALARY88`. The automatic blank is preserved for `MOHIRED`, `YRHIRED`, and `DEPT`, but the blank is suppressed for `SALARY85` to `SALARY88` by the format specification. The `1X` format

element is therefore specified with `DOLLAR8` to add one blank after each value of `SALARY85` to `SALARY88`.

- `NAME` uses the default dictionary format.

## Strings

You can specify string values within the variable list. Strings must be enclosed in quotes.

- If a format is specified for a variable list, the application of the format is interrupted by a specified string. Thus, the string has the same effect within a variable list as an asterisk.
- Strings can be used to create column headings for the displayed variables. The `PRINT` command that specifies the column headings must be used within a `DO IF-END IF` structure. If you want the column headings to begin a new page in the output, use a `PRINT EJECT` command (rather than `PRINT`) to specify the headings (see `PRINT EJECT`).

### Including Strings in the Output

```
PRINT / NAME 'HIRED=' MOHIRED(F2) '/' YRHIRED
        ' SALARY=' SALARY (DOLLAR8).
EXECUTE.
```

- Three strings are specified. The strings `HIRED=` and `SALARY=` label the values being displayed. The slash that is specified between `MOHIRED` and `YRHIRED` creates a composite hiring date. The `F2` format is supplied for variable `MOHIRED` in order to suppress the blank that would follow it if the dictionary format were used.
- `NAME` and `YRHIRED` are displayed with default formats. The `'HIRED='` specification prevents the `F2` format from applying to `NAME`, and the `'SALARY='` specification prevents the `DOLLAR8` format from applying to `YRHIRED`.

### Setting Up Column Headers

```
DO IF $CASENUM EQ 1.
PRINT /   NAME ' 1 'DEPT' 25 'HIRED' 30 ' SALARY' 35.
END IF.
PRINT / NAME DEPT *
        MOHIRED 30-31 '/' YRHIRED *
        SALARY 35-42(DOLLAR).
EXECUTE.
```

- The first `PRINT` command specifies strings only. The integer after each string specifies the beginning column number of the string. The strings will be used as column headings for the variables. `DO IF $CASENUM EQ 1` causes the first `PRINT` command to be executed only once, as the first case is processed. `END IF` closes the structure.
- The second `PRINT` command specifies the variables to be displayed. This command is executed once for each case in the data. Column locations are specified to align the values with the column headings. In this example, the `T` format element could also have been used

to align the variables and the column headings. For example, `MOHIRED (T30,F2)` begins the display of values for variable *MOHIRED* in column 30.

- The asterisk after *DEPT* prevents the format that is specified for *MOHIRED* from applying to *NAME* and *DEPT*. The asterisk after *YRHIRED* prevents the format that is specified for *SALARY* from applying to *YRHIRED*.

## **RECORDS Subcommand**

`RECORDS` indicates the total number of lines that are displayed per case. The number that is specified on `RECORDS` is informational only. The actual specification that causes variables to be displayed on a new line is a slash within the variable specifications. Each new line is requested by another slash.

- `RECORDS` must be specified before the slash that precedes the start of the variable specifications.
- The only specification on `RECORDS` is an integer to indicate the number of records for the output. If the number does not agree with the actual number of records that are indicated by slashes, the program issues a warning and ignores the specification on `RECORDS`.
- Specifications for each line of output must begin with a slash. An integer can follow the slash, indicating the line on which values are to be displayed. The integer is informational only and cannot be used to rearrange the order of records in the output. If the integer does not agree with the actual record number that is indicated by the number of slashes in the variable specifications, the integer is ignored.
- A slash that is not followed by a variable list generates a blank line in the output.

### **Examples**

```
PRINT RECORDS=3 /EMPLOYID NAME DEPT
                /EMPLOYID TENURE SALARY
                /.
```

EXECUTE.

- `PRINT` displays the values of an individual's name and department on one line, displays tenure and salary on the next line, and displays the employee identification number on both lines, followed by a blank third line. Two lines are displayed for each case, and cases in the output are separated by a blank line.

```
PRINT RECORDS=3 /1 EMPLOYID NAME DEPT
                /2 EMPLOYID TENURE SALARY
                /3.
```

- This `PRINT` command is equivalent to the command in the preceding example.

```
PRINT / EMPLOYID NAME DEPT / EMPLOYID TENURE SALARY /.
```

- This `PRINT` command is equivalent to the commands in the two preceding examples.

## ***OUTFILE Subcommand***

OUTFILE specifies a file for the output from the PRINT command. By default, PRINT output is included with the rest of the output from the session.

- OUTFILE must be specified before the slash that precedes the start of the variable specifications.
- The output from PRINT cannot exceed 132 characters, even if the external file is defined with a longer record length.

### ***Example***

```
PRINT OUTFILE=PRINTOUT
  /1 EMPLOYID DEPT SALARY /2 NAME.
EXECUTE.
```

- OUTFILE specifies *PRINTOUT* as the file that receives the PRINT output.

## ***ENCODING Subcommand***

ENCODING specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is UTF8. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), UTF8, UTF16, UTF16BE (big endian), UTF16LE (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- If there is no OUTFILE subcommand, the ENCODING subcommand is ignored.

## ***TABLE Subcommand***

TABLE requests a table that shows how the variable information is formatted. NOTABLE, which suppresses the format table, is the default.

- TABLE must be specified before the slash that precedes the start of the variable specifications.

### ***Example***

```
PRINT TABLE /1 EMPLOYID DEPT SALARY /2 NAME.
EXECUTE.
```

- TABLE requests a summary table that describes the PRINT specifications. The table is included with the PRINT output.

# PRINT EJECT

```
PRINT EJECT [OUTFILE='file'] [ENCODING='encoding specification']
  [RECORDS={1}] [{NOTABLE}]
    {n}   {TABLE }

  /{1    } varlist [{col location [(format)]}] [varlist...]
    {rec #}      {(format list)      }
                {*                    }

  [{2    }...]
    {rec #}
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 16.0

- ENCODING subcommand added for Unicode support.

## Example

```
DO IF $CASENUM EQ 1.
PRINT EJECT / '   NAME ' 1 'DEPT' 25 'HIRED' 30 ' SALARY' 35.
END IF.
PRINT / NAME DEPT *
      MOHIRED(T30,F2) '/' YRHIRED *
      SALARY (T35,DOLLAR8) .
EXECUTE.
```

## Overview

PRINT EJECT displays specified information at the top of a new page of the output. Each time that it is executed, PRINT EJECT causes a page ejection. If not used in a DO IF-END IF structure, PRINT EJECT is executed for each case in the data, and each case is displayed on a separate page.

PRINT EJECT is designed to be used with the PRINT command to insert titles and column headings above the values that are displayed by PRINT. PRINT can also generate titles and headings, but PRINT cannot be used to control page ejections.

PRINT EJECT and PRINT can be used for writing simple reports.

## Options

The options that are available for PRINT EJECT are identical to the options that are available for PRINT:

- You can specify formats for the variables.
- You can specify string values within the variable specifications. With PRINT EJECT, the strings are usually used as titles or column headings and often include a specification for column location.



- You can use the `RECORDS` subcommand to display each case on more than one line.
- You can use the `OUTFILE` subcommand to direct the output to a specified file.
- You can use the `TABLE` subcommand to display a table that summarizes variable formats.
- You can use `ENCODING` to specify the file encoding. If there is no `OUTFILE` subcommand, the `ENCODING` subcommand is ignored.

For additional information, refer to `PRINT`.

### **Basic Specification**

The basic specification is a slash followed by a variable list and/or a list of string values that will be used as column headings or titles. The values for each variable or string are displayed on the top line of a new page in the output. `PRINT EJECT` is usually used within a `DO IF-END IF` structure to control the page ejections.

### **Operations**

- `PRINT EJECT` is a transformation and is not executed unless it is followed by a procedure or the `EXECUTE` command.
- If not used within a `DO IF-END IF` structure, `PRINT EJECT` is executed for each case in the data and displays the values for each case on a separate page.
- Values are displayed with a blank space between them. However, if a format is specified for a variable, the blank space for that variable's values is suppressed.
- Values are displayed in the output as the data are read. The `PRINT` output appears before the output from the first procedure.
- If more variables are specified than can be displayed in 132 columns or within the width that is specified on `SET WIDTH`, the program displays an error message. You must reduce the number of variables or split the output into several records.
- User-missing values are displayed exactly like valid values. System-missing values are represented by a period.

## **Examples**

### **Displaying Column Headings on the First Output Page Only**

```
DO IF $CASENUM EQ 1.
PRINT EJECT / '   NAME ' 1 'DEPT' 25 'HIRED' 30 '   SALARY' 35.
END IF.
PRINT / NAME DEPT *
      MOHIRED(T30,F2) '/' YRHIRED *
      SALARY (T35,DOLLAR8).
EXECUTE.
```

- `PRINT EJECT` specifies strings to be used as column headings and causes a page ejection. `DO IF-END IF` causes the `PRINT EJECT` command to be executed only once, when the system variable `$CASENUM` equals 1 (the value that is assigned to the first case in the file). Thus, column headings are displayed on the first page of the output only. The next example shows how to display column headings at the top of every page of the output.

- If a PRINT command were used in place of PRINT EJECT, the column headings would begin immediately after the command printback.

### ***Displaying Column Headings on Each Output Page***

```
DO IF MOD($CASENUM,50) = 1.  
PRINT EJECT FILE=OUT / ' NAME ' 1 'DEPT' 25 'HIRED' 30 ' SALARY' 35.  
END IF.  
PRINT FILE=OUT / NAME DEPT *  
          MOHIRED 30-31 '/' YRHIRED *  
          SALARY 35-42 (DOLLAR).  
EXECUTE.
```

- In this example, DO IF specifies that PRINT EJECT is executed if MOD (the remainder) of \$CASENUM divided by 50 equals 1 (see Arithmetic functions on p. 73 for information on the MOD function). Thus, column headings are displayed on a new page after every 50th case.
- If PRINT were used instead of PRINT EJECT, column headings would be displayed after every 50th case but would not appear at the top of a new page.
- Both PRINT EJECT and PRINT specify the same file for the output. If the FILE subcommands on PRINT EJECT and PRINT do not specify the same file, the column headings and the displayed values end up in different files.

# PRINT FORMATS

```
PRINT FORMATS varlist(format) [varlist...]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
PRINT FORMATS SALARY (DOLLAR8) / HOURLY (DOLLAR7.2)
/ RAISE BONUS (PCT2) .
```

## Overview

PRINT FORMATS changes variable print formats. Print formats are output formats and control the form in which values are displayed by a procedure or by the PRINT command.

PRINT FORMATS changes only print formats. To change write formats, use the WRITE FORMATS command. To change both the print and write formats with a single specification, use the FORMATS command. For information about assigning input formats during data definition, see DATA LIST. [For more information, see the topic Variable Types and Formats on p. 54.](#)

## Basic Specification

The basic specification is a variable list followed by the new format specification in parentheses. All specified variables receive the new format.

## Syntax Rules

- You can specify more than one variable or variable list, followed by a format in parentheses. Only one format can be specified after each variable list. For clarity, each set of specifications can be separated by a slash.
- You can use keyword TO to refer to consecutive variables in the active dataset.
- The specified width of a format must include enough positions to accommodate any punctuation characters, such as decimal points, commas, dollar signs, or date and time delimiters. (This situation differs from assigning an *input* format on DATA LIST, where the program automatically expands the input format to accommodate punctuation characters in output.)
- Custom currency formats (CCw, CCw.d) must first be defined on the SET command before they can be used on PRINT FORMATS.
- For string variables, you can only use PRINT FORMATS to switch between A and AHX formats. PRINT FORMATS cannot be used to change the length of string variables. To change the defined length of a string variable, use the ALTER TYPE command.

**Operations**

- Unlike most transformations, `PRINT FORMATS` takes effect as soon as it is encountered in the command sequence. Special attention should be paid to the position of `PRINT FORMATS` among commands.
- Variables that are not specified on `PRINT FORMATS` retain their current print formats in the active dataset. To see the current formats, use the `DISPLAY` command.
- The new print formats are changed only in the active dataset and are in effect for the duration of the session or until changed again with a `PRINT FORMATS` or `FORMATS` command. Print formats in the original data file (if this file exists) are not changed, unless the file is resaved with the `SAVE` or `XSAVE` command.
- New numeric variables that are created with transformation commands are assigned default print formats of `F8.2` (or the format that is specified on the `FORMAT` subcommand of `SET`). The `FORMATS` command can be used to change the new variable's print formats.
- New string variables that are created with transformation commands are assigned the format that is specified on the `STRING` command that declares the variable. `PRINT FORMATS` cannot be used to change the format of a new string variable.
- If a numeric data value exceeds its width specification, the program still attempts to display some value. First, the program rounds decimal values, then removes punctuation characters, and then tries scientific notation. Finally, if there is still not enough space, the program produces asterisks indicating that a value is present but cannot be displayed in the assigned width.

**Examples****Basic Example**

```
PRINT FORMATS SALARY (DOLLAR8) / HOURLY (DOLLAR7.2)
          / RAISE BONUS (PCT2) .
```

- The print format for `SALARY` is changed to `DOLLAR` with eight positions, including the dollar sign and comma when appropriate. The value 11550 is displayed as \$11,550. An eight-digit number requires a `DOLLAR11` format specification: eight characters for digits, two characters for commas, and one character for the dollar sign.
- The print format for `HOURLY` is changed to `DOLLAR` with seven positions, including the dollar sign, decimal point, and two decimal places. The number 115 is displayed as \$115.00. If `DOLLAR6.2` had been specified, the value 115 would be displayed as \$115.0. The program would truncate the last 0 because a width of 6 is not enough to display the full value.
- The print format for both `RAISE` and `BONUS` is changed to `PCT` with two positions: one position for the percentage and one position for the percent sign. The value 9 is displayed as 9%. Because the width allows for only two positions, the value 10 is displayed as 10.

**Changing Default Formats**

```
COMPUTE V3=V1 + V2 .
PRINT FORMATS V3 (F3.1) .
```

- COMPUTE creates the new numeric variable *V3*. By default, *V3* is assigned an F8.2 format (or the default format that is specified on SET).
- PRINT FORMATS changes the print format for *V3* to F3.1.

### **Working With Custom Currency Formats**

```
SET CCA='-/- .Df1 ..-' .  
PRINT FORMATS COST (CCA14.2) .
```

- SET defines a European currency format for the custom currency format type CCA.
- PRINT FORMATS assigns the print format CCA to variable *COST*. With the format defined for CCA on SET, the value 37419 is displayed as Dfl'37.419,00. See the SET command for more information about custom currency formats.

# PRINT SPACE

```
PRINT SPACE [OUTFILE='file'] [ENCODING='encoding specification']  
           [numeric expression]
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- ENCODING subcommand added for Unicode support.

## **Example**

```
PRINT SPACE.
```

## **Overview**

PRINT SPACE displays blank lines in the output and is generally used with a PRINT or WRITE command. Because PRINT SPACE displays a blank line each time that the command is executed, it is often used in a DO IF-END IF structure.

## **Basic Specification**

The basic specification is the command PRINT SPACE.

## **Syntax Rules**

- To display more than one blank line, specify a numeric expression after PRINT SPACE. The expression can be an integer or a complex expression.
- OUTFILE directs the output to a specified file. OUTFILE should be specified if an OUTFILE subcommand is specified on the PRINT or WRITE command that is used with PRINT SPACE. The OUTFILE subcommand on PRINT SPACE and PRINT or WRITE should specify the same file.
- ENCODING specifies the encoding for the output file. If there is no OUTFILE subcommand, the ENCODING subcommand is ignored.

## **Operations**

- If not used in a DO IF-END IF structure, PRINT SPACE is executed for each case in the data and displays a blank line for every case.

## Examples

### Inserting a Blank Line after the Output for Each Case

```
PRINT / NAME DEPT82 *
      MOHIRED(T30,F2) '/' YRHIRED *
      SALARY82 (T35,DOLLAR8) .
PRINT SPACE.
EXECUTE.
```

- Each time that it is executed, PRINT SPACE displays one blank line. Because PRINT SPACE is not used in a DO IF-END IF structure, PRINT SPACE is executed once for each case. In effect, the output is double-spaced.

### Using PRINT SPACE Inside a DO IF-END IF Structure

```
NUMERIC #LINE.
DO IF MOD(#LINE,5) = 0.
PRINT SPACE 2.
END IF.
COMPUTE #LINE=#LINE + 1.
PRINT / NAME DEPT *
      MOHIRED 30-31 '/' YRHIRED *
      SALARY 35-42(DOLLAR) .
EXECUTE.
```

- DO IF specifies that PRINT SPACE will be executed if MOD (the remainder) of #LINE divided by 5 equals 0. Because #LINE is incremented by 1 for each case, PRINT SPACE is executed once for every five cases. (See Arithmetic functions on p. 73 for information about the MOD function.)
- PRINT SPACE specifies two blank lines. Cases are displayed in groups of five with two blank lines between each group.

### Using an Expression to Specify the Number of Blank Lines

```
* Printing addresses on labels.

COMPUTE #LINES=0.           /*Initiate #LINES to 0
DATA LIST FILE=ADDRESS/RECORD 1-40 (A). /*Read a record
COMPUTE #LINES=#LINES+1.   /*Bump counter and print
WRITE OUTFILE=LABELS /RECORD.

DO IF RECORD EQ ' '.      /*Blank between addresses
+ PRINT SPACE OUTFILE=LABELS 8 - #LINES. /*Add extra blank #LINES
+ COMPUTE #LINES=0.
END IF.
EXECUTE.
```

- PRINT SPACE uses a complex expression for specifying the number of blank lines to display. The data contain a variable number of input records for each name and address, which must be printed in a fixed number of lines for mailing labels. The goal is to know when the last line for each address has been printed, how many lines have printed, and therefore how many blank records must be printed in order for the next address to fit on the next label. The example assumes that there is already one blank line between each address on input and that you want to print eight lines per label.

- The `DATA LIST` command defines the data. Each line of the address is contained in columns 1–40 of the data file and is assigned the variable name *RECORD*. For the blank line between each address, *RECORD* is blank.
- Variable *#LINES* is initialized to 0 as a scratch variable and is incremented for each record that is written. When the program encounters a blank line (`RECORD EQ ' '`), `PRINT SPACE` prints a number of blank lines that is equal to 8 minus the number already printed, and *#LINES* is then reset to 0.
- `OUTFILE` on `PRINT SPACE` specifies the same file that is specified by `OUTFILE` on `WRITE`.



# PROBIT

PROBIT is available in the Regression option.

```
PROBIT response-count varname OF observation-count varname
      WITH varlist [BY varname(min,max)]

[/MODEL={PROBIT**}]
      {LOGIT  }
      {BOTH   }

[/LOG={ {10** } }
      {2.718 }
      {value }
      {NONE  }

[/CRITERIA={ {OPTOL  } ({epsilon**0.8}) } [P({0.15**})] [STEPLIMIT({0.1**})]
      {CONVERGE} {n      }           {p      }           {n      }

      [ITERATE({max(50,3(p+1)**)})]
      {n      }           }

[/NATRES[=value]]

[/PRINT={ [CI**] [FREQ**] [RMP**] } [PARALL] [NONE] [ALL]]
      {DEFAULT** }

[/MISSING={ {EXCLUDE**} } ]
      {INCLUDE  } ]
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
PROBIT R OF N BY ROOT(1,2) WITH X
      /MODEL = BOTH.
```

## Overview

PROBIT can be used to estimate the effects of one or more independent variables on a dichotomous dependent variable (such as dead or alive, employed or unemployed, product purchased or not). The program is designed for dose-response analyses and related models, but PROBIT can also estimate logistic regression models.

### Options

**The Model.** You can request a probit or logit response model, or both, for the observed response proportions with the MODEL subcommand.

**Transform Predictors.** You can control the base of the log transformation applied to the predictors or request no log transformation with the LOG subcommand.

**Natural Response Rates.** You can instruct `PROBIT` to estimate the natural response rate (threshold) of the model or supply a known natural response rate to be used in the solution with the `NATRES` subcommand.

**Algorithm Control Parameters.** You can specify values of algorithm control parameters, such as the limit on iterations, using the `CRITERIA` subcommand.

**Statistics.** By default, `PROBIT` calculates frequencies, fiducial confidence intervals, and the relative median potency. It also produces a plot of the observed probits or logits against the values of a single independent variable. Optionally, you can use the `PRINT` subcommand to request a test of the parallelism of regression lines for different levels of the grouping variable or to suppress any or all of these statistics.

### ***Basic Specification***

- The basic specification is the response-count variable, keyword `OF`, the observation-count variable, keyword `WITH`, and at least one independent variable.
- `PROBIT` calculates maximum-likelihood estimates for the parameters of the default probit response model and automatically displays estimates of the regression coefficient and intercept terms, their standard errors, a covariance matrix of parameter estimates, and a Pearson chi-square goodness-of-fit test of the model.

### ***Subcommand Order***

- The variable specification must be first.
- Subcommands can be named in any order.

### ***Syntax Rules***

- The variables must include a response count, an observation count, and at least one predictor. A categorical grouping variable is optional.
- All subcommands are optional and each can appear only once.
- Generally, data should not be entered for individual observations. `PROBIT` expects predictor values, response counts, and the total number of observations as the input case.
- If the data are available only in a case-by-case form, use `AGGREGATE` first to compute the required response and observation counts.

### ***Operations***

- The transformed response variable is predicted as a linear function of other variables using the nonlinear-optimization method. Note that the previous releases used the iteratively weighted least-squares method, which has a different way of transforming the response variables. [For more information, see the topic MODEL Subcommand on p. 1607.](#)
- If individual cases are entered in the data, `PROBIT` skips the plot of transformed response proportions and predictor values.
- If individual cases are entered, the chi-square goodness-of-fit statistic and associated degrees of freedom are based on the individual cases. The case-based chi-square goodness-of-fit statistic generally differs from that calculated for the same data in aggregated form.

**Limitations**

- Only one prediction model can be tested on a single `PROBIT` command, although both probit and logit response models can be requested for that prediction.
- Confidence limits, the plot of transformed response proportions and predictor values, and computation of relative median potency are necessarily limited to single-predictor models.

**Examples**

```
PROBIT
  response OF nsubj BY site(1 3) WITH value
  /LOG
  /MODEL PROBIT
  /PRINT FREQ CI RMP PARALLEL
  /NATRES
  /CRITERIA P(.15) ITERATE(100) STEPLIMIT(.1) .
```

- The fits a probit model to *response* given *nsubj* observations by *site* and the predictor *value*.
- The empty `LOG` subcommand means that the natural logarithm will be used.
- The empty `NATRES` subcommand asks the procedure to estimate the natural response rate.
- The `PRINT` subcommand requests the table of observed and expected frequencies, fiducial confidence intervals, relative mean potencies, and parallelism tests.
- All other options are set to their default values.

**Example: Using data in a case-by-case form**

```
DATA LIST FREE / PREPARTN DOSE RESPONSE.
BEGIN DATA
1 1.5 0
...
4 20.0 1
END DATA.
COMPUTE SUBJECT = 1.
PROBIT RESPONSE OF SUBJECT BY PREPARTN(1,4) WITH DOSE.
```

- This dose-response model (Finney, 1971) illustrates a case-by-case analysis. A researcher tests four different preparations at varying doses and observes whether each subject responds. The data are individually recorded for each subject, with 1 indicating a response and 0 indicating no response. The number of observations is always 1 and is stored in variable *SUBJECT*.
- `PROBIT` warns that the data are in a case-by-case form and that the plot is therefore skipped.
- The goodness-of-fit test and associated degrees of freedom are based on individual cases, not dosage groups.
- `PROBIT` displays predicted and observed frequencies for all individual input cases.

**Example: Aggregating case-by-case data**

```
DATA LIST FREE/PREPARTN DOSE RESPONSE.
BEGIN DATA
  1.00      1.50      .00
  ...
  4.00      20.00     1.00
END DATA.
AGGREGATE OUTFILE=*
```

```

/BREAK=PREPARTN DOSE
/SUBJECTS=N(RESPONSE)
/NRESP=SUM(RESPONSE) .
PROBIT NRESP OF SUBJECTS BY PREPARTN(1,4) WITH DOSE .

```

- This example analyzes the same dose-response model as the previous example, but the data are first aggregated.
- AGGREGATE summarizes the data by cases representing all subjects who received the same preparation (*PREPARTN*) at the same dose (*DOSE*).
- The number of cases having a nonmissing response is recorded in the aggregated variable *SUBJECTS*.
- Because *RESPONSE* is coded 0 for no response and 1 for a response, the sum of the values gives the number of observations with a response.
- PROBIT requests a default analysis.
- The parameter estimates for this analysis are the same as those calculated for individual cases in the example above. The chi-square test, however, is based on the number of dosages.

## Variable Specification

The variable specification on PROBIT identifies the variables for response count, observation count, groups, and predictors. The variable specification is required.

- The variables must be specified first. The specification must include the response-count variable, followed by the keyword OF and then the observation-count variable.
- If the value of the response-count variable exceeds that of the observation-count variable, a procedure error occurs and PROBIT is not executed.
- At least one predictor (covariate) must be specified following the keyword WITH. The number of predictors is limited only by available workspace. All predictors must be continuous variables.
- You can specify a grouping variable (factor) after the keyword BY. Only one variable can be specified. It must be numeric and can contain only integer values. You must specify, in parentheses, a range indicating the minimum and maximum values for the grouping variable. Each integer value in the specified range defines a group.
- Cases with values for the grouping variable that are outside the specified range are excluded from the analysis.
- Keywords BY and WITH can appear in either order. However, both must follow the response-and-observation-count variables.

### Example

```
PROBIT R OF N WITH X .
```

- The number of observations having the measured response appears in variable *R*, and the total number of observations is in *N*. The predictor is *X*.

**Example**

```
PROBIT R OF N BY ROOT(1,2) WITH X.
PROBIT R OF N WITH X BY ROOT(1,2).
```

- Because keywords `BY` and `WITH` can be used in either order, these two commands are equivalent. Each command specifies  $X$  as a continuous predictor and  $ROOT$  as a categorical grouping variable.
- Groups are identified by the levels of variable  $ROOT$ , which may be 1 or 2.
- For each combination of predictor and grouping variables, the variable  $R$  contains the number of observations with the response of interest, and  $N$  contains the total number of observations.

**MODEL Subcommand**

`MODEL` specifies the form of the dichotomous-response model. Response models can be thought of as transformations ( $T$ ) of response rates, which are proportions or probabilities ( $p$ ). Note the difference in the transformations between the current version and the previous versions.

- A **probit** is the inverse of the cumulative standard normal distribution function. Thus, for any proportion, the probit transformation returns the value below which that proportion of standard normal deviates is found. For the probit response model, the program uses  $T(p) = \text{PROBIT}(p)$ . Hence:

$$T(0.025) = \text{PROBIT}(0.025) = -1.96$$

$$T(0.400) = \text{PROBIT}(0.400) = -0.25$$

$$T(0.500) = \text{PROBIT}(0.500) = 0.00$$

$$T(0.950) = \text{PROBIT}(0.950) = 1.64$$

- A **logit** is simply the natural log of the odds ratio,  $p/(1-p)$ . In the Probit procedure, the response function is given as  $T(p) = \log_e(p/(1-p))$ . Hence:

$$T(0.025) = \text{LOGIT}(0.025) = -3.66$$

$$T(0.400) = \text{LOGIT}(0.400) = -0.40$$

$$T(0.500) = \text{LOGIT}(0.500) = 0.00$$

$$T(0.950) = \text{LOGIT}(0.950) = 2.94$$

You can request one or both of the models on the `MODEL` subcommand. The default is `PROBIT` if the subcommand is not specified or is specified with no keyword.

**PROBIT**     *Probit response model.* This is the default.

**LOGIT**     *Logit response model.*

**BOTH**     *Both probit and logit response models.* `PROBIT` displays all the output for the logit model followed by the output for the probit model.

- If subgroups and multiple-predictor variables are defined, `PROBIT` estimates a separate intercept,  $a_j$ , for each subgroup and a regression coefficient,  $b_i$ , for each predictor.

## LOG Subcommand

LOG specifies the base of the logarithmic transformation of the predictor variables or suppresses the default log transformation.

- LOG applies to all predictors.
- To transform only selected predictors, use COMPUTE commands before the Probit procedure. Then specify NONE on the LOG subcommand.
- If LOG is omitted, a logarithm base of 10 is used.
- If LOG is used without a specification, the natural logarithm base  $e$  (2.718) is used.
- If you have a control group in your data and specify NONE on the LOG subcommand, the control group is included in the analysis. [For more information, see the topic NATRES Subcommand on p. 1609.](#)

You can specify one of the following on LOG:

<b>value</b>	<i>Logarithm base to be applied to all predictors.</i>
<b>NONE</b>	<i>No transformation of the predictors.</i>

### Example

```
PROBIT R OF N BY ROOT (1,2) WITH X
  /LOG = 2.
```

- LOG specifies a base-2 logarithmic transformation.

## CRITERIA Subcommand

Use CRITERIA to specify the values of control parameters for the PROBIT algorithm. You can specify any or all of the keywords below. Defaults remain in effect for parameters that are not changed.

<b>OPTOL(n)</b>	<i>Optimality tolerance.</i> Alias CONVERGE. If an iteration point is a feasible point and the next step will not produce a relative change in either the parameter vector or the log-likelihood function of more than the square root of $n$ , an optimal solution has been found. OPTOL can also be thought of as the number of significant digits in the log-likelihood function at the solution. For example, if OPTOL=10 <sup>-6</sup> , the log-likelihood function should have approximately six significant digits of accuracy. The default value is machine epsilon*0.8.
<b>ITERATE(n)</b>	<i>Iteration limit.</i> Specify the maximum number of iterations. The default is max(50, 3( $p$ + 1)), where $p$ is the number of parameters in the model.
<b>P(p)</b>	<i>Heterogeneity criterion probability.</i> Specify a cutoff value between 0 and 1 for the significance of the goodness-of-fit test. The cutoff value determines whether a heterogeneity factor is included in calculations of confidence levels for effective levels of a predictor. If the significance of chi-square is greater than the cutoff, the heterogeneity factor is not included. If you specify 0, this criterion is disabled; if you specify 1, a heterogeneity factor is automatically included. The default is 0.15.

- STEPLIMIT(n)**      *Step limit.* The PROBIT algorithm does not allow changes in the length of the parameter vector to exceed a factor of  $n$ . This limit prevents very early steps from going too far from good initial estimates. Specify any positive value. The default value is 0.1.
- CONVERGE(n)**      *Alias of OPTOL.*

## NATRES Subcommand

You can use NATRES either to supply a known natural response rate to be used in the solution or to instruct PROBIT to estimate the natural (or threshold) response rate of the model.

- To supply a known natural response rate as a constraint on the model solution, specify a value less than 1 on NATRES.
- To instruct PROBIT to estimate the natural response rate of the model, you can indicate a control group by giving a 0 value to any of the predictor variables. PROBIT displays the estimate of the natural response rate and the standard error and includes the estimate in the covariance/correlation matrix as NAT RESP.
- If no control group is indicated and NATRES is specified without a given value, PROBIT estimates the natural response rate from the entire data and informs you that no control group has been provided. The estimate of the natural response rate and the standard error are displayed and NAT RESP is included in the covariance/correlation matrix.
- If you have a control group in your data and specify NONE on the LOG subcommand, the control group is included in the analysis.

### Example

```
DATA LIST FREE / SOLUTION DOSE NOBSN NRESP.
BEGIN DATA
1 5 100 20
1 10 80 30
1 0 100 10
...
END DATA.

PROBIT NRESP OF NOBSN BY SOLUTION(1,4) WITH DOSE
/NATRES.
```

- This example reads four variables and requests a default analysis with an estimate of the natural response rate.
- The predictor variable, *DOSE*, has a value of 0 for the third case.
- The response count (10) and the observation count (100) for this case establish the initial estimate of the natural response rate.
- Because the default log transformation is performed, the control group is not included in the analysis.

### Example

```
DATA LIST FREE / SOLUTION DOSE NOBSN NRESP.
BEGIN DATA
1 5 100 20
1 10 80 30
```

```

1  0 100 10
...
END DATA.

PROBIT NRESP OF NOBSN BY SOLUTION(1,4) WITH DOSE
/NATRES = 0.10.

```

- This example reads four variables and requests an analysis in which the natural response rate is set to 0.10. The values of the control group are ignored.
- The control group is excluded from the analysis because the default log transformation is performed.

## PRINT Subcommand

Use PRINT to control the statistics calculated by PROBIT.

- PROBIT always displays the plot (for a single-predictor model) and the parameter estimates and covariances for the probit model.
- If PRINT is used, the requested statistics are calculated and displayed in addition to the parameter estimates and plot.
- If PRINT is not specified or is specified without any keyword, *FREQ*, *CI*, and *RMP* are calculated and displayed in addition to the parameter estimates and plot.

<b>DEFAULT</b>	<i>FREQ</i> , <i>CI</i> , and <i>RMP</i> . This is the default if PRINT is not specified or is specified by itself.
<b>FREQ</b>	<i>Frequencies</i> . Display a table of observed and predicted frequencies with their residual values. If observations are entered on a case-by-case basis, this listing can be quite lengthy.
<b>CI</b>	<i>Fiducial confidence intervals</i> . Print fiducial confidence intervals (Finney et al., 1971) for the levels of the predictor needed to produce each proportion of responses. PROBIT displays this default output for single-predictor models only. If a categorical grouping variable is specified, PROBIT produces a table of confidence intervals for each group. If the Pearson chi-square goodness-of-fit test is significant ( $p < 0.15$ by default), PROBIT uses a heterogeneity factor to calculate the limits.
<b>RMP</b>	<i>Relative median potency</i> . Display the relative median potency (RMP) of each pair of groups defined by the grouping variable. PROBIT displays this default output for single-predictor models only. For any pair of groups, the RMP is the ratio of the stimulus tolerances in those groups. <b>Stimulus tolerance</b> is the value of the predictor necessary to produce a 50% response rate. If the derived model for one predictor and two groups estimates that a predictor value of 21 produces a 50% response rate in the first group, and that a predictor value of 15 produces a 50% response rate in the second group, the relative median potency would be $21/15 = 1.40$ . In biological assay analyses, RMP measures the comparative strength of preparations.
<b>PARALL</b>	<i>Parallelism test</i> . Produce a test of the parallelism of regression lines for different levels of the grouping variable. This test displays a chi-square value and its associated probability. It requires an additional pass through the data and, thus, additional processing time.
<b>NONE</b>	<i>Display only the unconditional output</i> . This option can be used to override any other specification on the PRINT subcommand for PROBIT.
<b>ALL</b>	<i>All available output</i> . This is the same as requesting <i>FREQ</i> , <i>CI</i> , <i>RMP</i> , and <i>PARALL</i> .



## ***MISSING Subcommand***

PROBIT always deletes cases having a missing value for any variable. In the output, PROBIT indicates how many cases it rejected because of missing data. This information is displayed with the DATA Information that prints at the beginning of the output. You can use the MISSING subcommand to control the treatment of user-missing values.

- EXCLUDE**            *Delete cases with user-missing values.* This is the default. You can also make it explicit by using the keyword DEFAULT.
- INCLUDE**            *Include user-missing values.* PROBIT treats user-missing values as valid. Only cases with system-missing values are rejected.

## ***References***

Finney, D. J. 1971. *Probit analysis*. Cambridge: Cambridge University Press.

# PROCEDURE OUTPUT

```
PROCEDURE OUTPUT OUTFILE=file.
```

## **Example**

```
PROCEDURE OUTPUT OUTFILE=CELLDATA.
```

## **Overview**

PROCEDURE OUTPUT specifies the files to which CROSSTABS and SURVIVAL (included in the Advanced Statistics option) can write procedure output. PROCEDURE OUTPUT has no other applications.

### **Basic Specification**

The only specification is OUTFILE and the file specification. PROCEDURE OUTPUT must precede the command to which it applies.

### **Operations**

Commands with the WRITE subcommand or keyword write to the output file that is specified on the most recent PROCEDURE OUTPUT command. If only one output file has been specified, the output from the last such procedure overwrites all previous ones.

## **Examples**

### **Using PROCEDURE OUTPUT with CROSSTABS**

```
PROCEDURE OUTPUT OUTFILE=CELLDATA.  
CROSSTABS VARIABLES=FEAR SEX (1,2)  
  /TABLES=FEAR BY SEX  
  /WRITE=ALL.
```

- PROCEDURE OUTPUT precedes CROSSTABS and specifies *CELLDATA* as the file to receive the cell frequencies.
- The WRITE subcommand on CROSSTABS is required for writing cell frequencies to a procedure output file.

### **Using PROCEDURE OUTPUT with SURVIVAL**

```
PROCEDURE OUTPUT OUTFILE=SURVTBL.  
SURVIVAL TABLES=ONSSURV,RECSURV BY TREATMNT(1,3)  
  /STATUS = RECURSIT(1,9) FOR RECSURV  
  /STATUS = STATUS(3,4) FOR ONSSURV  
  /INTERVAL=THRU 50 BY 5 THRU 100 BY 10/PLOTS/COMPARE  
  /CALCULATE=CONDITIONAL PAIRWISE  
  /WRITE=TABLES.
```

- `PROCEDURE OUTPUT` precedes `SURVIVAL` and specifies *SURVTBL* as the file to receive the survival tables.
- The `WRITE` subcommand on `SURVIVAL` is required for writing survival tables to a procedure output file.

# PROXIMITIES

PROXIMITIES is available in the Statistics Base option.

```
PROXIMITIES varlist [/VIEW={CASE** }]
                    {VARIABLE}

[/STANDARDIZE={ {VARIABLE} } [ {NONE** }]]
                    {CASE } {Z }
                    {SD }
                    {RANGE }
                    {MAX }
                    {MEAN }
                    {RESCALE}

[/MEASURE={ {EUCLID** } } [ABSOLUTE] [REVERSE] [RESCALE]
            {SEUCLID }
            {COSINE }
            {CORRELATION }
            {BLOCK }
            {CHEBYCHEV }
            {POWER(p,r) }
            {MINKOWSKI(p) }
            {CHISQ }
            {PH2 }
            {RR[ (p[, np]) ] }
            {SM[ (p[, np]) ] }
            {JACCARD[ (p[, np]) ] }
            {DICE[ (p[, np]) ] }
            {SS1[ (p[, np]) ] }
            {RT[ (p[, np]) ] }
            {SS2[ (p[, np]) ] }
            {K1[ (p[, np]) ] }
            {SS3[ (p[, np]) ] }
            {K2[ (p[, np]) ] }
            {SS4[ (p[, np]) ] }
            {HAMANN[ (p[, np]) ] }
            {OCHIAI[ (p[, np]) ] }
            {SS5[ (p[, np]) ] }
            {PHI[ (p[, np]) ] }
            {LAMBDA[ (p[, np]) ] }
            {D[ (p[, np]) ] }
            {Y[ (p[, np]) ] }
            {Q[ (p[, np]) ] }
            {BEUCLID[ (p[, np]) ] }
            {SIZE[ (p[, np]) ] }
            {PATTERN[ (p[, np]) ] }
            {BSEUCLID[ (p[, np]) ] }
            {BSHAPE[ (p[, np]) ] }
            {DISPER[ (p[, np]) ] }
            {VARIANCE[ (p[, np]) ] }
            {BLWMN[ (p[, np]) ] }
            {NONE }

[/PRINT={ {PROXIMITIES**} } [/ID=varname]
        {NONE }

[/MISSING={EXCLUDE** } [INCLUDE]]

[/MATRIX=[IN({'savfile' | 'dataset'})] [OUT({'savfile' | 'dataset'})]]
          {* } {* }
```

\*\*Default if subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

**Example**

PROXIMITIES A B C.

**Overview**

PROXIMITIES computes a variety of measures of similarity, dissimilarity, or distance between pairs of cases or pairs of variables for moderate-sized datasets (see “Limitations” below). PROXIMITIES matrix output can be used as input to procedures ALSICAL, CLUSTER, and FACTOR.

**Options**

**Standardizing Data.** With the STANDARDIZE subcommand, you can use several different methods to standardize the values for each variable or for each case.

**Proximity Measures.** You can use the MEASURE subcommand to compute a variety of similarity, dissimilarity, and distance measures. (Similarity measures increase with greater similarity; dissimilarity and distance measures decrease.) MEASURE can compute measures for interval data, frequency-count data, and binary data. Only one measure can be requested in any one PROXIMITIES procedure. With the VIEW subcommand, you can control whether proximities are computed between variables or between cases.

**Output.** You can use the PRINT subcommand to display a computed matrix.

**Matrix Input and Output.** You can use the MATRIX subcommand to write a computed proximities matrix to IBM® SPSS® Statistics data files. This matrix can be used as input to procedures CLUSTER, ALSICAL, and FACTOR. You can also use MATRIX to read a similarity, dissimilarity, or distance matrix. This option lets you rescale or transform existing proximity matrices.

**Basic Specification**

The basic specification is a variable list, which obtains Euclidean distances between cases based on the values of each specified variable.

**Subcommand Order**

- The variable list must be first.
- Subcommands can be named in any order.

**Operations**

- PROXIMITIES ignores case weights when computing coefficients.

**Limitations**

- PROXIMITIES keeps the raw data for the current split-file group in memory. Storage requirements increase rapidly with the number of cases and the number of items (cases or variables) for which PROXIMITIES computes coefficients.

## Example

PROXIMITIES A B C.

- PROXIMITIES computes Euclidean distances between cases based on the values of variables *A*, *B*, and *C*.

## Variable Specification

- The variable list must be specified first.
- The variable list can be omitted when an input matrix data file is specified. A slash must then be specified before the first subcommand to indicate that the variable list is omitted.

## STANDARDIZE Subcommand

Use STANDARDIZE to standardize data values for either cases or variables before computing proximities. One of two options can be specified to control the direction of standardization:

**VARIABLE**            *Standardize the values for each variable.* This setting is the default.  
**CASE**                    *Standardize the values within each case.*

Several standardization methods are available. These methods allow you to equalize selected properties of the values. All methods can be used with either VARIABLE or CASE. Only one standardization method can be specified.

- If STANDARDIZE is omitted, proximities are computed by using the original values (keyword NONE).
- If STANDARDIZE is used without specifications, proximities are computed by using *Z* scores (keyword Z).
- STANDARDIZE cannot be used with binary measures.

**NONE**                    *Do not standardize.* Proximities are computed by using the original values. This setting is the default if STANDARDIZE is omitted.

**Z**                         *Standardize values to Z scores, with a mean of 0 and a standard deviation of 1.* PROXIMITIES subtracts the mean value for the variable or case from each value that is being standardized, and then PROXIMITIES divides by the standard deviation. If the standard deviation is 0, PROXIMITIES sets all values for the case or variable to 0. This setting is the default if STANDARDIZE is used without specifications.

**RANGE**                 *Standardize values to have a range of 1.* PROXIMITIES divides each value that is being standardized by the range of values for the variable or case. If the range is 0, PROXIMITIES leaves all values unchanged.

**RESCALE**              *Standardize values to have a range from 0 to 1.* From each value that is being standardized, PROXIMITIES subtracts the minimum value and then divides by the range for the variable or case. If a range is 0, PROXIMITIES sets all values for the case or variable to 0.50.

**MAX**                     *Standardize values to a maximum magnitude of 1.* PROXIMITIES divides each value that is being standardized by the maximum value for the variable or case. If the maximum of the values is 0, PROXIMITIES divides each value by the absolute magnitude of the smallest value and adds 1.

<b>MEAN</b>	<i>Standardize values to a mean of 1.</i> PROXIMITIES divides each value that is being standardized by the mean of the values for the variable or case. If the mean is 0, PROXIMITIES adds 1 to all values for the case or variable to produce a mean of 1.
<b>SD</b>	<i>Standardize values to unit standard deviation.</i> PROXIMITIES divides each value that is being standardized by the standard deviation of the values for the variable or case. PROXIMITIES does not change the values if their standard deviation is 0.

**Example**

```
PROXIMITIES A B C
  /STANDARDIZE=CASE RANGE.
```

- Within each case, values are standardized to have ranges of 1.

**VIEW Subcommand**

VIEW indicates whether proximities are computed between cases or between variables.

<b>CASE</b>	<i>Compute proximity values between cases.</i> This is the default.
<b>VARIABLE</b>	<i>Compute proximity values between variables.</i>

**MEASURE Subcommand**

MEASURE specifies the similarity, dissimilarity, or distance measure that PROXIMITIES computes. Three transformations are available:

<b>ABSOLUTE</b>	<i>Take the absolute values of the proximities.</i> Use ABSOLUTE when the sign of the values indicates the direction of the relationship (as with correlation coefficients) but only the magnitude of the relationship is of interest.
<b>REVERSE</b>	<i>Transform similarity values into dissimilarities, or vice versa.</i> Use this specification to reverse the ordering of the proximities by negating the values.
<b>RESCALE</b>	<i>Rescale the proximity values to a range of 0 to 1.</i> RESCALE standardizes the proximities by first subtracting the value of the smallest proximity and then dividing by the range. You would not usually use RESCALE with measures that are already standardized on meaningful scales, as are correlations, cosines, and many binary coefficients.

PROXIMITIES can compute any one of a number of measures between items. You can choose among measures for interval data, frequency-count data, or binary data. Available keywords for each type of measures are defined in the following sections.

- Only one measure can be specified. However, each measure can be specified with any of the transformations ABSOLUTE, REVERSE, or RESCALE. To apply a transformation to an existing matrix of proximity values without computing any measures, use keyword NONE (see Transforming Measures in Proximity Matrix on p. 1624).
- If more than one transformation is specified, PROXIMITIES handles them in the order listed above: ABSOLUTE, REVERSE, and then RESCALE (regardless of the order in which they are specified).
- Each entry in the resulting proximity matrix represents a pair of items. The items can be either cases or variables, whichever is specified on the VIEW subcommand.

- When the items are cases, the computation for each pair of cases involves pairs of values for the specified variables.
- When the items are variables, the computation for each pair of variables involves pairs of values for the variables across all cases.

**Example**

```
PROXIMITIES A B C
  /MEASURE=EUCLID REVERSE.
```

- MEASURE specifies a EUCLID measure and a REVERSE transformation.

**Measures for Interval Data**

To obtain proximities for interval data, use one of the following keywords on MEASURE:

<b>EUCLID</b>	<i>Euclidean distance.</i> The distance between two items, $x$ and $y$ , is the square root of the sum of the squared differences between the values for the items. This setting is the default. $EUCLID(x, y) = \sqrt{\sum_i (x_i - y_i)^2}$
<b>SEUCLID</b>	<i>Squared Euclidean distance.</i> The distance between two items is the sum of the squared differences between the values for the items. $SEUCLID(x, y) = \sum_i (x_i - y_i)^2$
<b>CORRELATION</b>	<i>Correlation between vectors of values.</i> This measure is a pattern-similarity measure. $CORRELATION(x, y) = \frac{\sum_i (Z_{xi} Z_{yi})}{N-1}$ where $Z_{xi}$ is the $Z$ -score (standardized) value of $x$ for the $i$ th case or variable, and $N$ is the number of cases or variables.
<b>COSINE</b>	<i>Cosine of vectors of values.</i> This measure is a pattern-similarity measure. $COSINE(x, y) = \frac{\sum_i (x_i y_i)}{\sqrt{(\sum_i x_i^2)(\sum_i y_i^2)}}$
<b>CHEBYCHEV</b>	<i>Chebychev distance metric.</i> The distance between two items is the maximum absolute difference between the values for the items. $CHEBYCHEV(x, y) = \max_i  x_i - y_i $
<b>BLOCK</b>	<i>City-block or Manhattan distance.</i> The distance between two items is the sum of the absolute differences between the values for the items. $BLOCK(x, y) = \sum_i  x_i - y_i $



- MINKOWSKI(p)** *Distance in an absolute Minkowski power metric.* The distance between two items is the  $p$ th root of the sum of the absolute differences to the  $p$ th power between the values for the items. Appropriate selection of the integer parameter  $p$  yields Euclidean and many other distance metrics.
- $$MINKOWSKI(x, y) = (\sum_i |x_i - y_i|^p)^{1/p}$$
- POWER(p,r)** *Distance in an absolute power metric.* The distance between two items is the  $r$ th root of the sum of the absolute differences to the  $p$ th power between the values for the items. Appropriate selection of the integer parameters  $p$  and  $r$  yields Euclidean, squared Euclidean, Minkowski, city-block, and many other distance metrics.
- $$POWER(x, y) = (\sum_i |x_i - y_i|^p)^{1/r}$$

### Measures for Frequency-Count Data

To obtain proximities for frequency-count data, use either of the following keywords on MEASURE:

- CHISQ** *Based on the chi-square test of equality for two sets of frequencies.* The magnitude of this dissimilarity measure depends on the total frequencies of the two cases or variables whose dissimilarity is computed. Expected values are from the model of independence of cases or variables  $x$  and  $y$ .
- $$CHISQ(x, y) = \sqrt{\frac{\sum_i (x_i - E(x_i))^2}{E(x_i)} + \frac{\sum_i (y_i - E(y_i))^2}{E(y_i)}}$$
- PH2** *Phi-square between sets of frequencies.* This measure is the CHISQ measure normalized by the square root of the combined frequency. Therefore, its value does not depend on the total frequencies of the two cases or variables whose dissimilarity is computed.
- $$PH2(x, y) = \sqrt{\frac{\frac{\sum_i (x_i - E(x_i))^2}{E(x_i)} + \frac{\sum_i (y_i - E(y_i))^2}{E(y_i)}}{N}}$$

### Measures for Binary Data

Different binary measures emphasize different aspects of the relationship between sets of binary values. However, all measures are specified in the same way. Each measure has two optional integer-valued parameters,  $p$  (present) and  $np$  (not present).

- If both parameters are specified, PROXIMITIES uses the value of the first parameter as an indicator that a characteristic is present, and PROXIMITIES uses the value of the second parameter as an indicator that a characteristic is absent. PROXIMITIES skips all other values.
- If only the first parameter is specified, PROXIMITIES uses that value to indicate presence and uses all other values to indicate absence.
- If no parameters are specified, PROXIMITIES assumes that 1 indicates presence and 0 indicates absence.

Using the indicators for presence and absence within each item (case or variable), PROXIMITIES constructs a  $2 \times 2$  contingency table for each pair of items and uses this table to compute a proximity measure for the pair.

	Item 2 characteristics	
	Present	Absent
Item 1 characteristics		
Present	<i>a</i>	<i>b</i>
Absent	<i>c</i>	<i>d</i>

PROXIMITIES computes all binary measures from the values of *a*, *b*, *c*, and *d*. These values are tallied across variables (when the items are cases) or cases (when the items are variables). For example, if variables *V*, *W*, *X*, *Y*, *Z* have values 0, 1, 1, 0, 1 for case 1 and have values 0, 1, 1, 0, 0 for case 2 (where 1 indicates presence and 0 indicates absence), the contingency table is as follows:

	Case 2 characteristics	
	Present	Absent
Case 1 characteristics		
Present	2	1
Absent	0	2

The contingency table indicates that both cases are present for two variables (*W* and *X*), both cases are absent for two variables (*V* and *Y*), and case 1 is present and case 2 is absent for one variable (*Z*). There are no variables for which case 1 is absent and case 2 is present.

The available binary measures include matching coefficients, conditional probabilities, predictability measures, and other measures.

**Matching Coefficients.** The following table shows a classification scheme for PROXIMITIES matching coefficients. In this scheme, *matches* are joint presences (value *a* in the contingency table) or joint absences (value *d*). *Nonmatches* are equal in number to value *b* plus value *c*. Matches and nonmatches may be weighted equally or not. The three coefficients JACCARD, DICE, and SS2 are related monotonically, as are SM, SS1, and RT. All coefficients in the table are similarity measures, and all coefficients except  $\kappa_1$  and SS3 range from 0 to 1.  $\kappa_1$  and SS3 have a minimum value of 0 and have no upper limit.

Table 193-1  
Binary matching coefficients in PROXIMITIES

	Joint absences excluded from numerator	Joint absences included in numerator
All matches included in denominator		
Equal weight for matches and nonmatches	RR	SM
Double weight for matches		SS1
Double weight for nonmatches		RT
Joint absences excluded from denominator		
Equal weight for matches and nonmatches	JACCARD	
Double weight for matches	DICE	
Double weight for nonmatches	SS2	

	Joint absences excluded from numerator	Joint absences included in numerator
All matches excluded from denominator		
Equal weight for matches and nonmatches	K1	SS3

**RR**[(p],[np])]

*Russell and Rao similarity measure.* This measure is the binary dot product.

$$RR(x, y) = \frac{a}{a+b+c+d}$$

**SM**[(p],[np])]

*Simple matching similarity measure.* This measure is the ratio of the number of matches to the total number of characteristics.

$$SM(x, y) = \frac{a+d}{a+b+c+d}$$

**JACCARD**[(p],[np])]

*Jaccard similarity measure.* This measure is also known as the *similarity ratio*.

$$JACCARD(x, y) = \frac{a}{a+b+c}$$

**DICE**[(p],[np])]

*Dice (or Czekanowski or Sorenson) similarity measure.*

$$DICE(x, y) = \frac{2a}{2a+b+c}$$

**SS1**[(p],[np])]

*Sokal and Sneath similarity measure 1.*

$$SS1(x, y) = \frac{2(a+d)}{2(a+d)+b+c+d}$$

**RT**[(p],[np])]

*Rogers and Tanimoto similarity measure.*

$$RT(x, y) = \frac{a+d}{a+d+2(b+c)}$$

**SS2**[(p],[np])]

*Sokal and Sneath similarity measure 2.*

$$SS2(x, y) = \frac{a}{a+2(b+c)}$$

**K1**[(p],[np])]

*Kulczynski similarity measure 1.* This measure has a minimum value of 0 and no upper limit. The measure is undefined when there are no nonmatches ( $b=0$  and  $c=0$ ).

$$K1(x, y) = \frac{a}{b+c}$$

**SS3**[(p],[np])]

*Sokal and Sneath similarity measure 3.* This measure has a minimum value of 0 and no upper limit. The measure is undefined when there are no nonmatches ( $b=0$  and  $c=0$ ).

$$SS3(x, y) = \frac{a+d}{b+c}$$

**Conditional Probabilities.** The following binary measures yield values that can be interpreted in terms of conditional probability. All three measures are similarity measures.

**K2[(p[,np])]**

*Kulczynski similarity measure 2.* This measure yields the average conditional probability that a characteristic is present in one item given that the characteristic is present in the other item. The measure is an average over both items that are acting as predictors. The measure has a range of 0 to 1.

$$K2(x, y) = \frac{a/(a+b)+a/(a+c)}{2}$$

**SS4[(p[,np])]**

*Sokal and Sneath similarity measure 4.* This measure yields the conditional probability that a characteristic of one item is in the same state (presence or absence) as the characteristic of the other item. The measure is an average over both items that are acting as predictors. The measure has a range of 0 to 1.

$$SS4(x, y) = \frac{a/(a+b)+a/(a+c)+d/(b+d)+d/(c+d)}{4}$$

**HAMANN[(p[,np])]**

*Hamann similarity measure.* This measure gives the probability that a characteristic has the same state in both items (present in both or absent from both) minus the probability that a characteristic has different states in the two items (present in one and absent from the other). HAMANN has a range of -1 to +1 and is monotonically related to SM, SS1, and RT.

$$HAMANN(x, y) = \frac{(a+d)-(b+c)}{a+b+c+d}$$

**Predictability Measures.** The following four binary measures assess the association between items as the predictability of one item given the other item. All four measures yield similarities.

**LAMBDA[(p[,np])]**

*Goodman and Kruskal's lambda (similarity).* This coefficient assesses the predictability of the state of a characteristic on one item (present or absent) given the state on the other item. Specifically, LAMBDA measures the proportional reduction in error, using one item to predict the other item when the directions of prediction are of equal importance. LAMBDA has a range of 0 to 1.

$$LAMBDA(x, y) = \frac{t_1 - t_2}{2(a+b+c+d) - t_2}$$

where

$$t_1 = \max(a, b) + \max(c, d) + \max(a, c) + \max(b, d)$$

$$t_2 = \max(a + c, b + d) + \max(a + d, c + d).$$

**D[(p[,np])]**

*Anderberg's D (similarity).* This coefficient assesses the predictability of the state of a characteristic on one item (present or absent) given the state on the other item. D measures the actual reduction in the error probability when one item is used to predict the other item. The range of D is 0 to 1.

$$D(x, y) = \frac{t_1 - t_2}{2(a+b+c+d)}$$

where

$$t_1 = \max(a, b) + \max(c, d) + \max(a, c) + \max(b, d)$$

$$t_2 = \max(a + c, b + d) + \max(a + d, c + d)$$

**Y[(p[,np])]** *Yule's Y coefficient of colligation (similarity)*. This measure is a function of the cross ratio for a 2×2 table and has a range of -1 to +1.

$$Y(x, y) = \frac{\sqrt{ad} - \sqrt{bc}}{\sqrt{ad} + \sqrt{bc}}$$

**Q[(p[,np])]** *Yule's Q (similarity)*. This measure is the 2×2 version of Goodman and Kruskal's ordinal measure *gamma*. Like Yule's Y, Q is a function of the cross ratio for a 2×2 table and has a range of -1 to +1.

**Other Binary Measures.** The remaining binary measures that are available in PROXIMITIES are either binary equivalents of association measures for continuous variables or measures of special properties of the relationship between items.

**OCHIAI[(p[,np])]** *Ochiai similarity measure*. This measure is the binary form of the cosine and has a range of 0 to 1.

$$OCHIAI(x, y) = \sqrt{\frac{a}{a+b} \frac{a}{a+c}}$$

**SS5[(p[,np])]** *Sokal and Sneath similarity measure 5*. The range is 0 to 1.

$$SS5(x, y) = \frac{ad}{\sqrt{(a+b)(a+c)(b+d)(c+d)}}$$

**PHI[(p[,np])]** *Fourfold point correlation (similarity)*. This measure is the binary form of the Pearson product-moment correlation coefficient.

$$PHI(x, y) = \frac{ad - bc}{\sqrt{(a+b)(a+c)(b+d)(c+d)}}$$

**BEUCLID[(p[,np])]** *Binary Euclidean distance*. This measure is a distance measure. Its minimum value is 0, and it has no upper limit.

$$BEUCLID(x, y) = \sqrt{b + c}$$

**BSEUCLID[(p[,np])]** *Binary squared Euclidean distance*. This measure is a distance measure. Its minimum value is 0, and it has no upper limit.

$$BSEUCLID(x, y) = b + c$$

**SIZE[(p[,np])]** *Size difference*. This measure is a dissimilarity measure with a minimum value of 0 and no upper limit.

$$SIZE(x, y) = \frac{(b-c)^2}{(a+b+c+d)^2}$$

**PATTERN[(p[,np])]** *Pattern difference*. This measure is a dissimilarity measure. The range is 0 to 1.

$$PATTERN(x, y) = \frac{bc}{(a+b+c+d)^2}$$

**BSHAPE[(p[,np])]** *Binary shape difference*. This dissimilarity measure has no upper limit or lower limit.

$$BSHAPE(x, y) = \frac{(a+b+c+d)(b+c) - (b-c)^2}{(a+b+c+d)^2}$$

**DISPER[(p[,np])]** *Dispersion similarity measure*. The range is -1 to +1.

$$DISPER(x, y) = \frac{ad - bc}{(a+b+c+d)^2}$$

**VARIANCE**[(p[,np])] *Variance dissimilarity measure.* This measure has a minimum value of 0 and no upper limit.

$$VARIANCE(x, y) = \frac{b+c}{4(a+b+c+d)}$$

**BLWMN**[(p[,np])] *Binary Lance-and-Williams nonmetric dissimilarity measure.* This measure is also known as the Bray-Curtis nonmetric coefficient. The range is 0 to 1.

$$BLWMN(x, y) = \frac{b+c}{2a+b+c}$$

**Example**

```
PROXIMITIES A B C
  /MEASURE=RR (1, 2) .
```

- MEASURE computes Russell and Rao coefficients from data in which 1 indicates the presence of a characteristic and 2 indicates the absence. Other values are ignored.

**Example**

```
PROXIMITIES A B C
  /MEASURE=SM (2) .
```

- MEASURE computes simple matching coefficients from data in which 2 indicates presence and all other values indicate absence.

**Transforming Measures in Proximity Matrix**

Use keyword **NONE** to apply the **ABSOLUTE**, **REVERSE**, and/or **RESCALE** transformations to an existing matrix of proximity values without computing any proximity measures.

**NONE** *Do not compute proximity measures.* Use **NONE** only if you have specified an existing proximity matrix on keyword **IN** on the **MATRIX** subcommand.

**PRINT Subcommand**

**PROXIMITIES** always prints the name of the measure that it computes and the number of cases. Use **PRINT** to control printing of the proximity matrix.

**PROXIMITIES** *Print the matrix of the proximities between items.* This setting is the default. The matrix may have been either read or computed. When the number of cases or variables is large, this specification produces a large volume of output and uses significant CPU time.

**NONE** *Do not print the matrix of proximities.*

**ID Subcommand**

By default, **PROXIMITIES** identifies cases by case number alone. Use **ID** to specify an identifying string variable for cases.

- Any string variable in the active dataset can be named as the identifier. PROXIMITIES uses the first eight characters of this variable to identify cases in the output.
- When used with the MATRIX IN subcommand, the variable that is specified on the ID subcommand identifies the labeling variable in the matrix file.

## **MISSING Subcommand**

MISSING controls the treatment of cases with missing values.

- PROXIMITIES deletes cases with missing values listwise. By default, PROXIMITIES excludes user-missing values from the analysis.

**EXCLUDE**        *Exclude cases with user-missing values.* This setting is the default.

**INCLUDE**        *Include cases with user-missing values.* Only cases with system-missing values are deleted.

## **MATRIX Subcommand**

MATRIX reads and writes matrix data files.

- Either IN or OUT and the matrix file in parentheses are required. When both IN and OUT are used on the same PROXIMITIES command, they can be specified on separate MATRIX subcommands or on the same subcommand.

**OUT ('savfile'|'dataset')**        *Write a matrix data file or dataset.* Specify either a filename, a previously declared dataset name, or an asterisk, enclosed in parentheses. Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. If you specify an asterisk (\*), the matrix data file replaces the active dataset. If you specify an asterisk or a dataset name, the file is not stored on disk unless you use SAVE or XSAVE.

**IN ('savfile'|'dataset')**        *Read a matrix data file or dataset.* Specify either a filename, dataset name created during the current session, or an asterisk enclosed in parentheses. An asterisk reads the matrix data from the active dataset. Filenames should be enclosed in quotes and are read from the working directory unless a path is included as part of the file specification.

When a matrix is produced by using the MATRIX OUT subcommand, the matrix corresponds to a unique dataset. All subsequent analyses that are performed on this matrix would match the corresponding analysis on the original data. However, if the data file is altered in any way, this matching process would no longer occur.

For example, if the original file is edited or rearranged it would no longer correspond to the initially produced matrix. You need to make sure that the data match the matrix whenever inferring the results from the matrix analysis. Specifically, when the cluster membership is saved into an active dataset in the CLUSTER procedure, the proximity matrix in the MATRIX IN statement must match the current active dataset.

### **Matrix Output**

- PROXIMITIES writes a variety of proximity matrices, each matrix with *ROWTYPE\_* values of PROX. PROXIMITIES neither reads nor writes additional statistics with its matrix materials. See [Format of the Matrix Data File on p. 1627](#) for a description of the file.
- The matrices that PROXIMITIES writes can be used by PROXIMITIES or other procedures. Procedures CLUSTER and ALSICAL can read a proximity matrix directly. Procedure FACTOR can read a correlation matrix that is written by PROXIMITIES, but RECODE must first be used to change the *ROWTYPE\_* value PROX to *ROWTYPE\_* value CORR. Also, the ID subcommand cannot be used on PROXIMITIES if the matrix will be used in FACTOR.
- If VIEW=VARIABLE, the variables in the matrix file will have the names and labels of the original variables.
- If VIEW=CASE (the default), the variables in the matrix file are named *VARI*, *VAR2*, ...*VARn*, where *n* is the sequential number of the variable in the new file. The numeric suffix *n* is consecutive and does not necessarily match the number of the actual case. If there are no split files, the case number appears in the variable label in the form *CASE m*. The numeric suffix *m* is the actual case number and may not be consecutive (for example, if cases were selected before PROXIMITIES was executed).
- If VIEW=CASE, a numeric variable *CASENO\_* is added to the matrix file. Values of *CASENO\_* are the case numbers in the original file.
- The new file preserves the names and values of any split-file variables that are in effect. When split-file processing is in effect, no labels are generated for variables in the new file. The actual case number is retained by the variable *ID*.
- Any documents that are contained in the active dataset are not transferred to the matrix file.

### **Matrix Input**

- PROXIMITIES can read a matrix file that is written by a previous PROXIMITIES procedure.
- Values for split-file variables should precede values for *ROWTYPE\_*, *CASENO\_* and the labeling variable (if present) should come after *ROWTYPE\_* and before *VARNAME\_*.
- If *CASENO\_* is of type string rather than numeric, it is considered unavailable and a warning is issued.
- If *CASENO\_* appears on a variable list, a syntax error results.
- PROXIMITIES ignores unrecognized *ROWTYPE\_* values. In addition, PROXIMITIES ignores variables that are present in the matrix file that are not specified (or used by default) on the PROXIMITIES variable list.
- The program reads variable names, variable and value labels, and print and write formats from the dictionary of the matrix data file.
- MATRIX=IN cannot be used unless an active dataset has already been defined. To read an existing matrix data file at the beginning of a session, use GET to retrieve the matrix file, and then specify IN(\*) on MATRIX.
- When you read a matrix that is created with MATRIX DATA, you should supply a value label for PROX of either *SIMILARITY* or *DISSIMILARITY* so that the matrix is correctly identified. If you do not supply a label, PROXIMITIES assumes *DISSIMILARITY*. See [Format of the Matrix Data File on p. 1627](#).



- The variable list on PROXIMITIES can be omitted when a matrix file is used as input. When the variable list is omitted, all variables in the matrix data file are used in the analysis. If a variable list is specified, the specified variables can be a subset of the variables in the matrix file.
- With a large number of variables, the matrix data file will wrap when displayed (as with LIST) and will be difficult to read. Nonetheless, the matrix values are accurate and can be used as matrix input.

### **Format of the Matrix Data File**

- The matrix data file includes three special variables created by the program: ROWTYPE\_, VARNAME\_, and CASENO\_. Variable ROWTYPE\_ is a short string variable with value PROX (for proximity measure). PROX is assigned value labels containing the distance measure that is used to create the matrix and either SIMILARITY or DISSIMILARITY as an identifier. Variable VARNAME\_ is a short string variable whose values are the names of the new variables. Variable CASENO\_ is a numeric variable with values equal to the original case numbers.
- The matrix file includes the string variable that is named on the ID subcommand. This variable is used to identify cases. Up to 20 characters can be displayed for the identifier variable; longer values are truncated. The identifier variable is present only when VIEW=CASE (the default) and when the ID subcommand is used.
- The remaining variables in the matrix file are the variables that are used to form the matrix.

### **Split Files**

- When split-file processing is in effect, the first variables in the matrix system file are the split variables, followed by ROWTYPE\_, the case-identifier variable (if VIEW=CASE and ID are used), VARNAME\_, and the variables that form the matrix.
- A full set of matrix materials is written for each split-file group that is defined by the split variables.
- A split variable cannot have the same name as any other variable that is written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split file must be in effect when that matrix is read by any procedure.

### **Example: Matrix Output to IBM SPSS Statistics External File**

```
PROXIMITIES V1 TO V20
/MATRIX=OUT(DISTOUT) .
```

- PROXIMITIES produces a default Euclidean distance matrix for cases by using variables V1 through V20 and saves the matrix in the IBM® SPSS® Statistics data file DISTOUT.
- The names of the variables on the matrix file will be VARI, VAR2, ...VARn.

**Example: Matrix Output to External File**

```
GET FILE='/data/crime.sav'.
PROXIMITIES MURDER TO MOTOR
  /ID=CITY
  /MEASURE=EUCLID
  /MATRIX=OUT(PROXMTX).
```

- PROXIMITIES reads data from the IBM® SPSS® Statistics data file *crime.sav* and writes one set of matrix materials to file *PROXMTX*.
- The active dataset is still *crime.sav*. Subsequent commands are executed on this file.

**Example: Matrix Output to Working File**

```
GET FILE='/data/crime.sav'.
PROXIMITIES MURDER TO MOTOR
  /ID=CITY
  /MEASURE=EUCLID
  /MATRIX=OUT(*).
LIST.
```

- PROXIMITIES writes the same matrix as in the example above. However, the matrix data file replaces the active dataset. The LIST command is executed on the matrix file, not on the *crime.sav* file.

**Example: Matrix Input from External File**

```
GET FILE PRSNNL.
FREQUENCIES VARIABLE=AGE.

PROXIMITIES CASE1 TO CASE8
  /ID=CITY
  /MATRIX=IN(PROXMTX).
```

- This example performs a frequencies analysis on file *PRSNNL* and then uses a different file that contains matrix data for PROXIMITIES.
- MATRIX=IN specifies the matrix data file *PROXMTX*. *PROXMTX* does not replace *PRSNNL* as the active dataset.

**Example: Matrix Input from Working File**

```
GET FILE PROXMTX.
PROXIMITIES CASE1 TO CASE8
  /ID=CITY
  /MATRIX=IN(*).
```

- This example assumes that you are starting a new session and want to read an existing matrix data file. GET retrieves the matrix file *PROXMTX*.
- MATRIX=IN specifies an asterisk because the matrix data file is the active dataset. If MATRIX=IN(PROXMTX) is specified, the program issues an error message.
- If the GET command is omitted, the program issues an error message.

### Example: Matrix Output to and Then Input from Working File

```
GET FILE='/data/crime.sav'.
PROXIMITIES MURDER TO MOTOR
  /ID=CITY
  /MATRIX=OUT(*).
PROXIMITIES
  /MATRIX=IN(*)
  /STANDARDIZE.
```

- GET retrieves the IBM® SPSS® Statistics data file *crime.sav*.
- The first PROXIMITIES command specifies variables for the analysis and reads data from file *crime.sav*. ID specifies *CITY* as the case identifier. MATRIX writes the resulting matrix to the active dataset.
- The second PROXIMITIES command uses the matrix file that is written by the first PROXIMITIES command as input. The asterisk indicates that the matrix file is the active dataset. The variable list is omitted, indicating that all variables in the matrix are to be used.
- The slash preceding the MATRIX subcommand on the second PROXIMITIES command is required. Without the slash, PROXIMITIES attempts to interpret MATRIX as a variable name rather than as a subcommand.

### Example: Q-factor Analysis

In this example, PROXIMITIES and FACTOR are used for a *Q*-factor analysis, in which factors account for variance shared among observations rather than among variables. Procedure FACTOR does not perform *Q*-factor analysis without some preliminary transformation such as what is provided by PROXIMITIES. Because the number of cases exceeds the number of variables, the model is not of full rank, and FACTOR will print a warning. This result is a common occurrence when case-by-case matrices from PROXIMITIES are used as input to FACTOR.

```
* Recoding a PROXIMITIES matrix for procedure FACTOR.

GET FILE='/data/crime.sav'.
PROXIMITIES MURDER TO MOTOR
  /MEASURE=CORR
  /MATRIX=OUT('/data/tempfile.sav').
GET FILE='/data/tempfile.sav' /DROP=ID.
RECODE ROWTYPE_ ('PROX' = 'CORR').
FACTOR MATRIX IN(COR=*) .
```

- The MATRIX subcommand on PROXIMITIES writes the correlation matrix to the active dataset. Because the matrix materials will be used in procedure FACTOR, the ID subcommand is not specified.
- RECODE recodes *ROWTYPE\_* value PROX to CORR so that procedure FACTOR can read the matrix.
- When FACTOR reads matrix materials, it reads all variables in the file. The MATRIX subcommand on FACTOR indicates that the matrix is a correlation matrix and that data are in the active dataset.

**References**

Anderberg, M. R. 1973. *Cluster analysis for applications*. New York: Academic Press.

Romesburg, H. C. 1984. *Cluster analysis for researchers*. Belmont, Calif.: Lifetime Learning Publications.

# PROXSCAL

PROXSCAL is available in the Categories option.

PROXSCAL varlist

```
[/TABLE = {rowid BY columnid [BY sourceid]]
           {sourceid
           }

[/SHAPE = [{LOWER**}]
           {UPPER
           }
           {BOTH
           }

[/INITIAL = [{SIMPLEX**
             }]]
            {TORGERSON
            }
            {RANDOM[({1})]
            }
            {n
            }
            {(['file'|'dataset']) [varlist] }

[/WEIGHTS = varlist]

[/CONDITION = [{MATRIX**
               }]]
              {UNCONDITIONAL
              }

[/TRANSFORMATION = [{RATIO**
                    }]]
                  {INTERVAL
                  }
                  {ORDINAL[({UNTIE
                             })]
                  }
                  {KEEPTIES
                  }
                  {SPLINE [DEGREE = {2}] [INKNOT = {1}]
                  }
                  {n
                  }

[/PROXIMITIES = [{DISSIMILARITIES**}]
                 {SIMILARITIES
                 }

[/MODEL = [{IDENTITY**
            }]]
           {WEIGHTED
           }
           {GENERALIZED
           }
           {REDUCED[({2})]
           }
           {n
           }

[/RESTRICTIONS = {COORDINATES('file'|'dataset') [{ALL
                                                    }]]
                  {varlist
                  }
                  {VARIABLES('file'|'dataset') [{ALL
                                                  }][({INTERVAL
                                                  })]
                  }
                  {varlist
                  }
                  {NOMINAL
                  }
                  {ORDINAL[({UNTIE
                             })]
                  }
                  {KEEPTIES
                  }
                  {SPLINE[DEGREE={2}] [INKNOT={1}]
                  }
                  {n
                  }
                  {n
                  }

[/ACCELERATION = NONE]

[/CRITERIA = [DIMENSIONS({2**
                        })]
             {min[,max]}
             [MAXITER({100**})]
             {n
             }
             [DIFFSTRESS({0.0001**})]
             {value
             }
             [MINSTRESS({0.0001**}) ]
             {value
             }

[/PRINT = [NONE] [INPUT] [RANDOM] [HISTORY] [STRESS**] [DECOMPOSITION]
           [COMMON**] [DISTANCES] [WEIGHTS**] [INDIVIDUAL]
           [TRANSFORMATIONS] [VARIABLES**] [CORRELATIONS**]

[/PLOT = [NONE] [STRESS] [COMMON**] [WEIGHTS**] [CORRELATIONS**]
         [INDIVIDUAL({varlist})]
         {ALL
         }
         [TRANSFORMATIONS({varlist}) [({varlist})[...]] ]
         {ALL
         }
         {ALL
         }]
```

```

[RESIDUALS({varlist}) [({varlist})[...]] ]
  {ALL      }      {ALL      }
[VARIABLES({varlist})]
  {ALL      }

[/OUTFILE = [COMMON('file'|'dataset')] [WEIGHTS('file'|'dataset')] [DISTANCES('file'|'dataset')]
  [TRANSFORMATIONS('file'|'dataset')] [VARIABLES('file'|'dataset')] ]

[/MATRIX = IN('file'|'dataset')].

```

\*\* Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Overview

PROXSCAL performs multidimensional scaling of proximity data to find a least-squares representation of the objects in a low-dimensional space. Individual differences models are allowed for multiple sources. A majorization algorithm guarantees monotone convergence for optionally transformed metric and nonmetric data under a variety of models and constraints.

### Options

**Data Input.** You can read one or more square matrices of proximities that can either be symmetrical or asymmetrical. Alternatively, you can provide specifications with the TABLE subcommand for matrices with proximities in a stacked format. You can read proximity matrices created by PROXIMITIES and CLUSTER with the MATRIX subcommand. Additionally, you can read weights, initial configurations, fixed coordinates, and independent variables.

**Methodological Assumptions.** You can specify transformations considering all sources (unconditional) or separate transformations for each source (matrix-conditional) on the CONDITION subcommand. You can treat proximities as nonmetric (ordinal) or as metric (numerical or splines) using the TRANSFORMATION subcommand. Ordinal transformations can treat tied observations as tied (discrete) and untied (continuous). You can specify whether your proximities are similarities or dissimilarities on the PROXIMITIES subcommand.

**Model Selection.** You can specify multidimensional scaling models by selecting a combination of PROXSCAL subcommands, keywords, and criteria. The subcommand MODEL offers, besides the identity model, three individual differences models. You can specify other selections on the CRITERIA subcommand.

**Constraints.** You can specify fixed coordinates or independent variables to restrict the configuration(s) on the RESTRICTIONS subcommand. You can specify transformations (numerical, nominal, ordinal, and splines) for the independent variables on the same subcommand.

**Output.** You can produce output that includes the original and transformed proximities, history of iterations, common and individual configurations, individual space weights, distances, and decomposition of the stress. Plots can be produced of common and individual configurations, individual space weights, transformations, and residuals.

**Basic Specification**

The basic specification is `PROXSCAL` followed by a variable list. By default, `PROXSCAL` produces a two-dimensional metric Euclidean multidimensional scaling solution (identity model). Input is expected to contain one or more square matrices with proximities that are dissimilarities. The ratio transformation of the proximities is matrix-conditional. The analysis uses a simplex start as an initial configuration. By default, output includes fit and stress values, the coordinates of the common space, and a chart of the common space configuration.

**Syntax Rules**

- The number of dimensions (both minimum and maximum) may not exceed the number of proximities minus one.
- Dimensionality reduction is omitted if combined with multiple random starts.
- If there is only one source, then the model is always assumed to be identity.

**Limitations**

- `PROXSCAL` needs at least three objects, which means that at least three variables must be specified in the variable list. In the case of the `TABLE` subcommand, the minimum value for `rowid` and `columnid` must be at least three.
- `PROXSCAL` recognizes data weights created by the `WEIGHT` command but only in combination with the `TABLE` subcommand.
- Split-file has no implications for `PROXSCAL`.

**Variable List Subcommand**

The variable list identifies the columns in the proximity matrix or matrices that `PROXSCAL` reads. Each variable identifies one column of the proximity matrix, with each case in the active dataset representing one row, unless specified otherwise with the `TABLE` subcommand. In this case, the variable list identifies whole matrices or sources.

- Only numeric variables can be specified.
- The total number of cases must be divisible by the number of variables. This is not applicable when the `TABLE` subcommand is used.
- `PROXSCAL` reads data row by row; the columns are represented by the variables on the variable list. The order of the variables on the list is crucial.

**Example**

```
DATA LIST
  /object01 object02 object03 object04.

BEGIN DATA
  0 2 6 3
  2 0 5 4
  6 5 0 1
  3 4 1 0
END DATA.
```

```
PROXSCAL VARIABLES=object01 TO object04.
```

- This example specifies an analysis on a 4×4 proximity matrix.
- The total number of cases must be divisible by 4.

## TABLE Subcommand

The `TABLE` subcommand specifies the row identifier `rowid` and the column identifier `columnid`. Using `TABLE`, the proximities of separate sources are given in separate variables on the `PROXSCAL` variable list.

In the same manner, sources are identified by `sourceid`. In combination with `rowid` and `columnid`, the proximities are stacked in one single variable, containing the proximities of all sources, where sources are distinguished by the values of `sourceid`.

Using `sourceid` as the only variable on the `TABLE` subcommand indicates the use of stacked matrices, where individual stacked matrices are recognized by different values of `sourceid`.

- `Rowid`, `columnid`, and `sourceid` should not be specified on the variable list.
- When specifying both upper- and lower-triangular parts of the matrix, the `SHAPE` subcommand will determine the handling of the data.
- If a cell's value is specified multiple times, the final specification is used.
- `Rowid`, `columnid`, and `sourceid` must appear in that order.
- Omitting `sourceid` causes `PROXSCAL` to use the sources specified on the `PROXSCAL` variable list. Each variable is assumed to contain the proximities of one source.
- Specifying multiple sources on the `PROXSCAL` variable list in conjunction with specifying `rowid`, `columnid`, and `sourceid` is not possible and causes `PROXSCAL` to ignore `sourceid`.

<b>rowid</b>	<i>Row identifying variable.</i> The values of this variable specify the row object of a proximity. The values must be integers between 1 and the number of objects, inclusive.
<b>columnid</b>	<i>Column identifying variable.</i> The values specify the column object of a proximity. The values must be integers between 1 and the number of objects, inclusive.
<b>sourceid</b>	<i>Source identifying variable.</i> The values specify the source number and must be integers between 1 and the number of sources, inclusive. The value labels of this variable are used to identify sources on other subcommands. These value labels must comply with variable name conventions. Omitting a value label causes <code>PROXSCAL</code> to use the default label <code>SRC_n</code> , where <code>n</code> is the number of the source.

### Example

```
DATA LIST
  /r_id c_id men women.
```

```
BEGIN DATA
2 1 1.08 1.14
3 1 0.68 1.12
3 2 0.95 0.75
4 1 0.96 0.32
```



```

4 2 0.76 0.98
4 3 0.47 0.69
. . . . .
. . . . .
13 10 0.55 0.86
13 11 0.61 0.97
13 12 0.46 0.83
END DATA.

```

```

PROXSCAL men women
  /TABLE=r_id BY c_id
  /PLOT = INDIVIDUAL (women).

```

- PROXSCAL reads two proximity matrices (*men* and *women*), where the row objects are specified by *r\_id* and the column objects by *c\_id*.
- A chart of the individual space for *women* is plotted.

This is one way to proceed. Another way is to add the proximities of the additional source below the proximities of the first source and specify *sourceid* on the TABLE subcommand, containing values distinguishing the first and the additional source (see the next example).

### Example

```

DATA LIST
  /r_id c_id s_id prox.

```

```

BEGIN DATA
2 1 1 1.08
3 1 1 0.68
3 2 1 0.95
4 1 1 0.96
4 2 1 0.76
4 3 1 0.47
. . . . .
. . . . .
13 10 1 0.55
13 11 1 0.61
13 12 1 0.46
2 1 2 1.14
3 1 2 1.12
3 2 2 0.75
4 1 2 0.32
4 2 2 0.98
4 3 2 0.69
. . . . .
. . . . .
13 10 2 0.86
13 11 2 0.97
13 12 2 0.83
END DATA.

```

```

VALUE LABELS s_id 1 'men' 2 'women'.

```

```

PROXSCAL prox
  /TABLE=r_id BY c_id BY s_id
  /PLOT = INDIVIDUAL (women).

```

- PROXSCAL reads two proximity matrices. The row objects are identified by *r\_id* and the column objects, by *c\_id*. The proximity matrices are gathered in one variable, *source01*, where each source is distinguished by a value of the source identifying variable *s\_id*.
- A chart of the individual space for *women* is plotted.

### Example

```
DATA LIST LIST
  /obj_1 obj_2 obj_3 obj_4 s_id

BEGIN DATA
0 0 0 0 1
1 0 0 0 1
2 3 0 0 1
4 5 6 0 1
7 0 0 0 2
0 0 0 0 2
8 9 0 0 2
12 11 12 0 2
END DATA.

VALUE LABELS s_id 1 'women' 2 'men'.

PROXSCAL obj_1 obj_2 obj_3 obj_4
  /TABLE = s_id
  /MODEL = WEIGHTED
  /PLOT = INDIVIDUAL (women).
```

- PROXSCAL reads two proximity matrices. The objects are given on the PROXSCAL variable list. Each source is distinguished by a value of the source identifying variable *s\_id*, which is also used for labeling.
- A chart of the individual space for *women* is plotted.

## SHAPE Subcommand

The SHAPE subcommand specifies the structure of the proximity matrix.

- |              |  |
|--------------|--|
| <b>LOWER</b> | <i>Lower-triangular data matrix.</i> For a lower-triangular matrix, PROXSCAL expects a square matrix of proximities of which the lower-triangular elements are used under the assumption that the full matrix is symmetric. The diagonal is ignored but must be included.              |
| <b>UPPER</b> | <i>Upper-triangular data matrix.</i> For an upper-triangular matrix, PROXSCAL expects a square matrix of proximities of which the upper-triangular elements are used under the assumption that the full matrix is symmetric. The diagonal is ignored but must be included.             |
| <b>BOTH</b>  | <i>Full data matrix.</i> The values in the corresponding cells in the upper and lower triangles may be different. PROXSCAL reads the complete square matrix and, after obtaining symmetry, continues with the lower-triangular elements. The diagonal is ignored but must be included. |

- System or other missing values on the (virtual) diagonal are ignored.

**Example**

```
PROXSCAL object01 TO object07
  /SHAPE=UPPER.
```

- PROXSCAL reads square matrices of seven columns per matrix of which the upper-triangular parts are used in computations.
- Although specified, the diagonal and lower-triangular part of the matrix are not used.

**INITIAL Subcommand**

INITIAL defines the initial or starting configuration of the common space for the analysis. When a reduction in dimensionality is specified on the CRITERIA subcommand, a derivation of coordinates in the higher dimensionality is used as a starting configuration in the lower dimensionality.

- You can specify one of the three keywords listed below.
- You can specify a variable list containing the initial configuration.

**SIMPLEX**

*Simplex start.* This specification is the default. PROXSCAL starts by placing the objects in the configuration all at the same distance of each other and taking one iteration to improve this high-dimensional configuration, followed by a dimension-reduction operation to obtain the user-provided maximum dimensionality specified in the CRITERIA subcommand with the keyword DIMENSIONS.

**TORGERSON**

*Torgerson start.* A classical scaling solution is used as initial configuration.

**RANDOM**

*(Multiple) random start.* You can specify the number of random starts ( $n$ ).  $n$  is any positive integer. The random sequence can be controlled by the RANDOM SEED command and not by a subcommand within the PROXSCAL command. Each analysis starts with a different random configuration. In the output, all  $n$  final stress values are reported, as well as the initial seeds of each analysis (for reproduction purposes), followed by the full output of the analysis with the lowest stress value. The default number of random starts is 1. Reduction of dimensionality—that is, using a maximum dimensionality that is larger than the minimum dimensionality—is not allowed within this option and the minimum dimensionality is used, if reduction is specified anyway.

Instead of these keywords, a parenthesized IBM® SPSS® Statistics data file can be specified containing the coordinates of the initial configuration. If the variable list is omitted, the first MAXDIM variables are automatically selected, where MAXDIM is the maximum number of dimensions requested for the analysis on the CRITERIA subcommand. Only nonmissing values are allowed as initial coordinates.

**Example**

```
PROXSCAL object01 TO object17
  /INITIAL=RANDOM(100).
```

- This example performs 100 analyses each, starting with different random configurations. The results of the analysis with the lowest final stress are displayed in the output.

## **WEIGHTS Subcommand**

The `WEIGHTS` subcommand specifies non-negative weights on the proximities included in the active dataset.

- The number and order of the variables in the variable list is important. The first variable on the `WEIGHTS` variable list corresponds to the first variable on the `PROXSCAL` variable list. This is repeated for all variables on the variable lists. Every proximity has its own weight. The number of variables on the `WEIGHTS` subcommand must therefore be equal to the number of variables on the `PROXSCAL` variable list.
- Negative weights are not allowed. If specified, a warning will be issued and the procedure will abort.

### **Example**

```
DATA LIST FILE='cola.dat' FREE
  /object01 TO object14 weight01 TO weight14.
```

```
PROXSCAL object01 TO object14
  /WEIGHTS=weight01 TO weight14.
```

- In this example, the `VARIABLES` subcommand indicates that there are 14 columns per matrix of which the weights can be found in *weight01* to *weight14*.
- *weight01* contains the weights for *object01*, etc.

## **CONDITION Subcommand**

`CONDITION` specifies how transformations among sources are compared. The `TRANSFORMATION` subcommand specifies the type of transformation.

<b>MATRIX</b>	<i>Matrix conditional.</i> Only the proximities within each source are compared with each other. This is the default.
<b>UNCONDITIONAL</b>	<i>Unconditional.</i> This specification is appropriate when the proximities in all sources can be compared with each other and result in a single transformation of all sources simultaneously.

- Note that if there is only one source, then `MATRIX` and `UNCONDITIONAL` give the same results.

### **Example**

```
PROXSCAL object01 TO object15
  /CONDITION=UNCONDITIONAL
  /TRANSFORMATION=ORDINAL (UNTIE) .
```

- In this example, the proximities are ordinally transformed, where tied proximities are allowed to be untied. The transformations are performed simultaneously over all possible sources.

## **TRANSFORMATION Subcommand**

TRANSFORMATION offers four different options for optimal transformation of the original proximities. The resulting values are called transformed proximities. The distances between the objects in the configuration should match these transformed proximities as closely as possible.

<b>RATIO</b>	<i>No transformation.</i> Omitting the entire subcommand is equivalent to using this keyword. In both cases, the transformed proximities are proportional to the original proximities. This “transformation” is only allowed for positive dissimilarities. In all other cases, a warning is issued and the transformation is set to INTERVAL.
<b>INTERVAL</b>	<i>Numerical transformation.</i> In this case, the transformed proximities are proportional to the original proximities, including free estimation of the intercept. The inclusion of the intercept assures that all transformed proximities are positive.
<b>ORDINAL</b>	<i>Ordinal transformation.</i> The transformed proximities have the same order as the original proximities. In parentheses, the approach to tied proximities can be specified. Keeping tied proximities tied, also known as secondary approach to ties, is default. Specification may be implicit, ORDINAL, or explicit, ORDINAL (KEEPTIES). Allowing tied proximities to be untied, also known as the primary approach to ties, is specified as ORDINAL (UNTIE).
<b>SPLINE</b>	<i>Monotone spline transformation.</i> The transformed proximities are a smooth nondecreasing piecewise polynomial transformation of the original proximities of the chosen degree. The pieces are specified by the number and placement of the interior knots.

### **SPLINE Keyword**

SPLINE has the following keywords:

<b>DEGREE</b>	<i>The degree of the polynomial.</i> If DEGREE is not specified, the degree is assumed to be 2. The range of DEGREE is between 1 and 3 (inclusive).
<b>INKNOT</b>	<i>The number of interior knots.</i> If INKNOT is not specified, the number of interior knots is assumed to be 1. The range of INKNOT is between 1 and the number of different proximities.

#### **Example**

```
PROXSCAL object01 TO object05
  /TRANSFORMATION=ORDINAL (UNTIE) .
```

- In this example, the proximities are ordinally transformed, where tied proximities are allowed to be untied.
- The default conditionality (MATRIX) implies that the transformation is performed for each source separately.

## **PROXIMITIES Subcommand**

The PROXIMITIES subcommand specifies the type of proximities used in the analysis. The term proximity is used for either similarity or dissimilarity data.

<b>DISSIMILARITIES</b>	<i>Dissimilarity data.</i> This specification is the default when PROXIMITIES is not specified. Small dissimilarities correspond to small distances, and large dissimilarities correspond to large distances.
<b>SIMILARITIES</b>	<i>Similarity data.</i> Small similarities correspond to large distances and large similarities correspond to small distances.

### **Example**

```
PROXSCAL object01 TO object12
  /PROXIMITIES=SIMILARITIES.
```

- In this example, PROXSCAL expects the proximities to be similarities.

## **MODEL Subcommand**

MODEL defines the scaling model for the analysis if more than one source is present. IDENTITY is the default model. The three other models are individual differences models.

<b>IDENTITY</b>	<i>Identity model.</i> All sources have the same configuration. This is the default model, and it is not an individual differences model.
<b>WEIGHTED</b>	<i>Weighted Euclidean model.</i> This model is an individual differences model and equivalent to the INDSCAL model in the ALSCAL procedure. Each source has an individual space, in which every dimension of the common space is weighted differentially.
<b>GENERALIZED</b>	<i>Generalized Euclidean model.</i> This model is equivalent to the GEMSCAL model in the ALSCAL procedure. Each source has an individual space that is equal to a rotation of the common space, followed by a differential weighting of the dimensions.
<b>REDUCED</b>	<i>Reduced rank model.</i> This model is similar to GENERALIZED, but the rank of the individual space is equal to $n$ . This number is always smaller than the maximum number of dimensions and equal to or greater than 1. The default is 2.

- If IDENTITY is specified for only one source, this subcommand is silently ignored.
- If an individual differences model is specified for only one source, a warning is issued, and the model is set to IDENTITY.

### **Example**

```
PROXSCAL object01 TO object07
  /MODEL=WEIGHTED.
```

- A weighted Euclidean model is fitted, but only when the number of cases in the active dataset is a multiple of 7, starting from 14 (14, 21, 28, and so on). Otherwise, there is only one source, and the model is set to IDENTITY.

## **RESTRICTIONS Subcommand**

PROXSCAL provides two types of restrictions for the user to choose from. The first type fixes (some) coordinates in the configuration. The second type specifies that the common space is a weighted sum of independent variables.

<b>COORDINATES</b>	<i>Fixed coordinates.</i> A parenthesized IBM® SPSS® Statistics data filename must be specified containing the fixed coordinates for the common space. A variable list may be given, if some specific variables need to be selected from the external file. If the variable list is omitted, the procedure automatically selects the first MAXDIM variables in the external file, where MAXDIM is the maximum number of dimensions requested for the analysis on the CRITERIA subcommand. A missing value indicates that a coordinate on a dimension is free. The coordinates of objects with nonmissing values are kept fixed during the analysis. The number of cases for each variable must be equal to the number of objects.
<b>VARIABLES</b>	<i>Independent variables.</i> The common space is restricted to be a linear combination of the independent variables in the variable list. A parenthesized SPSS Statistics data file must be specified containing the independent variables. If the variable list is omitted, the procedure automatically selects all variables in the external file. Instead of the variable list, the user may specify the keyword FIRST( <i>n</i> ), where <i>n</i> is a positive integer, to select the first <i>n</i> variables in the external file. The number of cases for each variable must be equal to the number of objects. After the variable selection specification, we may provide a list of keywords (in number equal to the number of the independent variables) indicating the transformations for the independent variables.

## **VARIABLES Keyword**

The following keywords may be specified:

<b>INTERVAL</b>	<i>Numerical transformation.</i> In this case, the transformed values of a variable are proportional to the original values of the variable, including free estimation of the intercept.
<b>NOMINAL</b>	<i>Nominal transformation.</i> The values are treated as unordered. The same values will obtain the same transformed values.
<b>ORDINAL</b>	<i>Ordinal transformation.</i> The values of the transformed variable have the same order as the values of the original variable. In parenthesis, the approach to tied values can be specified. Keeping tied values tied, also known as secondary approach to ties, is default. Specification may be implicit, ORDINAL, or explicit, ORDINAL(KEEPTIES). Allowing tied values to be untied, also known as the primary approach to ties, is specified as ORDINAL(UNTIE).
<b>SPLINE</b>	<i>Monotone spline transformation.</i> The transformed values of the variable are a smooth nondecreasing piecewise polynomial transformation of the original values of the chosen degree. The pieces are specified by the number and placement of the interior knots.

**SPLINE Keyword**

SPLINE has the following keywords:

<b>DEGREE</b>	<i>The degree of the polynomial.</i> If DEGREE is not specified, the degree is assumed to be 2. The range of DEGREE is between 1 and 3 (inclusive).
<b>INKNOT</b>	<i>The number of interior knots.</i> If INKNOT is not specified, the number of interior knots is assumed to be 1. The range of INKNOT is between 0 and the number of different values of the variable.

**Example**

```
PROXSCAL aunt TO uncle
  /RESTRICTIONS=VARIABLES(ivars.sav) degree generation gender
  (ORDINAL ORDINAL NOMINAL) .
```

- In this example, there are three independent variables specified: *degree*, *generation*, and *gender*.
- The variables are specified in the data file *ivars.sav*.
- On both *degree* and *generation*, ordinal transformations are allowed. By default, tied values in ordinal variables are kept tied. *Gender* is allowed to be nominally transformed.

**ACCELERATION Subcommand**

By default, a fast majorization method is used to minimize stress.

**NONE**                    *The standard majorization update.* This turns off the fast method.

- If the subcommand RESTRICTION is used with fixed coordinates or independent variables, ACCELERATION=NONE is in effect.
- If an individual differences model is specified on the MODEL subcommand, ACCELERATION=NONE is in effect.

**Example**

```
PROXSCAL VARIABLES=object01 TO object12
  /ACCELERATION=NONE .
```

- Here, relaxed updates are switched off through the specification of the keyword NONE after ACCELERATION.



## CRITERIA Subcommand

Use `CRITERIA` to set the dimensionality and criteria for terminating the algorithm, or minimization process. You can specify one or more of the following keywords:

- DIMENSIONS** *Minimum and maximum number of dimensions.* By default, PROXSCAL computes a solution in two dimensions ( $\text{min}=2$  and  $\text{max}=2$ ). The minimum and maximum number of dimensions can be any integers inclusively between 1 and the number of objects minus 1, as long as the minimum is less than or equal to the maximum. PROXSCAL starts computing a solution in the largest dimensionality and reduces the dimensionality in steps, until the lowest dimensionality is reached. Specifying a single value represents both minimum and maximum number of dimensions, thus `DIMENSIONS (4)` is equivalent to `DIMENSIONS (4, 4)`.
- MAXITER** *Maximum number of iterations.* By default,  $n=100$ , specifying the maximum number of iterations that is performed while one of the convergence criterion below (`CONVERGENCE` and `STRESSMIN`) is not yet reached. Decreasing this number might give less accurate results but will take less time.  $N$  must have a positive integer value.
- DIFFSTRESS** *Convergence criterion.* PROXSCAL minimizes the goodness-of-fit index normalized raw stress. By default, PROXSCAL stops iterating when the difference in consecutive stress values is less than 0.0001 ( $n=0.0001$ ). To obtain a more precise solution, you can specify a smaller value. The value specified must lie between 0.0 and 1.0, inclusively.
- MINSTRESS** *Minimum stress value.* By default, PROXSCAL stops iterating when the stress value itself is small, that is, less than 0.0001 ( $n=0.0001$ ). To obtain an even more precise solution, you can specify a smaller value. The value specified must lie between 0.0 and 1.0, inclusively.

### Example

```
PROXSCAL VARIABLES=object01 TO object24
  /CRITERIA=DIMENSIONS (2,4) MAXITER (200) DIFFSTRESS (0.00001) .
```

- The maximum number of dimensions equals 4 and the minimum number of dimensions equals 2. PROXSCAL computes a four-, three-, and two-dimensional solution, respectively.
- The maximum number of iteration is raised to 200.
- The convergence criterion is sharpened to 0.00001.

## PRINT Subcommand

`PRINT` specifies the optional output. By default, PROXSCAL displays the stress and fit values for each analysis, the coordinates of the common space, and, with appropriate specification on corresponding subcommands, the individual space weights and transformed independent variables, corresponding regression weights, and correlations.

- Omitting the `PRINT` subcommand or specifying `PRINT` without keywords is equivalent to specifying `COMMON`, `WEIGHTS`, and `VARIABLES`.
- If a keyword(s) is specified, only the output for that particular keyword(s) is displayed.
- In the case of duplicate or contradicting keyword specification, the last keyword applies.

- Inapplicable keywords are silently ignored. That is, specifying a keyword for which no output is available (for example, specifying `INDIVIDUAL` with only one source) will silently ignore this keyword.

<b>NONE</b>	<i>No output.</i> Display only the normalized raw stress and corresponding fit values.
<b>INPUT</b>	<i>Input data.</i> The display includes the original proximities, and, if present, the data weights, the initial configuration, and the fixed coordinates or the independent variables.
<b>RANDOM</b>	<i>Multiple random starts.</i> Displays the random number seed and stress value of each random start.
<b>HISTORY</b>	<i>History of iterations.</i> Displays the history of iterations of the main algorithm.
<b>STRESS</b>	<i>Stress measures.</i> Displays different stress values. The table contains values for normalized raw stress, Stress-I, Stress-II, S-Stress, dispersion accounted for (D.A.F.), and Tucker's coefficient of congruence. This is specified by default.
<b>DECOMPOSITION</b>	<i>Decomposition of stress.</i> Displays an object and source decomposition of stress, including row and column totals.
<b>COMMON</b>	<i>Common space.</i> Displays the coordinates of the common space. This is specified by default.
<b>DISTANCES</b>	<i>Distances.</i> Displays the distances between the objects in the configuration.
<b>WEIGHTS</b>	<i>Individual space weights.</i> Displays the individual space weights, only if one of the individual differences models is specified on the <code>MODEL</code> subcommand. Depending on the model, the space weights are decomposed in rotation weights and dimension weights, which are also displayed. This is specified by default.
<b>INDIVIDUAL</b>	<i>Individual spaces.</i> The coordinates of the individual spaces are displayed, only if one of the individual differences models is specified on the <code>MODEL</code> subcommand.
<b>TRANSFORMATION</b>	<i>Transformed proximities.</i> Displays the transformed proximities between the objects in the configuration.
<b>VARIABLES</b>	<i>Independent variables.</i> If <code>VARIABLES</code> was specified on the <code>RESTRICTIONS</code> subcommand, this keyword triggers the display of the transformed independent variables and the corresponding regression weights. This is specified by default.
<b>CORRELATIONS</b>	<i>Correlations.</i> The correlations between the independent variables and the dimensions of the common space are displayed. This is specified by default.

### Example

```
PROXSCAL VARIABLES=source01 TO source02
  /TABLE=row_id BY col_id
  /MODEL=WEIGHTED
  /PRINT=HISTORY COMMON STRESS.
```

- Here, a weighted Euclidean model is specified with two sources.
- The output consists of the history of iterations of the main algorithm, the coordinates of the common space, the individual space weights, and several measures of fit.

## ***PLOT Subcommand***

PLOT controls the display of plots. By default, PROXSCAL produces a scatterplot of object coordinates of the common space, the individual space weights, and the correlations between the independent variables (that is, equivalent to specifying COMMON, WEIGHTS, and CORRELATIONS).

- Specifying a keyword overrides the default output and only output is generated for that keyword.
- Duplicate keywords are silently ignored.
- In case of contradicting keywords, only the last keyword is considered.
- Inapplicable keywords (for example, stress with equal minimum and maximum number of dimensions on the CRITERIA subcommand) are silently ignored.
- Multiple variable lists are allowed for TRANSFORMATIONS and RESIDUALS. For each variable list, a separate plot will be displayed.

<b>NONE</b>	<i>No plots.</i> PROXSCAL does not produce any plots.
<b>STRESS</b>	<i>Stress plot.</i> A plot is produced of stress versus dimensions. This plot is only produced if the maximum number of dimensions is larger than the minimum number of dimensions.
<b>COMMON</b>	<i>Common space.</i> A scatterplot matrix of coordinates of the common space is displayed.
<b>WEIGHTS</b>	<i>Individual space weights.</i> A scatterplot is produced of the individual space weights. This is only possible if one of the individual differences models is specified on the MODEL subcommand. For the weighted Euclidean model, the weights are printed in plots with one dimension on each axis. For the generalized Euclidean model, one plot is produced per dimension, indicating both rotation and weighting of that dimension. The reduced rank model produces the same plot as the generalized Euclidean model does but reduces the number of dimensions for the individual spaces.
<b>INDIVIDUAL</b>	<i>Individual spaces.</i> For each source specified on the variable list, the coordinates of the individual spaces are displayed in scatterplot matrices. This is only possible if one of the individual differences models is specified on the MODEL subcommand.
<b>TRANSFORMATIONS</b>	<i>Transformation plots.</i> Plots are produced of the original proximities versus the transformed proximities. On the variable list, the sources can be specified of which the plot is to be produced.
<b>RESIDUALS</b>	<i>Residuals plots.</i> The transformed proximities versus the distances are plotted. On the variable list, the sources can be specified of which the plot is to be produced.
<b>VARIABLES</b>	<i>Independent variables.</i> Transformation plots are produced for the independent variables specified on the variable list.
<b>CORRELATIONS</b>	<i>Correlations.</i> A plot of correlations between the independent variables and the dimensions of the common space is displayed.

### ***Example***

```
PROXSCAL VARIABLES=source01 TO source02
  /TABLE=row_id BY col_id
  /MODEL=WEIGHTED
  /CRITERIA=DIMENSIONS(3)
  /PLOT=COMMON INDIVIDUAL(source02).
```

- Here, the syntax specifies a weighted Euclidean model with two sources in three dimensions.
- `COMMON` produces a scatterplot matrix defined by dimensions 1, 2, and 3.
- For the individual spaces, a scatterplot matrix with 3 dimensions is only produced for the individual space of *source02*.

## ***OUTFILE Subcommand***

`OUTFILE` saves coordinates of the common space, individual space weights, distances, transformed proximities, and transformed independent variables to an external IBM® SPSS® Statistics data file or previously declared dataset. The only specification required is a name for the output file.

<b>COMMON</b>	<i>Common space coordinates.</i> The coordinates of the common space are written to an external SPSS Statistics data file or previously declared dataset. The columns (variables) represent the dimensions <i>DIM_1</i> , <i>DIM_2</i> , ..., <i>DIM_n</i> of the common space. The number of cases (rows) in the SPSS Statistics data file equals the number of objects.
<b>WEIGHTS</b>	<i>Individual space weights.</i> The individual space weights are written to an external SPSS Statistics data file or previously declared dataset. The columns represent the dimensions <i>DIM_1</i> , <i>DIM_2</i> , ..., <i>DIM_n</i> of the space weights. The number of cases depends on the individual differences model specified on the <code>MODEL</code> subcommand. The weighted Euclidean model uses diagonal weight matrices. Only the diagonals are written to file and the number of cases is equal to the number of dimensions. The generalized Euclidean model uses full-rank nonsingular weight matrices. The matrices are written to the data file row by row. The reduced rank model writes matrices to the data file in the same way as the generalized Euclidean model does but does not write the reduced part.
<b>DISTANCES</b>	<i>Distances.</i> The matrices containing the distances for each source are stacked beneath each other and written to an external SPSS Statistics data file or previously declared dataset. The number of variables in the data file are equal to the number of objects ( <i>OBJ_1</i> , <i>OBJ_2</i> , ..., <i>OBJ_n</i> ) and the number of cases in the data file are equal to the number of objects times the number of sources.
<b>TRANSFORMATION</b>	<i>Transformed proximities.</i> The matrices containing the transformed proximities for each source are stacked beneath each other and written to an external SPSS Statistics data file or previously declared dataset. The number of variables in the file are equal to the number of objects ( <i>OBJ_1</i> , <i>OBJ_2</i> , ..., <i>OBJ_n</i> ) and the number of cases in the data file are equal to the number of objects times the number of sources.
<b>VARIABLES</b>	<i>Independent variables.</i> The transformed independent variables are written to an external SPSS Statistics data file or previously declared dataset. The variables are written to the columns ( <i>VAR_1</i> , <i>VAR_2</i> , ..., <i>VAR_n</i> ). The number of variables in the data file are equal to the number of independent variables and the number of cases are equal to the number of objects.

### ***Example***

```
PROXSCAL VARIABLES=source01 TO source04
/TABLE=row_id BY col_id
/OUTFILE=COMMON('/data/start.sav').
```

- Here, the coordinates of the common space are written to the SPSS Statistics data file *start.sav*.
- Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files.

## **MATRIX Subcommand**

MATRIX reads matrix data files. It can read a matrix written by either PROXIMITIES or CLUSTER.

- The specification on MATRIX is the keyword IN and the matrix file in parentheses.
- Generally, data read by PROXSCAL are already in matrix form, whether in square format, or in stacked format using the TABLE subcommand.
- The proximity matrices PROXSCAL reads have ROWTYPE\_ values of PROX.
- Using MATRIX=IN, PROXSCAL will ignore variables specified on the main variable list. All numerical variables from the matrix data file are processed.
- PROXSCAL ignores variables specified in the WEIGHTS subcommand in combination with the use of MATRIX=IN.
- With MATRIX=IN, only a source identifying variable can be specified on the TABLE subcommand. The sources are created as a result of a split file action.

**IN('file'|'dataset')**      *Read a matrix data file.* Specify a quoted file specification or previously declared dataset name, enclosed in parentheses. Data read through the MATRIX subcommand does not replace the active dataset.

### **Example**

```
GET FILE = '/data/proxmtx.SAV'.
```

```
PROXSCAL  
  /MATRIX=IN('/data/matrix.sav').
```

- MATRIX=IN specifies an external matrix data file called *matrix.sav*, of which all numerical variables are used for the current analysis.

# QUICK CLUSTER

QUICK CLUSTER is available in the Statistics Base option.

```
QUICK CLUSTER {varlist}
               {ALL      }

[/MISSING={ {LISTWISE**} } [INCLUDE] ]
           {PAIRWISE   }
           {DEFAULT    }

[/FILE='savfile' | 'dataset']

[/INITIAL=(value list)]

[/CRITERIA=[CLUSTER({2**})] [NOINITIAL] [MXITER({10**})] [CONVERGE({0**})]]
           {n      }
           {n      }
           {n      }

[/METHOD={ {KMEANS [ (NOUPDATE) ] **} }
           {KMEANS (UPDATE) }
           {CLASSIFY      }

[/PRINT=[INITIAL**] [CLUSTER] [ID(varname)] [DISTANCE] [ANOVA] [NONE]]

[/OUTFILE='savfile' | 'dataset']

[/SAVE=[CLUSTER( (varname) )] [DISTANCE( (varname) )]]
```

\*\*Default if subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
QUICK CLUSTER V1 TO V4
  /CRITERIA=CLUSTER(4)
  /SAVE=CLUSTER(GROUP).
```

## Overview

When the desired number of clusters is known, QUICK CLUSTER groups cases efficiently into clusters. It is not as flexible as CLUSTER, but it uses considerably less processing time and memory, especially when the number of cases is large.

### Options

**Algorithm Specifications.** You can specify the number of clusters to form with the CRITERIA subcommand. You can also use CRITERIA to control initial cluster selection and the criteria for iterating the clustering algorithm. With the METHOD subcommand, you can specify how to update cluster centers, and you can request classification only when working with very large data files.

**Initial Cluster Centers.** By default, QUICK CLUSTER chooses the initial cluster centers. Alternatively, you can provide initial centers on the INITIAL subcommand. You can also read initial cluster centers from IBM® SPSS® Statistics data files using the FILE subcommand.

**Optional Output.** With the `PRINT` subcommand, you can display the cluster membership of each case and the distance of each case from its cluster center. You can also display the distances between the final cluster centers and a univariate analysis of variance between clusters for each clustering variable.

**Saving Results.** You can write the final cluster centers to a data file using the `OUTFILE` subcommand. In addition, you can save the cluster membership of each case and the distance from each case to its classification cluster center as new variables in the active dataset using the `SAVE` subcommand.

### ***Basic Specification***

The basic specification is a list of variables. By default, `QUICK CLUSTER` produces two clusters. The two cases that are farthest apart based on the values of the clustering variables are selected as initial cluster centers and the rest of the cases are assigned to the nearer center. The new cluster centers are calculated as the means of all cases in each cluster, and if neither the minimum change nor the maximum iteration criterion is met, all cases are assigned to the new cluster centers again. When one of the criteria is met, iteration stops, the final cluster centers are updated, and the distance of each case is computed.

### ***Subcommand Order***

- The variable list must be specified first.
- Subcommands can be named in any order.

### ***Operations***

The procedure generally involves four steps:

- First, initial cluster centers are selected, either by choosing one case for each cluster requested or by using the specified values.
- Second, each case is assigned to the nearest cluster center, and the mean of each cluster is calculated to obtain the new cluster centers.
- Third, the maximum change between the new cluster centers and the initial cluster centers is computed. If the maximum change is not less than the minimum change value and the maximum iteration number is not reached, the second step is repeated and the cluster centers are updated. The process stops when either the minimum change or maximum iteration criterion is met. The resulting clustering centers are used as classification centers in the last step.
- In the last step, all cases are assigned to the nearest classification center. The final cluster centers are updated and the distance for each case is computed.

When the number of cases is large, directly clustering all cases may be impractical. As an alternative, you can cluster a sample of cases and then use the cluster solution for the sample to classify the entire group. This can be done in two phases:

- The first phase obtains a cluster solution for the sample. This involves all four steps of the `QUICK CLUSTER` algorithm. `OUTFILE` then saves the final cluster centers to a data file.

- The second phase requires only one pass through the data. First, the `FILE` subcommand specifies the file containing the final cluster centers from the first analysis. These final cluster centers are used as the initial cluster centers for the second analysis. `CLASSIFY` is specified on the `METHOD` subcommand to skip the second and third steps of the clustering algorithm, and cases are classified using the initial cluster centers. When all cases are assigned, the cluster centers are updated and the distance of each case is computed. This phase can be repeated until final cluster centers are stable.

## Example

```
QUICK CLUSTER
  zlnlong zlntoll zlnequi zlnCARD zlnwire zmultlin zvoice zpager zinterne
  zcallid zcallwai zforward zconfer zebill
  /MISSING=PAIRWISE
  /CRITERIA= CLUSTER(3) MXITER(20) CONVERGE(0)
  /METHOD=KMEANS(NOUPDATE)
  /PRINT INITIAL ANOVA DISTAN.
```

- The procedure clusters cases based upon their values for variables *zlnlong* through *zebill*.
- The `MISSING` subcommand specifies that a case is assigned to clusters based upon all clustering variables for which the case has nonmissing values.
- The `CRITERIA` subcommand specifies that three clusters will be formed, and that 20 iterations are allowed for updating cluster centers.
- The `PRINT` subcommand specifies that the initial cluster centers, distances between final cluster centers, and an ANOVA table of descriptive *F* tests for the clustering variables should be displayed.
- All other options are set to their default values.

## Variable List

The variable list identifies the clustering variables.

- The variable list is required and must be the first specification on `QUICK CLUSTER`.
- You can use keyword `ALL` to refer to all user-defined variables in the active dataset.
- `QUICK CLUSTER` uses squared Euclidean distances, which equally weight all clustering variables. If the variables are measured in units that are not comparable, the procedure will give more weight to variables with large variances. Therefore, you should standardize variables measured on different scales using procedure `DESCRIPTIVES` before performing `QUICK CLUSTER`.



## CRITERIA Subcommand

CRITERIA specifies the number of clusters to form and controls options for the clustering algorithm. You can use any or all of the keywords below.

- The NOINITIAL option followed by the remaining steps of the default QUICK CLUSTER algorithm makes QUICK CLUSTER equivalent to MacQueen's  $n$ -means clustering method.

<b>CLUSTER(n)</b>	<i>Number of clusters.</i> QUICK CLUSTER assigns cases to $n$ clusters. The default is 2.
<b>NOINITIAL</b>	<i>No initial cluster center selection.</i> By default, initial cluster centers are formed by choosing one case (with valid data for the clustering variables) for each cluster requested. The initial selection requires a pass through the data to ensure that the centers are well separated from one another. If NOINITIAL is specified, QUICK CLUSTER selects the first $n$ cases without missing values as initial cluster centers.
<b>MXITER(n)</b>	<i>Maximum number of iterations for updating cluster centers.</i> The default is 10. Iteration stops when the maximum number of iterations has been reached. MXITER is ignored when METHOD=CLASSIFY.
<b>CONVERGE(n)</b>	<i>Convergence criterion controlling minimum change in cluster centers.</i> The default value for $n$ is 0. The minimum change value equals the convergence value ( $n$ ) times the minimum distance between initial centers. Iteration stops when the largest change of any cluster center is less than or equal to the minimum change value. CONVERGE is ignored when METHOD=CLASSIFY.

## METHOD Subcommand

By default, QUICK CLUSTER recalculates cluster centers after assigning all the cases and repeats the process until one of the criteria is met. You can use the METHOD subcommand to recalculate cluster centers after each case is assigned or to suppress recalculation until after classification is complete. When METHOD=KMEANS is specified, QUICK CLUSTER displays the iteration history table.

<b>KMEANS(NOUPDATE)</b>	<i>Recalculate cluster centers after all cases are assigned for each iteration.</i> This is the default.
<b>KMEANS(UPDATE)</b>	<i>Recalculate a cluster center each time a case is assigned.</i> QUICK CLUSTER calculates the mean of cases currently in the cluster and uses this new cluster center in subsequent case assignment.
<b>CLASSIFY</b>	<i>Do not recalculate cluster centers.</i> QUICK CLUSTER uses the initial cluster centers for classification and computes the final cluster centers as the means of all the cases assigned to the same cluster. When CLASSIFY is specified, the CONVERGE or MXITER specifications on CRITERIA are ignored.

## INITIAL Subcommand

INITIAL specifies the initial cluster centers. Initial cluster centers can also be read from a data file (see [FILE Subcommand on p. 1652](#)).

- One value for each clustering variable must be included for each cluster requested. Values are specified in parentheses cluster by cluster.

**Example**

```
QUICK CLUSTER  A B C D
/CRITERIA = CLUSTER(3)
/INITIAL = (13 24 1 8
           7 12 5 9
           10 18 17 16) .
```

- This example specifies four clustering variables and requests three clusters. Thus, twelve values are supplied on INITIAL.
- The initial center of the first cluster has a value of 13 for variable *A*, 24 for variable *B*, 1 for *C*, and 8 for *D*.

**FILE Subcommand**

Use FILE to obtain initial cluster centers from an external IBM® SPSS® Statistics data file or currently open dataset. (DATASET DECLARE command).

- The only specification is the quoted file specification or dataset name.

**Example**

```
QUICK CLUSTER  A B C D
/FILE=' /data/init.sav'
/CRITERIA = CLUSTER(3) .
```

- In this example, the initial cluster centers are read from file *init.sav*. The file must contain cluster centers for the same four clustering variables specified (*A*, *B*, *C*, and *D*).

**PRINT Subcommand**

QUICK CLUSTER always displays in a Final Cluster Centers table listing the centers used to classify cases and the mean values of the cases in each cluster and a Number of Cases in Each Cluster table listing the number of weighted (if weighting is on) and unweighted cases in each cluster. Use PRINT to request other types of output.

- If PRINT is not specified or is specified without keywords, the default is INITIAL.

<b>INITIAL</b>	<i>Initial cluster centers.</i> When SPLIT FILES is in effect, the initial cluster center for each split file is displayed. This is the default.
<b>CLUSTER</b>	<i>Cluster membership.</i> Each case displays an identifying number or value, the number of the cluster to which it was assigned, and its distance from the center of that cluster. This output is extensive when the number of cases is large.
<b>ID(varname)</b>	<i>Case identification.</i> The value of the specified variable is used in addition to the case numbers to identify cases in output. Case numbers may not be sequential if cases have been selected.
<b>DISTANCE</b>	<i>Pairwise distances between all final cluster centers.</i> This output can consume a great deal of processing time when the number of clusters requested is large.

- ANOVA**            *Descriptive univariate F tests for the clustering variables.* Since cases are systematically assigned to clusters to maximize differences on the clustering variables, these tests are descriptive only and should not be used to test the null hypothesis that there are no differences between clusters. Statistics after clustering are also available through procedure DISCRIMINANT or GLM (GLM is available in the Advanced Statistics option).
- NONE**            *No additional output.* Only the default output is displayed. NONE overrides any other specifications on PRINT.

**Example**

```
QUICK CLUSTER A B C D E
  /CRITERIA=CLUSTERS(6)
  /PRINT=CLUSTER ID(CASEID) DISTANCE.
```

- Six clusters are formed on the basis of the five variables *A*, *B*, *C*, *D*, and *E*.
- For each case in the file, cluster membership and distance from cluster center are displayed. Cases are identified by the values of the variable *CASEID*.
- Distances between all cluster centers are printed.

**OUTFILE Subcommand**

OUTFILE saves the final cluster centers in an external IBM® SPSS® Statistics data file or a previously declared dataset in the current session. You can later use these final cluster centers as initial cluster centers for a different sample of cases that use the same variables. You can also cluster the final cluster centers themselves to obtain clusters of clusters.

- The only specification is a filename or previously declared dataset name for the file. Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. Datasets are available during the current session but are not available in subsequent sessions unless you explicitly save them as data files.
- The program displays the name of the saved file in the procedure information notes.

**Example**

```
QUICK CLUSTER A B C D
  /CRITERIA = CLUSTER(3)
  /OUTFILE = '/data/QC1.sav'.
```

- QUICK CLUSTER writes the final cluster centers to the file *QC1.sav*.

**SAVE Subcommand**

Use SAVE to save results of cluster analysis as new variables in the active dataset.

- You can specify a variable name in parentheses following either keyword. If no variable name is specified, QUICK CLUSTER forms unique variable names by appending an underscore and a sequential number to the rootname *QCL*. The number increments with each new variable saved.

---

*QUICK CLUSTER*

- The program displays the new variables and a short description of each in the procedure information notes.

**CLUSTER**[(varname)]      *The cluster number of each case. The value of the new variable is set to an integer from 1 to the number of clusters.*

**DISTANCE**[(varname)]      *The distance of each case from its classification cluster center.*

**Example**

```
QUICK CLUSTER A B C D
  /CRITERIA=CLUSTERS(6)
  /SAVE=CLUSTER DISTANCE.
```

- Six clusters of cases are formed on the basis of the variables *A*, *B*, *C*, and *D*.
- A new variable *QCL\_1* is created and set to an integer between 1 and 6 to indicate cluster membership for each case.
- Another new variable *QCL\_2* is created and set to the Euclidean distance between a case and the center of the cluster to which it is assigned.

**MISSING Subcommand**

MISSING controls the treatment of cases with missing values.

- LISTWISE, PAIRWISE, and DEFAULT are alternatives. However, each can be used with INCLUDE.

**LISTWISE**      *Delete cases with missing values listwise. A case with a missing value for any of the clustering variables is deleted from the analysis and will not be assigned to a cluster. This is the default.*

**PAIRWISE**      *Assign each case to the nearest cluster on the basis of the clustering variables for which the case has nonmissing values. Only cases with missing values for all clustering variables are deleted.*

**INCLUDE**      *Treat user-missing values as valid.*

**DEFAULT**      *Same as LISTWISE.*

# RANK

```
RANK VARIABLES= varlist [({A**})] [BY varlist]
                        {D }

[/TIES={MEAN** } ]
                        {LOW }
                        {HIGH }
                        {CONDENSE}

[/FRACTION={BLOM**} ]
                        {TUKEY }
                        {VW }
                        {RANKIT}

[/PRINT={YES**} ]
                        {NO }

[/MISSING={EXCLUDE**} ]
                        {INCLUDE }
```

*The following function subcommands can each be specified once:*

```
[/RANK**] [/NTILES(k)] [/NORMAL] [/PERCENT]
[/RFRACTION] [/PROPORTION] [/N] [/SAVAGE]
```

*The following keyword can be used with any function subcommand:*

```
[INTO varname]
```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
RANK VARIABLES=SALARY JOBTIME.
```

## **Overview**

RANK produces new variables containing ranks, normal scores, and Savage and related scores for numeric variables.

### **Options**

**Methods.** You can rank variables in ascending or descending order by specifying A or D on the VARIABLES subcommand. You can compute different rank functions and also name the new variables using the function subcommands. You can specify the method for handling ties on the TIES subcommand, and you can specify how the proportion estimate is computed for the NORMAL and PROPORTIONAL functions on the FRACTION subcommand.

**Format.** You can suppress the display of the summary table that lists the ranked variables and their associated new variables in the active dataset using the `PRINT` subcommand.

### **Basic Specification**

The basic specification is `VARIABLES` and at least one variable from the active dataset. By default, the ranking function is `RANK`. Direction is ascending, and ties are handled by assigning the mean rank to tied values. A summary table that lists the ranked variables and the new variables into which computed ranks have been stored is displayed.

### **Subcommand Order**

- `VARIABLES` must be specified first.
- The remaining subcommands can be specified in any order.

### **Operations**

- `RANK` does not change the way the active dataset is sorted.
- If new variable names are not specified with the `INTO` keyword on the function subcommand, `RANK` creates default names. (Note: The automatically generated new variable names are limited to a maximum length of 8 bytes.)
- `RANK` automatically assigns variable labels to the new variables. The labels identify the source variables. For example, the label for a new variable with the default name `RSALARY` is *RANK of SALARY*.

## **Example**

```
RANK VARIABLES=SALARY JOBTIME.
```

- `RANK` ranks `SALARY` and `JOBTIME` and creates two new variables in the active dataset, `RSALARY` and `RJOBTIME`, which contain the ranks.

## **VARIABLES Subcommand**

`VARIABLES` specifies the variables to be ranked.

- `VARIABLES` is required and must be the first specification on `RANK`. The minimum specification is a single numeric variable. To rank more than one variable, specify a variable list.
- After the variable list, you can specify the direction for ranking in parentheses. Specify `A` for ascending (smallest value gets smallest rank) or `D` for descending (largest value gets smallest rank). `A` is the default.
- To rank some variables in ascending order and others in descending order, use both `A` and `D` in the same variable list. `A` or `D` applies to all preceding variables in the list up to the previous `A` or `D` specification.
- To organize ranks into subgroups, specify keyword `BY` followed by the variable whose values determine the subgroups. The active dataset does not have to be sorted by this variable.
- String variables cannot be specified. Use `AUTORECODE` to recode string variables for ranking.

**Examples**

```
RANK VARIABLES=MURDERS ROBBERY (D) .
```

- RANK ranks *MURDERS* and *ROBBERY* and creates two new variables in the active dataset: *RMURDERS* and *RROBBERY*.
- D specifies descending order of rank. D applies to both *MURDERS* and *ROBBERY*.

```
RANK VARIABLES=MURDERS (D) ROBBERY (A) BY ETHNIC .
```

- Ranks are computed within each group defined by *ETHNIC*. *MURDERS* is ranked in descending order and *ROBBERY* in ascending order within each group of *ETHNIC*. The active dataset does not have to be sorted by *ETHNIC*.

**Function Subcommands**

The optional function subcommands specify different rank functions. RANK is the default function.

- Any combination of function subcommands can be specified for a RANK procedure, but each function can be specified only once.
- Each function subcommand must be preceded by a slash.
- The functions assign default names to the new variables unless keyword INTO is specified.

<b>RANK</b>	<i>Simple ranks.</i> The values for the new variable are the ranks. Rank can either be ascending or descending, as indicated on the VARIABLES subcommand. Rank values can be affected by the specification on the TIES subcommand.
<b>RFRACTION</b>	<i>Fractional ranks.</i> The values for the new variable equal the ranks divided by the sum of the weights of the nonmissing cases. If HIGH is specified on TIES, fractional rank values are an empirical cumulative distribution.
<b>NORMAL</b>	<i>Normal scores</i> (Lehmann, 1975). The new variable contains the inverse of the standard normal cumulative distribution of the proportion estimate defined by the FRACTION subcommand. The default for FRACTION is BLOM.
<b>PERCENT</b>	<i>Fractional ranks as a percentage.</i> The new variable contains fractional ranks multiplied by 100.
<b>PROPORTION</b>	<i>Proportion estimates.</i> The estimation method is specified by the FRACTION subcommand. The default for FRACTION is BLOM.
<b>N</b>	<i>Sum of case weights.</i> The new variable is a constant.
<b>SAVAGE</b>	<i>Savage scores</i> (Lehmann et al., 1975). The new variable contains Savage (exponential) scores.
<b>NTILES(k)</b>	<i>Percentile groups.</i> The new variable contains values from 1 to $k$ , where $k$ is the number of groups to be generated. Each case is assigned a group value, which is the integer part of $1+rk/(w+1)$ , where $r$ is the rank of the case, $k$ is the number of groups specified on NTILES, and $w$ is the sum of the case weights. Group values can be affected by the specification on TIES. There is no default for $k$ .

**INTO Keyword**

INTO specifies variable names for the new variable(s) added to the active dataset. INTO can be used with any of the function subcommands.

- INTO must follow a function subcommand. You must specify the INTO subcommand to assign names to the new variables created by the function.
- You can specify multiple variable names on INTO. The names are assigned to the new variables in the order they are created (the order the variables are specified on the VARIABLES subcommand).
- If you specify fewer names than the new variables, default names are used for the remaining new variables. If you specify more names, the program issues a message and the command is not executed.

If INTO is not specified on a function, RANK creates default names for the new variables according to the following rules:

- The first letter of the ranking function is added to the first seven characters of the original variable name.
- New variable names cannot duplicate variable names in the active dataset or names specified after INTO or generated by default.
- If a new default name is a duplicate, the scheme *XXXnmn* is used, where *XXX* represents the first three characters of the function and *nmn* is a three-digit number starting with 001 and increased by 1 for each variable. (If the ranking function is *N*, *XXX* is simply *N*.) If this naming scheme generates duplicate names, the duplicates are named *RNKXXnm*, where *XX* is the first two characters of the function and *nn* is a two-digit number starting with 01 and increased by 1 for each variable.
- If it is not possible to generate unique names, an error results.

### Example

```
RANK VARIABLES=SALARY
/NORMAL INTO SALNORM
/SAVAGE INTO SALSAV
/NTILES(4) INTO SALQUART.
```

- RANK generates three new variables from variable *SALARY*.
- NORMAL produces the new variable *SALNORM*. *SALNORM* contains normal scores for *SALARY* computed with the default formula BLOM.
- SAVAGE produces the new variable *SALSAV*. *SALSAV* contains Savage scores for *SALARY*.
- NTILES(4) produces the new variable *SALQUART*. *SALQUART* contains the value 1, 2, 3, or 4 to represent one of the four percentile groups of *SALARY*.

## TIES Subcommand

TIES determines the way tied values are handled. The default method is MEAN.

<b>MEAN</b>	<i>Mean rank of tied values is used for ties.</i> This is the default.
<b>LOW</b>	<i>Lowest rank of tied values is used for ties.</i>
<b>HIGH</b>	<i>Highest rank of tied values is used for ties.</i>
<b>CONDENSE</b>	<i>Consecutive ranks with ties sharing the same value.</i> Each distinct value of the ranked variable is assigned a consecutive rank. Ties share the same rank.



**Example**

```
RANK VARIABLES=BURGLARY /RANK INTO RMEAN /TIES=MEAN.
RANK VARIABLES=BURGLARY /RANK INTO RCONDS /TIES=CONDENSE.
RANK VARIABLES=BURGLARY /RANK INTO RHIGH /TIES=HIGH.
RANK VARIABLES=BURGLARY /RANK INTO RLOW /TIES=LOW.
```

- The values of *BURGLARY* and the four new ranking variables are shown below:

BURGLARY	RMEAN	RCONDS	RHIGH	RLOW
0	3	1	5	1
0	3	1	5	1
0	3	1	5	1
0	3	1	5	1
0	3	1	5	1
1	6.5	2	7	6
1	6.5	2	7	6
3	8	3	8	8

**FRACTION Subcommand**

FRACTION specifies the way to compute a proportion estimate  $P$  for the NORMAL and PROPORTION rank functions.

- FRACTION can be used only with function subcommands NORMAL or PROPORTION. If it is used with other function subcommands, FRACTION is ignored and a warning message is displayed.
- Only one formula can be specified for each RANK procedure. If more than one is specified, an error results.

In the following formulas,  $r$  is the rank and  $w$  is the sum of case weights.

**BLOM** *Blom's transformation, defined by the formula  $(r - 3/8) / (w + 1/4)$ . (Blom, 1958) This is the default.*

**RANKIT** *The formula is  $(r - 1/2) / w$ . (Chambers, Cleveland, Kleiner, and Tukey, 1983)*

**TUKEY** *Tukey's transformation, defined by the formula  $(r - 1/3) / (w + 1/3)$ . (Tukey, 1962)*

**VW** *Van der Waerden's transformation, defined by the formula  $r / (w + 1)$ . (Lehmann et al., 1975)*

**Example**

```
RANK VARIABLES=MORTGAGE VALUE /FRACTION=BLOM
/NORMAL INTO MORTNORM VALNORM.
```

- RANK generates new variables *MORTNORM* and *VALNORM*. *MORTNORM* contains normal scores for *MORTGAGE*, and *VALNORM* contains normal scores for *VALUE*.

## ***PRINT Subcommand***

PRINT determines whether the summary tables are displayed. The summary table lists the ranked variables and their associated new variables in the active dataset.

- YES**     *Display the summary tables.* This is the default.  
**NO**       *Suppress the summary tables.*

## ***MISSING Subcommand***

MISSING controls the treatment of user-missing values.

- INCLUDE**     *Include user-missing values.* User-missing values are treated as valid values.  
**EXCLUDE**     *Exclude all missing values.* User-missing values are treated as missing. This is the default.

### ***Example***

```
MISSING VALUE SALARY (0).  
RANK VARIABLES=SALARY /RANK INTO SALRANK /MISSING=INCLUDE.
```

- RANK generates the new variable *SALRANK*.
- INCLUDE causes the user-missing value 0 to be included in the ranking process.

## ***References***

- Blom, G. 1958. *Statistical estimates and transformed beta variables*. New York: John Wiley and Sons.
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- Frigge, M., D. C. Hoaglin, and B. Iglewicz. 1987. Some implementations for the boxplot. In: *Computer Science and Statistics Proceedings of the 19th Symposium on the Interface*, R. M. Heiberger, and M. Martin, eds. Alexandria, Virginia: American Statistical Association.
- Lehmann, E. L. 1975. *Nonparametrics: Statistical methods based on ranks*. San Francisco: Holden-Day.
- Tukey, J. W. 1962. The future of data analysis. *Annals of Mathematical Statistics*, 33:22, 1–67.

# RATIO STATISTICS

RATIO STATISTICS is available in the Statistics Base option.

```
RATIO STATISTICS numerator varname WITH denominator varname
  [BY group varname[({ASCENDING**})]]
                    {DESCENDING }
                    {NOSORT     }

[/MISSING = {EXCLUDE**}]
           {INCLUDE  }

[/OUTFILE('file'|'dataset') = [AAD] [BCOC((low,high) [(low,high)] ...)] [CIN[({95 } )]]]
                               {value}
                               [COD] [MAX] [MDCOV] [MEAN] [MEDIAN] [MIN] [MNCOV] [PRD]
                               [RANGE] [STDDEV] [WCOC(value list)] [WGTMEAN]]

[/PRINT = [AAD] [BCOC(low,high)...] [CIN[({95 } )]]]
          {value}
          [COD] [MAX] [MDCOV] [MEAN] [MEDIAN] [MIN] [MNCOV] [PRD]
          [RANGE] [STDDEV] [WCOC(value list)] [WGTMEAN]]
```

\*\* Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
RATIO STATISTICS appraise WITH price
  /PRINT = AAD BCOC((1,2) (3,4)) MEAN.
```

## Overview

RATIO STATISTICS provides a variety of descriptive statistics for the ratio between two variables.

### Basic Specification

The minimum specification is a numerator variable and a denominator variable, and either an OUTFILE subcommand or a PRINT subcommand.

### Subcommand Order

- The variable list must be specified first.
- Subcommands can be specified in any order.

### Syntax Rules

- Empty subcommands are silently ignored.
- All subcommands should be specified only once. If a subcommand is repeated, only the last specification will be used.

- The following words are reserved as keywords in this procedure: BY and WITH.

## Examples

```
RATIO STATISTICS lastval WITH saleval BY town (ASCENDING)
/MISSING = EXCLUDE
/PRINT = BCOC(0.8,1.2) COD MEDIAN WCOC( 20 ) .
```

- The procedure computes ratios by dividing *lastval* by *saleval*. Separate ratio statistics are computed for each category of *town*, with values displayed in ascending order.
- By default, user-missing values are excluded from the analysis.
- The PRINT subcommand requests the coefficient of concentration for the range 0.8-1.2, the coefficient of dispersion, the median, and the coefficient of concentration for ratios falling within 20 percent of the median.

## Case Frequency

- If a WEIGHT variable is specified, its values are used as frequency weights by this procedure.
- Cases with missing or nonpositive weights are not used for computing the ratio statistics.
- The weight values are rounded to the nearest whole numbers before use. For example, 0.5 is rounded to 1, and 2.4 is rounded to 2.

## Variable List

The variable list specifies the numerator variable, denominator variable, and optional group variable.

- The numerator variable must be the first specification after the procedure name.
- The denominator variable must be preceded by the keyword WITH.
- The group variable, if specified, must be preceded by the keyword BY.
- Both the numerator and the denominator variables must be numeric.
- The group variable can be of any type (numeric or string).
- By default or when the keyword ASCENDING is specified, values of the group variable are displayed in ascending order. Specify the keyword DESCENDING to display in descending order. Specify NOSORT to preserve the appearance order in the data.
- Only cases with no (system- or user-) missing values in both the numerator and the denominator variables will be used. Please note that this rule does not apply to the group variable.

### Example

```
RATIO STATISTICS appraise WITH price
/PRINT = AAD BCOC((1,2) (3,4)) MEAN.
```

- This is a typical analysis where *appraise* is the appraised value and *price* is the transaction price. The ratio is computed by dividing *appraise* by *price*.

**Example**

```
RATIO STATISTICS appraise WITH price BY county
  /PRINT = CIN(90) MEDIAN.
```

- The ratio is still computed by dividing *appraise* by *price*. However, separate ratio statistics are requested for each category of *county*.

**MISSING Subcommand**

MISSING specifies the way to handle cases with user-missing values.

- A case is never used if it contains system-missing values in the numerator and/or the denominator variables.
- If this subcommand is not specified, the default is EXCLUDE.
- Keywords EXCLUDE and INCLUDE are mutually exclusive. Only one of them can be specified once.

**EXCLUDE**      *Exclude both user-missing and system-missing values. This is the default.*

**INCLUDE**      *User-missing values are treated as valid. System-missing values cannot be included in the analysis.*

**OUTFILE Subcommand**

OUTFILE saves the requested statistics to an external IBM® SPSS® Statistics data file or a previously declared dataset (DATASET DECLARE command).

- The requested statistics are saved in a single record in the external file.
- If a group variable has been specified, the requested statistics at each category of the group variable will also be saved as additional records in the external file.
- The file specification or dataset name should be quoted and enclosed in quotes.

The following statistics are available.

<b>AAD</b>	<i>Average absolute deviation.</i> The result of summing the absolute deviations of the ratios about the median and dividing the result by the total number of ratios.
<b>BCOC (low,high) ...)</b>	<i>Coefficient of concentration.</i> The percentage of ratios that fall into an interval. Pairs of low and high values enclosed in parentheses specify the intervals.
<b>CIN(a)</b>	<i>Confidence interval.</i> Specifying this keyword displays confidence intervals for the mean, median, and weighted mean (if those statistics are requested). Specify a value greater than or equal to 0 and less than 100 as the confidence level.
<b>COD</b>	<i>Coefficient of dispersion.</i> The result of expressing the average absolute deviation as a percentage of the median.
<b>MAX</b>	<i>Maximum.</i> The largest ratio.
<b>MDCOV</b>	<i>Median-centered coefficient of variation.</i> The result of expressing the root mean squares of deviation from the median as a percentage of the median.

## RATIO STATISTICS

<b>MEAN</b>	<i>Mean.</i> The result of summing the ratios and dividing the result by the total number ratios.
<b>MEDIAN</b>	<i>Median.</i> The value such that number of ratios less than this value and the number of ratios greater than this value are the same.
<b>MIN</b>	<i>Minimum.</i> The smallest ratio.
<b>MNCOV</b>	<i>Mean-centered coefficient of variation.</i> The result of expressing the standard deviation as a percentage of the mean.
<b>PRD</b>	<i>Price-related differential.</i> Also known as the index of regressivity, the result of dividing the mean by the weighted mean.
<b>RANGE</b>	<i>Range.</i> The result of subtracting the minimum ratio from the maximum ratio.
<b>STDDEV</b>	<i>Standard deviation.</i> The result of summing the squared deviations of the ratios about the mean, dividing the result by the total number of ratios minus one, and taking the positive square root.
<b>WCOC(value list)</b>	<i>Coefficient of concentration.</i> The percentage of ratios that fall within the specified percent of the median. Specify a list of values that are greater than 0 and less than 100.
<b>WGTMEAN</b>	<i>Weighted mean.</i> The result of dividing the mean of the numerator by the mean of the denominator. It is also the mean of the ratios weighted by the denominator.

**Example**

```
RATIO STATISTICS appraise WITH price BY county
  /OUTFILE('/PropertyTax/Ratio.sav') = CIN(90) MEDIAN.
```

- The median ratios and their 90% confidence intervals at each category of *county* are saved to */PropertyTax/Ratio.sav*.
- The overall median ratio and its 90% confidence intervals are also saved.

**PRINT Subcommand**

PRINT displays optional output. If no PRINT subcommand is specified, only a case processing summary table is displayed by default.

<b>AAD</b>	<i>Average absolute deviation.</i> The result of summing the absolute deviations of the ratios about the median and dividing the result by the total number of ratios.
<b>BCOC(low,high) ...)</b>	<i>Coefficient of concentration.</i> The percentage of ratios that fall into an interval. Pairs of low and high values enclosed in parentheses specify the intervals.
<b>CIN(a)</b>	<i>Confidence interval.</i> Specifying this keyword displays confidence intervals for the mean, median, and weighted mean (if those statistics are requested). Specify a value greater than or equal to 0 and less than 100 as the confidence level.
<b>COD</b>	<i>Coefficient of dispersion.</i> The result of expressing the average absolute deviation as a percentage of the median.
<b>MAX</b>	<i>Maximum.</i> The largest ratio.
<b>MDCOV</b>	<i>Median-centered coefficient of variation.</i> The result of expressing the root mean squares of deviation from the median as a percentage of the median.

<b>MEAN</b>	<i>Mean.</i> The result of summing the ratios and dividing the result by the total number ratios.
<b>MEDIAN</b>	<i>Median.</i> The value such that number of ratios less than this value and the number of ratios greater than this value are the same.
<b>MIN</b>	<i>Minimum.</i> The smallest ratio.
<b>MNCOV</b>	<i>Mean-centered coefficient of variation.</i> The result of expressing the standard deviation as a percentage of the mean.
<b>PRD</b>	<i>Price-related differential.</i> Also known as the index of regressivity, the result of dividing the mean by the weighted mean.
<b>RANGE</b>	<i>Range.</i> The result of subtracting the minimum ratio from the maximum ratio.
<b>STDDEV</b>	<i>Standard deviation.</i> The result of summing the squared deviations of the ratios about the mean, dividing the result by the total number of ratios minus one, and taking the positive square root.
<b>WCOC(value list)</b>	<i>Coefficient of concentration.</i> The percentage of ratios that fall within the specified percentage of the median. Specify a list of values that are greater than 0 and less than 100.
<b>WGTMEAN</b>	<i>Weighted mean.</i> The result of dividing the mean of the numerator by the mean of the denominator. It is also the mean of the ratios weighted by the denominator.

### **Example**

```
RATIO STATISTICS appraise WITH price BY county
  /PRINT = BCOC((0.5,0.9) (1.3,1.5)) WCOC(15 30 45) MEDIAN PRD.
```

- The median ratios and priced related differentials at each category of *county* are displayed. The overall median ratio and the overall price-related differential are also displayed.
- Five coefficients of concentration are also displayed. The first two COC are percentages of ratios that fall into the intervals: (0.5, 0.9) and (1.3, 1.5). The next three COC are percentages of ratios that fall within 15% of the median, 30% of the median, and 45% of the median.

# RBF

RBF is available in the Neural Networks option.

```
RBF dependent variable [(MLEVEL = {S})] [dependent variable...]
                        {O}
                        {N}

    [BY factor list] [WITH covariate list]

[/EXCEPT VARIABLES = varlist]

[/RESCALE [COVARIATE = {STANDARDIZED**}]   [DEPENDENT = {STANDARDIZED**}]]
           {NORMALIZED   }                 {NORMALIZED   }
           {ADJNORMALIZED }                 {ADJNORMALIZED }
           {NONE         }                 {NONE         }

[/PARTITION {TRAINING = {70** } TESTING = {30** } HOLDOUT = {0** }}]
            {number}      {number}      {number}
            {VARIABLE = varname }

[/ARCHITECTURE [{[MINUNITS = {AUTO** } MAXUNITS = {AUTO** }]]]
               {integer}      {integer}
               {NUMUNITS = integer }

               [HIDDENFUNCTION = {NRBF**}]]
               {ORBF }

[/CRITERIA OVERLAP = {AUTO**}]
                   {number}

[/MISSING USERMISSING = {EXCLUDE**}]
                       {INCLUDE }

[/PRINT [CPS**] [NETWORKINFO**] [SUMMARY**] [CLASSIFICATION**]
        [SOLUTION] [IMPORTANCE] [NONE]]

[/PLOT [NETWORK**] [PREDICTED] [RESIDUAL] [ROC]
       [GAIN] [LIFT] [NONE]]

[/SAVE [PREDDVAL[(varname [varname...])]
       [PSEUDOPROB[(rootname[:{25 } ] [rootname...])]]]
       {integer}

[/OUTFILE MODEL = 'file' ['file'...]]
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- Command introduced.

## **Example**

```
RBF dep_var BY A B C WITH X Y Z.
```



## Overview

Neural networks are a data mining tool for finding unknown patterns in databases. Neural networks can be used to make business decisions by forecasting demand for a product as a function of price and other variables or by categorizing customers based on buying habits and demographic characteristics. The RBF procedure fits a radial basis function neural network, which is a feedforward, supervised learning network with an input layer, a hidden layer called the radial basis function layer, and an output layer. The hidden layer transforms the input vectors into radial basis functions. Like the MLP (multilayer perceptron) procedure, the RBF procedure performs prediction and classification.

The RBF procedure trains the network in two stages:

1. The procedure determines the radial basis functions using clustering methods. The center and width of each radial basis function are determined.
2. The procedure estimates the synaptic weights given the radial basis functions. The sum-of-squares error function with identity activation function for the output layer is used for both prediction and classification. Ordinary least squares regression is used to minimize the sum-of-squares error.

Because of this two-stage training approach, the RBF network is in general trained much faster than MLP.

### Options

**Prediction or classification.** One or more dependent variables may be specified, and they may be scale, categorical, or a combination. If a dependent variable has scale measurement level, then the neural network predicts continuous values that approximate the “true” value of some continuous function of the input data. If a dependent variable is categorical, then the neural network is used to classify cases into the “best” category based on the input predictors.

**Rescaling.** RBF optionally rescales covariates (predictors with scale measurement level) or scale dependent variables before training the neural network. There are three rescaling options: standardization, normalization, and adjusted normalization.

**Training, testing, and holdout data.** RBF optionally divides the dataset into training, testing, and holdout data. The neural network is trained using the training data. The testing data can be used to determine the “best” number of hidden units for the network. The holdout data is completely excluded from the training process and is used for independent assessment of the final network.

**Architecture selection.** The RBF procedure creates a neural network with one hidden layer and can perform automatic architecture selection to find the “best” number of hidden units. By default, the procedure automatically computes a reasonable range and finds the “best” number within the range. However, you can override these computations by providing your own range or a specific number of hidden units.

**Activation functions.** Units in the hidden layer can use the normalized radial basis function or the ordinary radial basis function.

**Missing values.** The RBF procedure has an option for treating user-missing values of categorical variables as valid. User-missing values of scale variables are always treated as invalid.

**Output.** RBF displays pivot table output but offers an option for suppressing most such output. Graphical output includes a network diagram (default) and a number of optional charts: predicted by observed values, residual by predicted values, ROC (Receiver Operating Characteristic) curves, cumulative gains, lift, and independent variable importance. The procedure also optionally saves predicted values in the active dataset. Hidden unit center and width vectors and synaptic weight estimates can be saved in XML files.

### ***Basic Specification***

The basic specification is the RBF command followed by one or more dependent variables, the BY keyword and one or more factors, and the WITH keyword and one or more covariates.

By default, the RBF procedure standardizes covariates and scale dependent variables and selects a training sample before training the neural network. Automatic architecture selection is used to find the “best” neural network architecture. User-missing values are excluded and default pivot table output is displayed.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### ***Syntax Rules***

- All subcommands are optional.
- Subcommands may be specified in any order.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not allowed.
- Any split variable defined on the SPLIT FILE command may not be used as a dependent variable, factor, covariate, or partition variable.

### ***Limitations***

Frequency weights specified on the WEIGHT command are ignored with a warning by the RBF procedure.

### ***Categorical Variables***

The RBF procedure temporarily recodes categorical predictors and dependent variables using one-of- $c$  coding for the duration of the procedure. If there are  $c$  categories of a variable, then the variable is stored as  $c$  vectors, with the first category denoted  $(1,0,\dots,0)$ , the next category,  $(0,1,0,\dots,0)$ , ..., and the final category,  $(0,0,\dots,0,1)$ .

Because of the one-of-*c* coding, the total number of input units is the number of scale predictors plus the number of categories across all categorical predictors. However, unlike the multilayer perceptron (MLP), this coding scheme does not increase the number of synaptic weights for categorical predictors and hence should not significantly increase the training time.

All one-of-*c* coding is based on the training data, even if a testing or holdout sample is defined (see [PARTITION Subcommand](#)). Thus, if the testing or holdout samples contain cases with predictor categories that are not present in the training data, then those cases are not used by the procedure or in scoring. If the testing or holdout samples contain cases with dependent variable categories that are not present in the training data, then those cases are not used by the procedure but they may be scored.

### **Replicating Results**

The RBF procedure uses random number generation during random assignment of partitions. To reproduce the same randomized results in the future, use the `SET` command to set the initialization value for the random number generator before each run of the RBF procedure.

RBF results are also dependent on data order because the two-step cluster algorithm is used to determine the radial basis functions. To minimize data order effects, randomly order the cases before running the RBF procedure. To verify the stability of a given solution, you may want to obtain several different solutions with cases sorted in different random orders. In situations with extremely large file sizes, multiple runs can be performed with a sample of cases sorted in different random orders.

In summary, if you want to exactly replicate RBF results in the future, use the same initialization value for the random number generator and the same data order, in addition to using the same RBF procedure settings.

## **Examples**

### **Basic specification with default neural network settings**

```
RBF DepVar BY A B C WITH X Y Z.
```

- The RBF procedure treats *DepVar* as the dependent variable. Predictors *A*, *B*, and *C* are factors, and *X*, *Y*, and *Z* are covariates.
- By default, covariates and the dependent variable (if it has scale measurement level) are standardized before training. Also, the active dataset is partitioned into training and testing data samples, with 70% going to the training data and 30% to the testing data sample.
- Automatic architecture selection is used to find the “best” neural network architecture.
- User-missing values are excluded and default output is displayed.

### **User-specified neural network with three hidden units**

```
RBF DepVar BY A B C WITH X Y Z
  /ARCHITECTURE
  NUMUNITS=3
  HIDDENFUNCTION=ORBF.
```

- The RBF procedure treats *DepVar* as the dependent variable. Predictors *A*, *B*, and *C* are factors, and *X*, *Y*, and *Z* are covariates.
- By default, covariates and the dependent variable (if it has scale measurement level) are standardized before training. Also, the active dataset is partitioned into training and testing data samples, with 70% going to the training data and 30% to the testing data sample.
- The ARCHITECTURE subcommand specifies a neural network with three hidden units, thereby overriding automatic architecture selection. In addition, the hidden layer uses the ordinary radial basis function.
- User-missing values are excluded and default output is displayed.

### ***Nominal dependent variable; user-specified rescaling and partitions***

```
*Radial Basis Function Network.
RBF custcat (MLEVEL=N) BY marital ed retire gender WITH age address
    income employ reside
  /RESCALE COVARIATE=ADJNORMALIZED
  /PARTITION TRAINING=6 TESTING=2 HOLDOUT=1
  /ARCHITECTURE MINUNITS=AUTO MAXUNITS=AUTO HIDDENFUNCTION=NRBF
  /CRITERIA OVERLAP=AUTO
  /PRINT CPS NETWORKINFO SUMMARY CLASSIFICATION
  /PLOT ROC GAIN LIFT PREDICTED
  /SAVE PREDVAL PSEUDOPROB
  /MISSING USERMISSING=EXCLUDE .
```

- The procedure constructs a network to predict the nominal-level variable *custcat*, using *marital*, *ed*, *retire*, and *gender* as factors and *age*, *address*, *income*, *employ*, and *reside* as covariates.
- The RESCALE subcommand specifies that covariates should be rescaled using the adjusted normalized method.
- The PARTITION subcommand specifies that the relative number of cases assigned to the training:testing:holdout samples should be 6:2:1. This assigned 2/3 of the cases to training, 2/9 to testing, and 1/9 to holdout.
- The PLOT subcommand does not request the default graphical output (the network diagram), but instead an ROC curve, cumulative gains chart, lift chart, and predicted-by-residuals chart will be produced.
- The SAVE subcommand requests that the predicted values and predicted pseudo-probabilities of *custcat* be saved to the active dataset.
- All other options are set to their default values.

## ***Variable Lists***

The command line variable lists specify the dependent variables, any categorical predictors (also known as factors), and any scale predictors (also known as covariates).

### ***Dependent Variables***

- A list of one or more dependent variables must be the first specification on the RBF command.

- Each dependent variable may be followed by the measurement level specification, which contains, in parentheses, the `MLEVEL` keyword followed by an equals sign and then `S` for scale, `O` for ordinal, or `N` for nominal. `RBF` treats ordinal and nominal dependent variables equivalently as categorical.
- If a measurement level is specified, then it temporarily overrides a dependent variable's setting in the data dictionary.
- If no measurement level is specified, then `RBF` defaults to the dictionary setting.
- If a measurement level is not specified and no setting is recorded in the data dictionary, then a numeric variable is treated as scale and a string variable is treated as categorical.
- Dependent variables can be numeric or string.
- A string variable may be defined as ordinal or nominal only.

### ***Predictor Variables***

- The names of the factors, if any, must be preceded by the keyword `BY`.
- If the keyword `BY` is specified with no factors, then a warning is issued and `BY` is ignored.
- The names of the covariates, if any, must be preceded by the keyword `WITH`.
- If the keyword `WITH` is specified with no covariates, then a warning is issued and `WITH` is ignored.
- If the dependent variable is specified within a factor list or a covariate list, then it is ignored in the list.
- All variables specified within a factor or covariate list must be unique. If duplicate variables are specified within a list, then the duplicates are ignored.
- If duplicate variables are specified across the factor and covariate lists, then an error is issued.
- Factor variables can be numeric or string.
- Covariates must be numeric.
- At least one predictor must be specified.

## ***EXCEPT Subcommand***

The `EXCEPT` subcommand lists any variables that the `RBF` procedure should exclude from the factor or covariate lists on the command line. This subcommand is useful if the factor or covariate lists contain a large number of variables—specified using the `TO` or `ALL` keyword, for example—but there are a few variables (for example, Case ID) that should be excluded.

The `EXCEPT` subcommand is introduced strictly for the purpose of simplifying syntax. Missing values on factors or covariates specified on `EXCEPT` do not affect whether a case is included in the analysis. For example, the following two `RBF` commands are equivalent. In both commands, listwise deletion is based on the dependent variable and factors *A*, *B*, and *C*.

```
RBF DepVar BY A B C.
```

```
RBF DepVar BY A B C D /EXCEPT VARIABLES=D.
```

- The `EXCEPT` subcommand ignores duplicate variables and variables that are not specified on the command line's factor or covariate lists.
- There is no default variable list on the `EXCEPT` subcommand.

## **RESCALE Subcommand**

The `RESCALE` subcommand is used to rescale covariates or scale dependent variables.

All rescaling is performed based on the training data, even if a testing or holdout sample is defined (see [PARTITION Subcommand](#)). That is, depending on the type of rescaling, the mean, standard deviation, minimum value, or maximum value of a covariate or dependent variable is computed using only the training data. It is important that these covariates or dependent variables have similar distributions across the training, testing, and holdout samples. If the data are partitioned by specifying percentages on the `PARTITION` subcommand, then the `RBF` procedure attempts to ensure this similarity by random assignment. However, if you use the `PARTITION` subcommand `VARIABLE` keyword to assign cases to the training, testing, and holdout samples, then we recommend that you confirm that the distributions are similar across samples before running the `RBF` procedure.

### **COVARIATE Keyword**

The `COVARIATE` keyword specifies the rescaling method to use for covariates specified following `WITH` on the command line.

If no covariates are specified on the command line, then the `COVARIATE` keyword is ignored.

<b>STANDARDIZED</b>	<i>Subtract the mean and divide by the standard deviation, <math>(x-\text{mean})/s</math>. This is the default rescaling method for covariates.</i>
<b>NORMALIZED</b>	<i>Subtract the minimum and divide by the range, <math>(x-\text{min})/(\text{max}-\text{min})</math>.</i>
<b>ADJNORMALIZED</b>	<i>Adjusted version of subtracting the minimum and dividing by the range, <math>[2*(x-\text{min})/(\text{max}-\text{min})]-1</math>.</i>
<b>NONE</b>	<i>No rescaling of covariates.</i>

### **DEPENDENT Keyword**

The `DEPENDENT` keyword specifies the rescaling method to use for the dependent variables.

This keyword is applied only for scale dependent variables; that is, either `MLEVEL=S` is specified on the command line or the variable has a scale measurement level based on its data dictionary setting. If a dependent variable is not scale, then the `DEPENDENT` keyword is ignored for that variable.

<b>STANDARDIZED</b>	<i>Subtract the mean and divide by the standard deviation, <math>(x-\text{mean})/s</math>. This is the default rescaling method for scale dependent variables.</i>
<b>NORMALIZED</b>	<i>Subtract the minimum and divide by the range, <math>(x-\text{min})/(\text{max}-\text{min})</math>.</i>
<b>ADJNORMALIZED</b>	<i>Adjusted version of subtracting the minimum and dividing by the range, <math>[2*(x-\text{min})/(\text{max}-\text{min})]-1</math>.</i>
<b>NONE</b>	<i>No rescaling of scale dependent variables.</i>

## ***PARTITION Subcommand***

The `PARTITION` subcommand specifies the method of partitioning the active dataset into training, testing, and holdout samples. The training sample comprises the data records used to train the neural network. The testing sample is an independent set of data records used to track prediction error during training in order to prevent overtraining. The holdout sample is another independent set of data records used to assess the final neural network.

- The partition can be defined by specifying the ratio of cases randomly assigned to each sample (training, testing, and holdout) or by a variable that assigns each case to the training, testing, or holdout sample.
- If the `PARTITION` subcommand is not specified, then the default partition randomly assigns 70% of the cases to the training sample, 30% to the testing sample, and 0% to the holdout sample. If you want to specify a different random assignment, then you must specify new values for the `TRAINING`, `TESTING`, and `HOLDOUT` keywords. The value specified on each keyword gives the relative number of cases in the active dataset to assign to each sample. For example, `/PARTITION TRAINING = 50 TESTING = 30 HOLDOUT = 20` is equivalent to `/PARTITION TRAINING = 5 TESTING = 3 HOLDOUT = 2`; both subcommands randomly assign 50% of the cases to the training sample, 30% to the testing sample, and 20% to the holdout sample.
- If you want to be able to reproduce results based on the `TRAINING`, `TESTING`, and `HOLDOUT` keywords later, use the `SET` command to set the initialization value for the random number generator before running the RBF procedure.
- Be aware of the relationship between rescaling and partitioning. [For more information, see the topic `RESCALE Subcommand` on p. 1672.](#)
- All partitioning is performed after listwise deletion of any cases with invalid data for any variable used by the procedure. See [MISSING Subcommand](#) for details about valid and invalid data.

### ***TRAINING Keyword***

The `TRAINING` keyword specifies the relative number of cases in the active dataset to randomly assign to the training sample. The value must be an integer greater than 0. The default (if the `PARTITION` subcommand is not specified) is 70.

### ***TESTING Keyword***

The `TESTING` keyword specifies the relative number of cases in the active dataset to randomly assign to the testing sample. The value must be an integer greater than 0. The default (if the `PARTITION` subcommand is not specified) is 30.

### ***HOLDOUT Keyword***

The `HOLDOUT` keyword specifies the relative number of cases in the active dataset to randomly assign to the holdout sample. The value must be an integer greater than 0. The default (if the `PARTITION` subcommand is not specified) is 0.

**VARIABLE Keyword**

The `VARIABLE` keyword specifies a variable that assigns each case in the active dataset to the training, testing, or holdout sample. Cases with a positive value on the variable are assigned to the training sample, cases with a value of 0, to the testing sample, and cases with a negative value, to the holdout sample. Cases with a system-missing value are excluded from the analysis. (Any user-missing values for the partition variable are always treated as valid.)

The variable may not be the dependent variable or any variable specified on the command line factor or covariate lists. The variable must be numeric.

**ARCHITECTURE Subcommand**

The `ARCHITECTURE` subcommand is used to specify the neural network architecture. `RBF` creates a neural network with one hidden layer. There are three ways of choosing the number of hidden units:

1. By default, `RBF` uses automatic architecture selection to find the “best” number of hidden units within a range. The procedure automatically computes the minimum and maximum and finds the best number of hidden units within the range.

If a testing sample is defined; that is, if `/PARTITION TESTING` is specified with a number greater than zero or via `/PARTITION VARIABLE`, then the procedure uses the testing data criterion: The best number of hidden units is the one that yields the smallest error in the testing data. If a testing sample is not defined, then the procedure uses the Bayesian information criterion (BIC): The best number of hidden units is the one that yields the smallest BIC based on the training data.

2. You can provide your own range by specifying the `MINUNITS` and `MAXUNITS` keywords with integers, and the procedure will find the “best” number of hidden units within that range. It is invalid to specify only `MINUNITS` or only `MAXUNITS`; both must be given. The best number of hidden units from the range is determined using the testing data criterion or the BIC.
3. You can override the use of a range and specify a particular number of units directly using the `NUMUNITS` keyword. It is invalid to specify `NUMUNITS` with either the `MINUNITS` or `MAXUNITS` keyword.

**MINUNITS Keyword**

The `MINUNITS` keyword specifies the minimum number of units to use in the range. Value `AUTO` indicates that the number should be computed automatically. Alternatively, a positive integer may be specified. The integer must be less than the number specified on the `MAXUNITS` keyword.

The default value is `AUTO`.

**MAXUNITS Keyword**

The `MAXUNITS` keyword specifies the maximum number of units to use in the range. Value `AUTO` indicates that the number should be computed automatically. Alternatively, a positive integer may be specified. The integer must be greater than the number specified on the `MINUNITS` keyword.



The default value is AUTO.

### **NUMUNITS Keyword**

The NUMUNITS keyword specifies a particular number of hidden units instead of a range. A positive integer may be specified.

- If NUMUNITS is specified when testing data are defined (see [PARTITION Subcommand](#)), then the testing data are not used in determining the network architecture and a warning is issued.
- There is no default value.

### **HIDDENFUNCTION Keyword**

The HIDDENFUNCTION keyword specifies the Gaussian radial basis function used in the hidden layer.

- |             |   |
|-------------|---|
| <b>NRBF</b> | <i>Normalized radial basis function.</i> Uses the softmax activation function so the activations of all hidden units are normalized to sum to 1. This is the default activation function for all units in the hidden layer. |
| <b>ORBF</b> | <i>Ordinary radial basis function.</i> Uses the exponential activation function so the activation of the hidden unit is a Gaussian “bump” as a function of the inputs.  |

## **CRITERIA Subcommand**

The CRITERIA subcommand specifies computational settings for the RBF procedure.

### **OVERLAP Keyword**

The OVERLAP keyword specifies the hidden-unit overlapping factor that controls how much overlap occurs among the hidden units.

- Specify AUTO to automatically compute the overlapping factor  $1+0.1d$ , where  $d$  is the number of input units. (The number of input units equals the sum of the number of categories across all factors and the number of covariates.) This is the default.
- Alternatively, specify a number greater than 0 to request a particular overlapping factor.

## **MISSING Subcommand**

The MISSING subcommand is used to control whether user-missing values for categorical variables; that is, factors and categorical dependent variables, are treated as valid values.

- By default, user-missing values for categorical variables are treated as invalid.
- User-missing values for scale variables are always treated as invalid. System-missing values for any variables are always treated as invalid.

### **USERMISSING=EXCLUDE**

*User-missing values for categorical variables are treated as invalid. This is the default.*

**USERMISSING=INCLUDE**

*User-missing values for categorical variables are treated as valid values.*

## **PRINT Subcommand**

The PRINT subcommand indicates the tabular output to display and can be used to request a sensitivity analysis. If PRINT is not specified, then the default tables are displayed. If PRINT is specified, then only the requested PRINT output is displayed.

### **CPS Keyword**

The CPS keyword displays the case processing summary table, which summarizes the number of cases included and excluded in the analysis, in total and by training, testing, and holdout samples. This table is shown by default.

### **NETWORKINFO Keyword**

The NETWORKINFO keyword displays information about the neural network, including the dependent variables, number of input and output units, number of hidden units, and activation functions. This table is shown by default.

### **SUMMARY Keyword**

The SUMMARY keyword displays a summary of the neural network results, including the error, the relative error or percentage of incorrect predictions, and the training time.

- The error is the sum-of-squares error. In addition, relative errors or percentages of incorrect predictions are displayed, depending on the dependent variable measurement levels. If any dependent variable has scale measurement level, then the average overall relative error (relative to the mean model) is displayed. If all dependent variables are categorical, then the average percentage of incorrect predictions is displayed. Relative errors or percentages of incorrect predictions are also displayed for individual dependent variables.
- Summary results are given for the training data, and for testing and holdout data if they exist.
- This table is shown by default.

### **CLASSIFICATION Keyword**

The CLASSIFICATION keyword displays a classification table for each categorical dependent variable. The table gives the number of cases classified correctly and incorrectly for each dependent variable category.

- In addition to classification tables, the CLASSIFICATION keyword reports the percentage of the total cases that were correctly classified. A case is correctly classified if its highest predicted probabilities correspond to the observed categories for that case.
- Classification results are given for the training data, and for testing and holdout data if they exist.

- Classification results are shown by default.
- The `CLASSIFICATION` keyword is ignored for scale dependent variables.

### ***SOLUTION Keyword***

The `SOLUTION` keyword displays the center and width vectors for each hidden unit, and the synaptic weight estimates (that is, the coefficient estimates) from the hidden units to the output units. These results are based on the training sample even if the active dataset is partitioned into training, testing, and holdout data.

This table is not shown by default because the number of synaptic weight estimates may be large, and these weights are generally not used for interpretation.

### ***IMPORTANCE Keyword***

The `IMPORTANCE` keyword performs a sensitivity analysis, which computes the importance of each predictor in determining the neural network. The analysis is based on the combined training and testing samples, or only the training sample if there is no testing sample. This keyword creates a table and a chart displaying importance and normalized importance for each predictor.

Sensitivity analysis is not performed by default because it is computationally expensive and time-consuming if there are a large number of predictors or cases.

### ***NONE Keyword***

The `NONE` keyword suppresses all `PRINT` output. This keyword may not be specified with any other `PRINT` keywords.

## ***PLOT Subcommand***

The `PLOT` subcommand indicates the chart output to display. If `PLOT` is not specified, then the default chart (the network diagram) is displayed. If `PLOT` is specified, then only the requested `PLOT` output is displayed.

### ***NETWORK Keyword***

The `NETWORK` keyword displays the network diagram. This chart is shown by default.

### ***PREDICTED Keyword***

The `PREDICTED` keyword displays a predicted-by-observed-value chart for each dependent variable. For categorical dependent variables, a boxplot of predicted pseudo-probabilities is displayed. For scale dependent variables, a scatterplot is displayed.

- Charts are based on the combined training and testing samples, or only the training sample if there is no testing sample.

**RESIDUAL Keyword**

The `RESIDUAL` keyword displays a residual-by-predicted-value chart for each scale dependent variable.

- This chart is available only for scale dependent variables. The `RESIDUAL` keyword is ignored for categorical dependent variables.
- Charts are based on the combined training and testing samples, or only the training sample if there is no testing sample.

**ROC Keyword**

The `ROC` keyword displays an ROC (Receiver Operating Characteristic) chart for each categorical dependent variable. It also displays a table giving the area under each curve in the chart.

- For a given dependent variable, the ROC chart displays one curve for each category. If the dependent variable has two categories, then each curve treats the category at issue as the positive state versus the other category. If the dependent variable has more than two categories, then each curve treats the category at issue as the positive state versus the aggregate of all other categories.
- This chart is available only for categorical dependent variables. The `ROC` keyword is ignored for scale dependent variables.
- Charts are based on the combined training and testing samples, or only the training sample if there is no testing sample.

**GAIN Keyword**

The `GAIN` keyword displays a cumulative gains chart for each categorical dependent variable.

- The display of one curve for each dependent variable category is the same as for the `ROC` keyword.
- This chart is available only for categorical dependent variables. The `GAIN` keyword is ignored for scale dependent variables.
- Charts are based on the combined training and testing samples, or only the training sample if there is no testing sample.

**LIFT Keyword**

The `LIFT` keyword displays a lift chart for each categorical dependent variable.

- The display of one curve for each dependent variable category is the same as for the `ROC` keyword.
- This chart is available only for categorical dependent variables. The `LIFT` keyword is ignored for scale dependent variables.
- Charts are based on the combined training and testing samples, or only the training sample if there is no testing sample.

### ***NONE* Keyword**

The NONE keyword suppresses all PLOT output. This keyword may not be specified with any other PLOT keywords.

## ***SAVE* Subcommand**

The SAVE subcommand writes optional temporary variables to the active dataset.

### **PREDVAL(varname varname...)**

*Predicted value or category.* If a dependent variable has scale measurement level, then this is the predicted value. If a dependent variable is categorical, then this is the predicted category.

Specify one or more unique, valid variable names. There should be as many variable names specified as there are dependent variables, and the names should be listed in the order of the dependent variables on the command line. If you do not specify enough variable names, then default names are used for any remaining variables. If you specify too many variable names, then any remaining names are ignored.

If there is only one dependent variable, then the default variable name is *RBF\_PredictedValue*. If there are multiple dependent variables, then the default variable names are *RBF\_PredictedValue\_1*, *RBF\_PredictedValue\_2*, etc., corresponding to the order of the dependent variables on the command line.

### **PSEUDOPROB(rootname:n rootname...)**

*Predicted pseudo-probability.* If a dependent variable is categorical, then this keyword saves the predicted pseudo-probabilities of the first *n* categories of that dependent variable.

Specify one or more unique, valid variable names. There should be as many variable names specified as there are categorical dependent variables, and the names should be listed in the order of the categorical dependent variables on the command line. The specified names are treated as rootnames. Suffixes are added to each rootname to get a group of variable names corresponding to the categories for a given dependent variable. If you do not specify enough variable names, then default names are used for any remaining categorical dependent variables. If you specify too many variable names, then any remaining names are ignored.

A colon and a positive integer giving the number of probabilities to save for a dependent variable can follow the rootname.

If there is only one dependent variable, then the default rootname is *RBF\_PseudoProbability*. If there are multiple dependent variables, then the default rootnames are *RBF\_PseudoProbability\_1*, *RBF\_PseudoProbability\_2*, etc., corresponding to the order of the categorical dependent variables on the command line and taking into account the position of any scale dependent variables.

The default *n* is 25.

This keyword is ignored for scale dependent variables.

### ***Probabilities and Pseudo-Probabilities***

Predicted pseudo-probabilities cannot be interpreted as probabilities because the RBF procedure uses the sum-of-squares error and identity activation function for the output layer. The SAVE subcommand saves these predicted pseudo-probabilities even if any are less than 0 or greater than 1, or the sum for a given dependent variable is not 1.

The ROC, cumulative gains, and lift charts (see `/PLOT ROC, GAIN, and LIFT`, respectively) are created based on pseudo-probabilities. In the event that any of the pseudo-probabilities are less than 0 or greater than 1, or the sum for a given variable is not 1, they are first rescaled to be between 0 and 1 and to sum to 1. The `SAVE` subcommand saves the original pseudo-probabilities, but the charts are based on rescaled pseudo-probabilities.

Pseudo-probabilities are rescaled by dividing by their sum. For example, if a case has predicted pseudo-probabilities of 0.50, 0.60, and 0.40 for a three-category dependent variable, then each pseudo-probability is divided by the sum 1.50 to get 0.33, 0.40, and 0.27.

If any of the pseudo-probabilities are negative, then the absolute value of the lowest is added to all pseudo-probabilities before the above rescaling. For example, if the pseudo-probabilities are -0.30, 0.50, and 1.30, then first add 0.30 to each value to get 0.00, 0.80, and 1.60. Next, divide each new value by the sum 2.40 to get 0.00, 0.33, and 0.67.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand saves XML-format (PMML) files containing the hidden unit center and width vectors, and the synaptic weight estimates. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#)

- Filenames must be specified in full. The `RBF` procedure does not supply extensions.
- The `MODEL` keyword is not honored if split-file processing is in effect (see [SPLIT FILE](#) ). If this keyword is specified when split-file processing is on, then a warning is displayed.

**MODEL = 'file' 'file'...**

*Writes the hidden unit center and width vectors and synaptic weight estimates to XML (PMML) files. Specify one or more unique, valid filenames. There should be as many filenames as there are dependent variables, and the names should be listed in the order of the dependent variables on the command line. If you do not specify enough filenames, then an error is issued. If you specify too many filenames, then any remaining names are ignored.*

If any 'file' specification refers to an existing file, then the file is overwritten. If any 'file' specifications refer to the same file, then only the last instance of this 'file' specification is honored.

# READ MODEL

```
READ MODEL FILE='filename'  
  
  [/KEEP={ALL**      }]  
        {model names}  
        {procedures }  
  
  [/DROP={model names}]  
        {procedures }  
  
  [/TYPE={MODEL**}]  
        {COMMAND}  
  
  [/TSET={CURRENT**}]  
        {RESTORE  }
```

\*\*Default if the subcommand is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
READ MODEL FILE='ACFMOD.DAT' .
```

## Overview

READ MODEL reads a model file that has been previously saved on the SAVE MODEL command (see SAVE MODEL).

### Options

You can restore a subset of models from the model file by using the DROP and KEEP subcommands. You can use the TYPE subcommand to control whether models are specified by model name or by the name of the procedure that generated them. With the TSET subcommand, you can restore the TSET settings that were in effect when the model file was created.

### Basic Specification

The basic specification is the FILE subcommand, specifying the name of a previously saved model file.

- By default, all models that are contained in the specified file are restored, replacing all models that are currently active. The restored models have their original *MOD\_n* default names or names that are assigned by the MODEL NAME command.

### Subcommand Order

- Subcommands can be specified in any order.

**Syntax Rules**

- If a subcommand is specified more than once, only the last subcommand is executed.

**Operations**

- READ MODEL is executed immediately.
- Models that are currently active are erased when READ MODEL is executed. To save these models for later use, specify the SAVE MODEL command before READ MODEL.
- Model files are designed to be read by specific procedures and should not be edited.
- DATE specifications are not saved in model files. Therefore, the DATE specifications from the current session are applied to the restored models.
- The following procedures can generate models that can be read by the READ MODEL command: AREG, ARIMA, EXSMOOTH, SEASON, and SPECTRA in the Forecasting module; ACF, CASEPLOT, CCF, CURVEFIT, PACF, PLOT, and TSLOT in the Core system; and WLS and 2SLS in the Regression module.

**Limitations**

- A maximum of one filename can be specified.
- The READ MODEL command can only read models created by a SAVE MODEL command on the same operating system/platform.

**Example**

```
READ MODEL FILE='ACFMOD.DAT'  
/DROP=MOD_1.
```

- In this example, all models are restored except MOD\_1 in the model file ACFMOD.DAT.

**FILE Subcommand**

FILE names the model file to be read and is the only required subcommand.

- The only specification on FILE is the name of the model file.
- The filename should be enclosed in quotes.
- Only one filename can be specified.
- Only files that are saved with the SAVE MODEL command can be read.
- You can specify files residing in other directories by supplying a fully qualified filename.

**KEEP and DROP Subcommands**

KEEP and DROP allow you to restore a subset of models. By default, all models in the model file are restored.

- KEEP specifies the models to be restored.
- DROP specifies the models to be excluded.



- Models can be specified by using individual model names or the names of the procedures that created them. To use procedure names, you must specify `COMMAND` on the `TYPE` subcommand.
- Model names are either the default `MOD_n` names or the names that are assigned with `MODEL NAME`.
- If a procedure name is specified on `KEEP`, all models that are created by that procedure are restored; on `DROP`, all models that are created by the procedure are dropped.
- Model names and procedure names cannot be mixed on a single `READ MODEL` command.
- If more than one `KEEP` or `DROP` subcommand is specified, only the last subcommand is executed.
- You can specify the keyword `ALL` on `KEEP` to restore all models in the model file. This setting is the default.
- The stored model file is not affected by the `KEEP` or `DROP` specification on `READ MODEL`.

### Example

```
READ MODEL FILE='ACFCCF.DAT'  
/KEEP=ACF1 ACF2.
```

- In this example, only models `ACF1` and `ACF2` are restored from model file `ACFCCF.DAT`.

## TYPE Subcommand

`TYPE` indicates whether models are specified by model name or procedure name on `DROP` and `KEEP`.

- One keyword, `MODEL` or `COMMAND`, can be specified after `TYPE`.
- `MODEL` is the default and indicates that models are specified as model names.
- `COMMAND` indicates that models are specified by procedure name.
- `TYPE` has no effect if `KEEP` or `DROP` is not specified.
- The `TYPE` specification applies only to the current `READ MODEL` command.

### Example

```
READ MODEL FILE='CURVE1.DAT'  
/KEEP=CURVEFIT  
/TYPE=COMMAND.
```

- In this example, all models that are created by `CURVEFIT` are restored from model file `CURVE1.DAT`.

## TSET Subcommand

`TSET` allows you to restore the `TSET` settings that were in effect when the model was created.

- The specification on `TSET` is either `CURRENT` or `RESTORE`.

- `CURRENT` (the default) indicates that you want to continue to use the current `TSET` settings.
- `RESTORE` indicates that you want to restore the `TSET` settings that were in effect when the model file was saved. The current `TSET` settings are replaced with the model file settings when the file is restored.

# RECODE

*For numeric variables:*

```
RECODE varlist (value list=value)...(value list=value) [INTO varlist]
[/varlist...]
```

*Input keywords:*

LO, LOWEST, HI, HIGHEST, THRU, MISSING, SYSMIS, ELSE

*Output keywords:*

COPY, SYSMIS

*For string variables:*

```
RECODE varlist [('string', ['string'...]='string')][INTO varlist]
[/varlist...]
```

*Input keywords:*

CONVERT, ELSE

*Output keyword:*

COPY

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Examples**

```
RECODE V1 TO V3 (0=1) (1=0) (2,3=-1) (9=9) (ELSE=SYSMIS).
```

```
RECODE STRNGVAR ('A', 'B', 'C'='A') ('D', 'E', 'F'='B') (ELSE=' ').
```

## **Overview**

RECODE changes, rearranges, or consolidates the values of an existing variable. RECODE can be executed on a value-by-value basis or for a range of values. Where it can be used, RECODE is much more efficient than the series of IF commands that produce the same transformation.

With RECODE, you must specify the new values. Use AUTORECODE to automatically recode the values of string or numeric variables to consecutive integers.

### **Options**

You can generate a new variable as the recoded version of an existing variable using the keyword `INTO`. You can also use `INTO` to recode a string variable into a new numeric variable for more efficient processing, or to recode a numeric variable into a new string variable to provide more descriptive values.

### **Basic Specification**

The basic specification is a variable name and, within parentheses, the original values followed by a required equals sign and a new value. `RECODE` changes the values to the left of the equals sign into the single value to the right of the equals sign.

## **Syntax Rules**

- The variables to be recoded must already exist and must be specified before the value specifications.
- Value specifications are enclosed in parentheses. The original value or values must be specified to the left of an equals sign. A single new value is specified to the right of the equals sign.
- Multiple values can be consolidated into a single recoded value by specifying, to the left of the equals sign, a list of values separated by blanks or commas. Only one recoded value per set is allowed to the right of the equals sign.
- Multiple sets of value specifications are permitted. Each set must be enclosed in parentheses and can result in only one new value.
- To recode multiple variables using the same set of value specifications, specify a variable list before the value specifications. Each variable in the list is recoded identically.
- To recode variables using different value specifications, separate each variable (or variable list) and its specifications from the others by a slash.
- Original values that are not specified remain unchanged unless the keyword `ELSE` or `INTO` is used to recode into a new variable. `ELSE` refers to all original values not previously mentioned, including the system-missing value. `ELSE` should be the last specification for the variable. When recoding `INTO` another variable, unspecified values are set to system-missing or blank for strings.
- `COPY` replicates original values without recoding them.
- `INTO` is required to recode a string variable into a numeric variable or a numeric variable into a string variable.

## **Numeric Variables**

- Keywords that can be used in the list of original values are `LO` (or `LOWEST`), `HI` (or `HIGHEST`), `THRU`, `MISSING`, `YSMIS`, and `ELSE`. Keywords that can be used in place of a new value are `COPY` and `YSMIS`.
- `THRU` specifies a value range and includes the specified end values.
- `LOWEST` and `HIGHEST` (`LO` and `HI`) specify the lowest and highest values encountered in the data. `LOWEST` and `HIGHEST` include user-missing values but not the system-missing value.

- `MISSING` specifies user-missing and system-missing values for recoding. `MISSING` can be used in the list of original values only.
- `SYSMIS` specifies the system-missing value and can be used as both an original value and a new value.

### **String Variables**

- Keywords that can be used in the list of original values are `CONVERT` and `ELSE`. The only keyword that can be used in place of a new value is `COPY`.
- Values must be enclosed in quotes.
- Blanks are significant characters.

### **Operations**

- Value specifications are scanned left to right.
- A value is recoded only once per `RECODE` command.
- Invalid specifications on a `RECODE` command that result in errors stop all processing of that `RECODE` command. No variables are recoded.

### **Numeric Variables**

- Blank fields for numeric variables are handled according to the `SET BLANKS` specification prior to recoding.
- When you recode a value that was previously defined as user-missing on the `MISSING VALUE` command, the new value is not missing.

### **String Variables**

- If the original or new value specified is shorter than the format width defined for the variable, the string is right-padded with blanks.
- If the original or recoded value specified is longer than the format width defined for that variable, the program issues an error message and `RECODE` is not executed.

## **Examples**

### **Recoding Numeric Variables**

```
RECODE V1 TO V3 (0=1) (1=0) (2,3=-1) (9=9) (ELSE=SYSMIS)
/QVAR(1 THRU 5=1) (6 THRU 10=2) (11 THRU HI=3) (ELSE=0) .
```

- The numeric variables between and including `V1` and `V3` are recoded: original values 0 and 1 are switched respectively to 1 and 0; 2 and 3 are changed to -1; 9 remains 9; and any other value is changed to the system-missing value.
- Variable `QVAR` is also recoded: original values 1 through 5 are changed to 1; 6 through 10 are changed to 2; 11 through the highest value in the data are changed to 3; and any other value, including system-missing, is changed to 0.

**Recoding String Variables**

```
RECODE STRNGVAR ('A', 'B', 'C'='A') ('D', 'E', 'F'='B') (ELSE=' ').
RECODE PET ('IGUANA', 'SNAKE' = 'WILD ').
```

- Values A, B, and C are changed to value A. Values D, E, and F are changed to value B. All other values are changed to a blank.
- Values IGUANA and SNAKE are changed to value WILD. The defined width of the variable *PET* is 6. Thus, values SNAKE and WILD include trailing blanks for a total of six characters. If blanks are not specified, the values are right-padded. In this example, the results will be the same.
- Each string value is enclosed within quotes.

**INTO Keyword**

*INTO* specifies a **target** variable to receive recoded values from the original, or **source**, variable. Source variables remain unchanged after being recoded.

- *INTO* must follow the value specifications for the source variables that are being recoded into the target variables.
- The number of target variables must equal the number of source variables.

**Numeric Variables**

- Target variables can be existing or new variables. For existing variables, cases with values not mentioned in the value specifications are not changed. For new variables, cases with values not mentioned are assigned the system-missing value.
- New numeric variables have default print and write formats of F8.2 (or the format specified on *SET FORMAT*).

**Recoding a Single Variable Into a Target Variable**

```
RECODE AGE (MISSING=9) (18 THRU HI=1) (0 THRU 18=0) INTO VOTER.
```

- The recoded *AGE* values are stored in target variable *VOTER*, leaving *AGE* unchanged.
- Value 18 and higher values are changed to value 1. Values between 0 and 18, but not including 18, are recoded to 0. If the specification 0 THRU 18 preceded the specification 18 THRU HI, value 18 would be recoded to 0.

**Recording Multiple Variables Into Target Variables**

```
RECODE V1 TO V3 (0=1) (1=0) (2=-1) INTO DEFENSE WELFARE HEALTH.
```

- Values for *V1* through *V3* are recoded and stored in *DEFENSE*, *WELFARE*, and *HEALTH*. *V1*, *V2*, and *V3* are not changed.

## String Variables

- Target variables must already exist. To create a new string variable, declare the variable with the `STRING` command before specifying it on `RECODE`.
- The new string values cannot be longer than the defined width of the target variable.
- If the new values are shorter than the defined width of the target variable, the values are right-padded with blanks.
- Multiple target variables are allowed. The target variables must all be the same defined width; the source variables can have different widths.
- If the source and target variables have different widths, the criterion for the width of the original values is the width defined for the source variable; the criterion for the width of the recoded values is the width defined for the target variable.

### Using Keyword *COPY* With Target Variables

```
STRING STATE1 (A2).
RECODE STATE ('IO'='IA') (ELSE=COPY) INTO STATE1.
```

- `STRING` declares the variable `STATE1` so that it can be used as a target variable on `RECODE`.
- `RECODE` specifies `STATE` as the source variable and `STATE1` as the target variable. The original value `IO` is recoded to `IA`. The keywords `ELSE` and `COPY` copy all other state codes over unchanged. Thus, `STATE` and `STATE1` are identical except for cases with the original value `IO`.

### Recoding a String Variable Into a Numeric Target

```
RECODE SEX ('M'=1) ('F'=2) INTO NSEX.
```

- `RECODE` recodes the string variable `SEX` into the numeric variable `NSEX`. Any value other than `M` or `F` becomes system-missing.
- The program can process a large number of cases more efficiently with the numeric variable `NSEX` than it can with the string variable `SEX`.

## CONVERT Keyword

`CONVERT` recodes the string representation of numbers to their numeric representation.

- If the keyword `CONVERT` precedes the value specifications, cases with numbers are recoded immediately and blanks are recoded to the system-missing value, even if you specifically recode blanks into a value.
- To recode blanks to a value other than system-missing or to recode a string value to a noncorresponding numeric value (for example, '0' to 10), you must specify a recode specification *before* the keyword `CONVERT`.
- `RECODE` converts numbers as if the variable were being reread using the `F` format.
- If `RECODE` encounters a value that cannot be converted, it scans the remaining value specifications. If there is no specific recode specification for that value, the target variable will be system-missing for that case.

**Examples**

```
RECODE #JOB (CONVERT) ('-'=11) ('&'=12) INTO JOB.
```

- RECODE first recodes all numbers in the string variable *#JOB* to numbers. The target variable is *JOB*.
- RECODE then specifically recodes the minus sign (the “eleven” punch) to 11 and the ampersand (or “twelve” punch in EBCDIC) to 12. The keyword `CONVERT` is specified first as an efficiency measure to recode cases with numbers immediately. Blanks are recoded to the system-missing value.

```
RECODE #JOB (' '=-99) (CONVERT) ('-'=11) ('&'=12) INTO JOB.
```

- The result is the same as in the above example except that blanks are changed to `-99`.



# RECORD TYPE

*For mixed file types:*

```
RECORD TYPE {value list} [SKIP]
             {OTHER      }
```

*For grouped file types:*

```
RECORD TYPE {value list} [SKIP] [CASE=col loc]
             {OTHER      }

[DUPLICATE={WARN  }] [MISSING={WARN  }]
             {NOWARN}             {NOWARN}
```

*For nested file types:*

```
RECORD TYPE {value list} [SKIP] [CASE=col loc]
             {OTHER      }

[SPREAD={YES}] [MISSING={WARN  }]
             {NO  }             {NOWARN}
```

## **Example**

```
FILE TYPE MIXED RECORD=RECID 1-2.
RECORD TYPE 23.
DATA LIST /SEX 5 AGE 6-7 DOSAGE 8-10 RESULT 12.
END FILE TYPE.
```

## **Overview**

RECORD TYPE is used with DATA LIST within a FILE TYPE–END FILE TYPE structure to define any one of the three types of complex raw data files: **mixed files**, which contain several types of records that define different types of cases; **hierarchical** or **nested files**, which contain several types of records with a defined relationship among the record types; or **grouped files**, which contain several records for each case with some records missing or duplicated (see FILE TYPE for more complete information). A fourth type of complex file, files with **repeating groups** of information, can be read with the REPEATING DATA command. REPEATING DATA can also be used to read mixed files and the lowest level of nested files.

Each type of complex file has varying types of records. One set of RECORD TYPE and DATA LIST commands is used to define each type of record in the data. The specifications available for RECORD TYPE vary according to whether MIXED, GROUPED, or NESTED is specified on FILE TYPE.

### **Basic Specification**

For each record type being defined, the basic specification is the value of the record type variable defined on the RECORD subcommand on FILE TYPE.

- RECORD TYPE must be followed by a DATA LIST command defining the variables for the specified records, unless SKIP is used.

- One pair of `RECORD TYPE` and `DATA LIST` commands must be used for each defined record type.

### **Syntax Rules**

- A list of values can be specified if a set of different record types has the same variable definitions. Each value must be separated by a space or comma.
- String values must be enclosed in quotes.
- For mixed files, each `DATA LIST` can specify variables with the same variable name, since each record type defines a separate case. For grouped and nested files, the variable names on each `DATA LIST` must be unique, since a case is built by combining all record types together onto a single record.
- For mixed files, if the same variable is defined for more than one record type, the format type and width of the variable should be the same on all `DATA LIST` commands. The program refers to the first `DATA LIST` command that defines a variable for the print and write formats to include in the dictionary of the active dataset.
- For nested files, the order of the `RECORD TYPE` commands defines the hierarchical structure of the file. The first `RECORD TYPE` defines the highest-level record type, the next `RECORD TYPE` defines the next highest-level record, and so forth. The last `RECORD TYPE` command defines a case in the active dataset.

### **Operations**

- If a record type is specified on more than one `RECORD TYPE` command, the program uses the `DATA LIST` command associated with the first specification and ignores all others.
- For `NESTED` files, the first record in the file should be the type specified on the first `RECORD TYPE` command—the highest-level record of the hierarchy. If the first record in the file is not the highest-level type, the program skips all records until it encounters a record of the highest-level type. If the `MISSING` or `DUPLICATE` subcommands have been specified on the `FILE TYPE` command, these records may produce warning messages but will not be used to build a case in the active dataset.

## **Examples**

### **Reading a Single Record Type From a Mixed File**

```
FILE TYPE MIXED RECORD=RECID 1-2.
RECORD TYPE 23.
DATA LIST /SEX 5 AGE 6-7 DOSAGE 8-10 RESULT 12.
END FILE TYPE.
```

```
BEGIN DATA
21 145010 1
22 257200 2
25 235 250 2
35 167 300 3
24 125150 1
23 272075 1
21 149050 2
25 134 035 3
30 138 300 3
```

```
32 229          500    3
END DATA.
```

- FILE TYPE begins the file definition, and END FILE TYPE indicates the end of file definition. FILE TYPE specifies a mixed file type. Since the data are included between BEGIN DATA–END DATA, the FILE subcommand is omitted. The record identification variable *RECID* is located in columns 1 and 2.
- RECORD TYPE indicates that records with value 23 for variable *RECID* will be copied into the active dataset. All other records are skipped. The program does not issue a warning when it skips records in mixed files.
- DATA LIST defines variables on records with the value 23 for variable *RECID*.

### Reading Multiple Record Types From a Mixed File

```
FILE TYPE MIXED FILE=TREATMNT RECORD=RECID 1-2.
+ RECORD TYPE 21,22,23,24.
+ DATA LIST /SEX 5 AGE 6-7 DOSAGE 8-10 RESULT 12.
+ RECORD TYPE 25.
+ DATA LIST /SEX 5 AGE 6-7 DOSAGE 10-12 RESULT 15.
END FILE TYPE.
```

- Variable *DOSAGE* is read from columns 8–10 for record types 21, 22, 23, and 24 and from columns 10–12 for record type 25. *RESULT* is read from column 12 for record types 21, 22, 23, and 24 and from column 15 for record type 25.
- The active dataset contains values for all variables defined on the DATA LIST commands for record types 21 through 25. All other record types are skipped.

### Working With Nested Files

\* A nested file of accident records.

```
FILE TYPE NESTED RECORD=6 CASE=ACCID 1-4.
RECORD TYPE 1.
DATA LIST /ACC_ID 9-11 WEATHER 12-13 STATE 15-16 (A) DATE 18-24 (A).
RECORD TYPE 2.
DATA LIST /STYLE 11 MAKE 13 OLD 14 LICENSE 15-16(A) INSURNCE 18-21 (A).
RECORD TYPE 3.
DATA LIST /PSNGR_NO 11 AGE 13-14 SEX 16 (A) INJURY 18 SEAT 20-21 (A)
          COST 23-24.
END FILE TYPE.
```

```
BEGIN DATA
0001 1 322 1 IL 3/13/88 /* Type 1:  accident record
0001 2   1 44MI 134M   /* Type 2:   vehicle record
0001 3   1 34 M 1 FR 3 /* Type 3:   person record
0001 2   2 16IL 322F   /*           vehicle record
0001 3   1 22 F 1 FR 11 /*           person record
0001 3   2 35 M 1 FR 5 /*           person record
0001 3   3 59 M 1 BK 7 /*           person record
0001 2   3 21IN 146M  /*           vehicle record
0001 3   1 46 M 0 FR 0 /*           person record
END DATA.
```

- FILE TYPE specifies a nested file type. The record identifier, located in column 6, is not assigned a variable name, so the default scratch variable name #####RECD is used. The case identification variable *ACCID* is located in columns 1–4.

- Because there are three record types, there are three RECORD TYPE commands. For each RECORD TYPE, there is a DATA LIST command to define variables on that record type. The order of the RECORD TYPE commands defines the hierarchical structure of the file.
- END FILE TYPE signals the end of file definition.
- The program builds a case for each lowest-level (type 3) record, representing each person in the file. There can be only one type 1 record for each type 2 record, and one type 2 record for each type 3 record. Each vehicle can be in only one accident, and each person can be in only one vehicle. The variables from the type 1 and type 2 records are spread to their corresponding type 3 records.

## ***OTHER Keyword***

OTHER specifies all record types that have not been mentioned on previous RECORD TYPE commands.

- OTHER can be specified only on the last RECORD TYPE command in the file definition.
- OTHER can be used with SKIP to skip all undefined record types.
- For nested files, OTHER can be used only with SKIP. Neither can be used separately.
- If WILD=WARN is in effect for the FILE TYPE command, OTHER cannot be specified on the RECORD TYPE command.

### ***Using Keyword OTHER With a Mixed File***

```
FILE TYPE MIXED FILE=TREATMNT RECORD=RECID 1-2.
RECORD TYPE 21,22,23,24.
DATA LIST /SEX 5 AGE 6-7 DOSAGE 8-10 RESULT 12.
RECORD TYPE 25.
DATA LIST /SEX 5 AGE 6-7 DOSAGE 10-12 RESULT 15.
RECORD TYPE OTHER.
DATA LIST /SEX 5 AGE 6-7 DOSAGE 18-20 RESULT 25.
END FILE TYPE.
```

- The first two RECORD TYPE commands specify record types 21–25. All other record types are specified by the third RECORD TYPE.

### ***Using Keyword OTHER With a Nested File***

```
FILE TYPE NESTED FILE=ACCIDENT RECORD=#RECID 6 CASE=ACCID 1-4.
RECORD TYPE 1. /* Accident record
DATA LIST /WEATHER 12-13.
RECORD TYPE 2. /* Vehicle record
DATA LIST /STYLE 16.
RECORD TYPE OTHER SKIP.
END FILE TYPE.
```

- The third RECORD TYPE specifies OTHER SKIP. Type 2 records are therefore the lowest-level records included in the active dataset. These commands build one case for each vehicle record. The person records are skipped.
- Because the data are in a nested file, OTHER can be specified only with SKIP.

## SKIP Subcommand

SKIP specifies record types to skip.

- To skip selected record types, specify the values for the types you want to skip and then specify SKIP. To skip all record types other than those specified on previous RECORD TYPE commands, specify OTHER and then SKIP.
- For nested files, SKIP can be used only with OTHER. Neither can be used separately.
- For grouped files, OTHER cannot be specified on SKIP if WILD=WARN (the default) is in effect for FILE TYPE.
- For mixed files, all record types that are not specified on a RECORD TYPE command are skipped by default. No warning is issued (WILD=NOWARN on FILE TYPE is the default for mixed files).
- For grouped files, a warning message is issued by default for all record types not specified on a RECORD TYPE command (WILD=WARN on FILE TYPE is the default for grouped files). If the record types are explicitly specified on SKIP, no warning is issued.

### Examples

```
FILE TYPE GROUPED FILE=HUBDATA RECORD=#RECID 80 CASE=ID 1-5
                                WILD=NOWARN.
RECORD TYPE 1.
DATA LIST /MOHIRED YRHIRED 12-15 DEPT79 TO DEPT82 SEX 16-20.
RECORD TYPE OTHER SKIP.
END FILE TYPE.
```

- The program reads variables from type 1 records and skips all other types.
- WILD=NOWARN on the FILE TYPE command suppresses the warning messages that is issued by default for undefined record types for grouped files. Keyword OTHER cannot be used when the default WILD=WARN specification is in effect.

```
FILE TYPE GROUPED FILE=HUBDATA RECORD=#RECID 80 CASE=ID 1-5.
RECORD TYPE 1.
DATA LIST /MOHIRED YRHIRED 12-15 DEPT79 TO DEPT82 SEX 16-20.
RECORD TYPE 2,3 SKIP.
END FILE TYPE.
```

- Record type 1 is defined for each case, and record types 2 and 3 are skipped.
- WILD=WARN (the default) on FILE TYPE GROUPED is in effect. The program therefore issues a warning message for any record types it encounters other than types 1, 2, and 3. No warning is issued for record types 2 and 3 because they are explicitly specified on a RECORD TYPE command.

## CASE Subcommand

CASE specifies the column locations of the case identification variable when that variable is not in the location defined by the CASE subcommand on FILE TYPE.

- CASE on RECORD TYPE applies only to those records specified by that RECORD TYPE command. The identifier for record types without CASE on RECORD TYPE must be in the location specified by CASE on FILE TYPE.

---

*RECORD TYPE*

- CASE can be used for nested and grouped files only. CASE cannot be used for mixed files.
- CASE can be used on RECORD TYPE only if a CASE subcommand is specified on FILE TYPE.
- The format type of the case identification variable must be the same on all records, and the same format must be assigned on the RECORD TYPE and FILE TYPE commands. For example, if the case identification variable is defined as a string on FILE TYPE, it cannot be defined as a numeric variable on RECORD TYPE.

**Example**

\* Specifying case on the record type command for a grouped file.

```
FILE TYPE GROUPED FILE=HUBDATA RECORD=#RECID 80 CASE=ID 1-5.
RECORD TYPE 1.
DATA LIST /MOHIRED YRHIRED 12-15 DEPT79 TO DEPT82 SEX 16-20.
RECORD TYPE 2.
DATA LIST /SALARY79 TO SALARY82 6-25
          HOURLY81 HOURLY82 40-53 (2)
          PROMO81 72 AGE 54-55 RAISE82 66-70.
RECORD TYPE 3 CASE=75-79.
DATA LIST /JOB CAT 6 NAME 25-48 (A).
END FILE TYPE.
```

- CASE on FILE TYPE indicates that the case identification variable is located in columns 1–5. On the third RECORD TYPE command, the CASE subcommand overrides the identifier location for type 3 records. For type 3 records, the case identification variable is located in columns 75–79.

**MISSING Subcommand**

MISSING controls whether the program issues a warning when it encounters a missing record type for a case. Regardless of whether the program issues the warning, it builds the case in the active dataset with system-missing values for the variables defined on the missing record.

- The only specification is a single keyword. NOWARN is the default for nested files. WARN is the default for grouped files. MISSING cannot be used with MIXED files.
- MISSING on RECORD TYPE applies only to those records specified by that RECORD TYPE command. The treatment of missing records for record types without the MISSING specification on RECORD TYPE is determined by the MISSING subcommand on FILE TYPE.
- For grouped files, the program checks whether there is a record for each case identification number. For nested files, the program verifies that each defined case includes one record of each type.

**WARN**            *Issue a warning message when a record type is missing for a case.* This is the default for grouped files.

**NOWARN**        *Suppress the warning message when a record type is missing for a case.* This is the default for nested files.

**Example**

```
FILE TYPE GROUPED FILE=HUBDATA RECORD=#RECID 80 CASE=ID 1-5.
RECORD TYPE 1.
DATA LIST /MOHIRED YRHIRED 12-15 DEPT79 TO DEPT82 SEX 16-20.
```

```

RECORD TYPE 2 MISSING=NOWARN.
DATA LIST /SALARY79 TO SALARY82 6-25
    HOURLY81 HOURLY82 40-53 (2) PROMO81 72 AGE 54-55 RAISE82 66-70.
RECORD TYPE 3.
DATA LIST /JOB CAT 6 NAME 25-48 (A).
END FILE TYPE.

```

- MISSING is not specified on FILE TYPE. Therefore the default MISSING=WARN is in effect for all record types.
- MISSING=NOWARN is specified on the second RECORD TYPE, overriding the default setting for type 2 records. WARN is still in effect for type 1 and type 3 records.

## ***DUPLICATE Subcommand***

DUPLICATE controls whether the program issues a warning when it encounters more than one record of each type for a single case.

- DUPLICATE on RECORD TYPE can be used for grouped files only. DUPLICATE cannot be used for mixed or nested files.
- The only specification is a single keyword. WARN is the default.
- DUPLICATE on RECORD TYPE applies only to those records specified by that RECORD TYPE command. The treatment of duplicate records for record types without DUPLICATE specification is determined by the DUPLICATE subcommand on FILE TYPE.
- Regardless of the specification on DUPLICATE, only the last record from a set of duplicates is included in the active dataset.

**WARN**            *Issue a warning message.* The program issues a message and the first 80 characters of the last record of the duplicate set of record types. This is the default.

**NOWARN**        *Suppress the warning message.*

### ***Example***

```

* Specifying DUPLICATE on RECORD TYPE for a grouped file.
FILE TYPE GROUPED FILE=HUBDATA RECORD=#RECID 80 CASE=ID 1-5.
RECORD TYPE 1.
DATA LIST /MOHIRED YRHIRED 12-15 DEPT79 TO DEPT82 SEX 16-20.
RECORD TYPE 2 DUPLICATE=NOWARN.
DATA LIST /SALARY79 TO SALARY82 6-25
    HOURLY81 HOURLY82 40-53 (2) PROMO81 72 AGE 54-55 RAISE82 66-70.
RECORD TYPE 3.
DATA LIST /JOB CAT 6 NAME 25-48 (A).
END FILE TYPE.

```

- DUPLICATE is not specified on FILE TYPE. Therefore the default DUPLICATE=WARN is in effect for all record types.
- DUPLICATE=NOWARN is specified on the second RECORD TYPE, overriding the FILE TYPE setting for type 2 records. WARN is still in effect for type 1 and type 3 records.

## SPREAD Subcommand

SPREAD controls whether the values for variables defined for a record type are spread to all related cases.

- SPREAD can be used for nested files only. SPREAD cannot be used for mixed or grouped files.
- The only specification is a single keyword. YES is the default.
- SPREAD=NO applies only to the record type specified on that RECORD TYPE command. The default YES is in effect for all other defined record types.

**YES** *Spread the values from the specified record type to all related cases.* This is the default.

**NO** *Spread the values from the specified type only to the first related case.* All other cases built from the same record are assigned the system-missing value for the variables defined on the record type.

### Example

\* A nested file.

```
FILE TYPE NESTED RECORD=#RECID 6 CASE=ACCID 1-4.
RECORD TYPE 1.
DATA LIST /ACC_NO 9-11 WEATHER 12-13
          STATE 15-16 (A) DATE 18-24 (A).
RECORD TYPE 2 SPREAD=NO.
DATA LIST /STYLE 11 MAKE 13 OLD 14
          LICENSE 15-16 (A) INSURNCE 18-21 (A).
RECORD TYPE 3.
DATA LIST /PSNGR_NO 11 AGE 13-14 SEX 16 (A)
          INJURY 18 SEAT 20-21 (A) COST 23-24.
END FILE TYPE.

BEGIN DATA
0001 1 322 1 IL 3/13/88 /* Type 1:  accident record
0001 2 1 44MI 134M /* Type 2:   vehicle record
0001 3 1 34 M 1 FR 3 /* Type 3:   person record
0001 2 2 16IL 322F /*           vehicle record
0001 3 1 22 F 1 FR 11 /*           person record
0001 3 2 35 M 1 FR 5 /*           person record
0001 3 3 59 M 1 BK 7 /*           person record
0001 2 3 21IN 146M /*           vehicle record
0001 3 1 46 M 0 FR 0 /*           person record
END DATA.
```

- The accident record (type 1) is spread to all related cases (in this example, all cases).
- The first vehicle record has one related person record. The values for *STYLE*, *MAKE*, *OLD*, *LICENSE*, and *INSURNCE* are spread to the case built for the person record.
- The second vehicle record has three related person records. The values for *STYLE*, *MAKE*, *OLD*, *LICENSE*, and *INSURNCE* are spread only to the case built from the first person record. The other two cases have the system-missing values for *STYLE*, *MAKE*, *OLD*, *LICENSE*, and *INSURNCE*.
- The third vehicle record has one related person record, and the values for type 2 records are spread to that case.



# ***REFORMAT***

This command is no longer supported.

# REGRESSION

REGRESSION is available in the Statistics Base option.

```
REGRESSION [MATRIX=[IN({file})] [OUT({file})]]
           { * } { * }

[/VARIABLES={varlist }
            { (COLLECT)** }
            { ALL } }

[/DESCRIPTIVES=[DEFAULTS] [MEAN] [STDEEV] [CORR] [COV]
               [VARIANCE] [XPROD] [SIG] [N] [BADCORR]
               [ALL] [NONE]**]

[/SELECT={varname relation value}

[/MISSING=[{LISTWISE** } ] [INCLUDE]]
          {PAIRWISE }
          {MEANSUBSTITUTION}

[/REGWGT=varname]

[/STATISTICS=[DEFAULTS**] [R**] [COEFF**] [ANOVA**] [OUTS**]
             [ZPP] [LABEL] [CHANGE] [CI({95 })] [F] [BCOV] [SES]
             {value}
             [XTX] [COLLIN] [TOL] [SELECTION] [ALL]]

[/CRITERIA=[DEFAULTS**] [TOLERANCE({0.0001**})] [MAXSTEPS(n)]
           {value }

           [PIN[({0.05**})] ] [POUT[({0.10**})] ]
           {value } {value }

           [FIN[({3.84 })] ] [FOUT[({2.71 })] ]
           {value } {value }

           [CIN[({ 95**})] ]
           {value }

[/{NOORIGIN**}]
 {ORIGIN }

/DEPENDENT=varlist

[/METHOD={STEPWISE [varlist] } [...] [/...]
         {FORWARD [varlist] }
         {BACKWARD [varlist] }
         {ENTER [varlist] }
         {REMOVE varlist }
         {TEST(varlist)(varlist)...}

[/RESIDUALS=[DEFAULTS] [DURBIN] [OUTLIERS({ZRESID })] [ID (varname)]
           {tempvars}

           [NORMPROB({ZRESID })] [HISTOGRAM({ZRESID })]
           {tempvars} {tempvars}

           [SIZE({SEPARATE})]
           {POOLED }

[/CASEWISE=[DEFAULTS] [{OUTLIERS({3 })}] [PLOT({ZRESID })]
             { {value} } {tempvar}
             {ALL }

             [{DEPENDENT PRED RESID}]
             {tempvars }

[/SCATTERPLOT [varname,varname]...]
```

```

[/PARTIALPLOT={ALL    }]
                {varlist}

[/OUTFILE={COVB ('savfile' | 'dataset')}}  [{MODEL('file'    )}]
                {CORB ('savfile' | 'dataset')}    {PARAMETER('file')}

[/SAVE=tempvar[(newname)]  [tempvar[(newname)]...]  [FITS]]

[/TEMPLATE='filename']

```

**\*\*Default if the subcommand is omitted.**

*Temporary residual variables are:*

*PRED, ADJPRED, SRESID, MAHAL, RESID, ZPRED, SDRESID, COOK, DRESID, ZRESID, SEPRED, LEVER, DFBETA, SDBETA, DFFIT, SDFFIT, COVRATIO, MCIN, ICIN*

*SAVE FITS saves:*

*DFFIT, SDFIT, DFBETA, SDBETA, COVRATIO*

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- PARAMETER keyword introduced on OUTFILE subcommand.

Release 16.0

- Added support for SET THREADS and SET MCACHE.

Release 17.0

- Added option to specify confidence level on CI keyword of STATISTICS subcommand.

Release 19.0

- TEMPLATE subcommand introduced.

### **Example**

```

REGRESSION VARIABLES=POP15 , POP75 , INCOME , GROWTH , SAVINGS
  /DEPENDENT=SAVINGS
  /METHOD=ENTER POP15 , POP75 , INCOME
  /METHOD=ENTER GROWTH .

```

## **Overview**

REGRESSION calculates multiple regression equations and associated statistics and plots. REGRESSION also calculates collinearity diagnostics, predicted values, residuals, measures of fit and influence, and several statistics based on these measures.

**Options**

**Input and Output Control Subcommands.** `DESCRIPTIVES` requests descriptive statistics on the variables in the analysis. `SELECT` estimates the model based on a subset of cases. `REGWGT` specifies a weight variable for estimating weighted least-squares models. `MISSING` specifies the treatment of cases with missing values. `MATRIX` reads and writes matrix data files.

**Equation-Control Subcommands.** These optional subcommands control the calculation and display of statistics for each equation. `STATISTICS` controls the statistics displayed for the equation(s) and the independent variable(s), `CRITERIA` specifies the criteria used by the variable selection method, and `ORIGIN` specifies whether regression is through the origin.

**Analysis of Residuals, Fit, and Influence.** `REGRESSION` creates temporary variables containing predicted values, residuals, measures of fit and influence, and several statistics based on these measures. These temporary variables can be analyzed within `REGRESSION` in Casewise Diagnostics tables (`CASEWISE` subcommand), scatterplots (`SCATTERPLOT` subcommand), histograms and normal probability plots (`RESIDUALS` subcommand), and partial regression plots (`PARTIALPLOT` subcommand). Any of the residuals subcommands can be specified to obtain descriptive statistics for the predicted values, residuals, and their standardized versions. Any of the temporary variables can be added to the active dataset with the `SAVE` subcommand.

**Templates.** You can specify a template, using the `TEMPLATE` subcommand, to override the default chart attribute settings on your system.

**Basic Specification**

The basic specification is `DEPENDENT`, which initiates the equation(s) and defines at least one dependent variable, followed by `METHOD`, which specifies the method for selecting independent variables.

- By default, all variables named on `DEPENDENT` and `METHOD` are used in the analysis.
- The default display for each equation includes a Model Summary table showing  $R^2$ , an ANOVA table, a Coefficients table displaying related statistics for variables in the equation, and an Excluded Variables table displaying related statistics for variables not yet in the equation.
- By default, all cases in the active dataset with valid values for all selected variables are used to compute the correlation matrix on which the regression equations are based. The default equations include a constant (intercept).
- All residuals analysis subcommands are optional. Most have defaults that can be requested by including the subcommand without any further specifications. These defaults are described in the discussion of each subcommand below.

**Subcommand Order**

The standard subcommand order for `REGRESSION` is

```
REGRESSION MATRIX=...
/VARIABLES=...
/DESCRIPTIVES=...
/SELECT=...
/MISSING=...
```

```

/REGWGT=...

      --Equation Block--
/STATISTICS=...
/CRITERIA=...
/ORIGIN
/DEPENDENT=...

      --Method Block(s)--
/METHOD=...
[/METHOD=...]

      --Residuals Block--
/RESIDUALS=...
/SAVE=...
/CASEWISE=...
/SCATTERPLOT=...
/PARTIALPLOT=...
/OUTFILE=...

```

- When used, `MATRIX` must be specified first.
- Subcommands listed before the equation block must be specified before any subcommands within the block.
- Only one equation block is allowed per `REGRESSION` command.
- An equation block can contain multiple `METHOD` subcommands. These methods are applied, one after the other, to the estimation of the equation for that block.
- The `STATISTICS`, `CRITERIA`, and `ORIGIN/NOORIGIN` subcommands must precede the `DEPENDENT` subcommand.
- The residuals subcommands `RESIDUALS`, `CASEWISE`, `SCATTERPLOT`, and `PARTIALPLOT` follow the last `METHOD` subcommand of any equation for which residuals analysis is requested. Statistics are based on this final equation.
- Residuals subcommands can be specified in any order. All residuals subcommands must follow the `DEPENDENT` and `METHOD` subcommands.

### **Syntax Rules**

- `VARIABLES` can be specified only once. If omitted, `VARIABLES` defaults to `COLLECT`.
- The `DEPENDENT` subcommand can be specified only once and must be followed immediately by one or more `METHOD` subcommands.
- `CRITERIA`, `STATISTICS`, and `ORIGIN` must be specified before `DEPENDENT` and `METHOD`. If any of these subcommands are specified more than once, only the last specified is in effect for all subsequent equations.
- More than one variable can be specified on the `DEPENDENT` subcommand. An equation is estimated for each.
- If no variables are specified on `METHOD`, all variables named on `VARIABLES` but not on `DEPENDENT` are considered for selection.

### **Operations**

- This procedure uses the multithreaded options specified by `SET THREADS` and `SET MCACHE`.

**Operations**

- REGRESSION calculates a correlation matrix that includes all variables named on VARIABLES. All equations requested on the REGRESSION command are calculated from the same correlation matrix.
- The MISSING, DESCRIPTIVES, and SELECT subcommands control the calculation of the correlation matrix and associated displays.
- If multiple METHOD subcommands are specified, they operate in sequence on the equations defined by the preceding DEPENDENT subcommand.
- Only independent variables that pass the tolerance criterion are candidates for entry into the equation. [For more information, see the topic CRITERIA Subcommand on p. 1711.](#)
- The temporary variables *PRED* (unstandardized predicted value), *ZPRED* (standardized predicted value), *RESID* (unstandardized residual), and *ZRESID* (standardized residual) are calculated and descriptive statistics are displayed whenever any residuals subcommand is specified. If any of the other temporary variables are referred to on the command, they are also calculated.
- Predicted values and statistics based on predicted values are calculated for every observation that has valid values for all variables in the equation. Residuals and statistics based on residuals are calculated for all observations that have a valid predicted value and a valid value for the dependent variable. The missing-values option therefore affects the calculation of residuals and predicted values.
- No residuals or predictors are generated for cases deleted from the active dataset with SELECT IF, a temporary SELECT IF, or SAMPLE.
- All variables are standardized before plotting. If the unstandardized version of a variable is requested, the standardized version is plotted.
- Residuals processing is not available when the active dataset is a matrix file or is replaced by a matrix file with MATRIX OUT (\*) on REGRESSION. If RESIDUALS, CASEWISE, SCATTERPLOT, PARTIALPLOT, or SAVE are used when MATRIX IN (\*) or MATRIX OUT (\*) is specified, the REGRESSION command is not executed.

For each analysis, REGRESSION can calculate the following types of temporary variables:

<b>PRED</b>	<i>Unstandardized predicted values.</i>
<b>RESID</b>	<i>Unstandardized residuals.</i>
<b>DRESID</b>	<i>Deleted residuals.</i>
<b>ADJPRED</b>	<i>Adjusted predicted values.</i>
<b>ZPRED</b>	<i>Standardized predicted values.</i>
<b>ZRESID</b>	<i>Standardized residuals.</i>
<b>SRESID</b>	<i>Studentized residuals.</i>
<b>SDRESID</b>	<i>Studentized deleted residuals. (Hoaglin and Welsch, 1978)</i>
<b>SEPREP</b>	<i>Standard errors of the predicted values.</i>
<b>MAHAL</b>	<i>Mahalanobis distances.</i>
<b>COOK</b>	<i>Cook's distances. (Cook, 1977)</i>
<b>LEVER</b>	<i>Centered leverage values. (Velleman and Welsch, 1981)</i>

<b>DFBETA</b>	<i>Change in the regression coefficient that results from the deletion of the <i>i</i>th case. A DFBETA value is computed for each case for each regression coefficient generated by a model. (Belsley, Kuh, and Welsch, 1980)</i>
<b>SDBETA</b>	<i>Standardized DFBETA. An SDBETA value is computed for each case for each regression coefficient generated by a model. (Belsley et al., 1980)</i>
<b>DFFIT</b>	<i>Change in the predicted value when the <i>i</i>th case is deleted. (Belsley et al., 1980)</i>
<b>SDFIT</b>	<i>Standardized DFFIT. (Belsley et al., 1980)</i>
<b>COVRATIO</b>	<i>Ratio of the determinant of the covariance matrix with the <i>i</i>th case deleted to the determinant of the covariance matrix with all cases included. (Belsley et al., 1980)</i>
<b>MCIN</b>	<i>Lower and upper bounds for the prediction interval of the mean predicted response. A lowerbound LMCIN and an upperbound UMCIN are generated. The default confidence interval is 95%. The confidence interval can be reset with the CIN subcommand. (Dillon and Goldstein, 1984)</i>
<b>ICIN</b>	<i>Lower and upper bounds for the prediction interval for a single observation. A lowerbound LICIN and an upperbound UICIN are generated. The default confidence interval is 95%. The confidence interval can be reset with the CIN subcommand. (Dillon et al., 1984)</i>

## Examples

### Multiple METHOD subcommands

```
REGRESSION VARIABLES=POP15 , POP75 , INCOME , GROWTH , SAVINGS
  /DEPENDENT=SAVINGS
  /METHOD=ENTER POP15 , POP75 , INCOME
  /METHOD=ENTER GROWTH .
```

- VARIABLES calculates a correlation matrix of five variables for use by REGRESSION.
- DEPENDENT defines a single equation, with SAVINGS as the dependent variable.
- The first METHOD subcommand enters POP15, POP75, and INCOME into the equation.
- The second METHOD subcommand adds GROWTH to the equation containing POP15 to INCOME.

### Specifying Residual Output

```
REGRESSION
  /MISSING LISTWISE
  /DESCRIPTIVES MEAN STDDEV
  /STATISTICS COEFF R ANOVA
  /CRITERIA=PIN(.05) POUT(.10)
  /NOORIGIN
  /DEPENDENT time
  /METHOD=ENTER diam
  /PARTIALPLOT ALL
  /SCATTERPLOT=(*SDRESID , *ZPRED )
  /RESIDUALS HIST(ZRESID) NORM(ZRESID) ID( type )
  /CASEWISE PLOT(ZRESID) OUTLIERS(2)
  /SAVE ZPRED ZRESID COOK LEVER .
```

- The procedure creates a model for time using diam as a predictor.

- RESIDUALS requests a histogram and normal probability plot of the standardized residuals. Cases are labeled by values of *type*. Because residuals processing has been requested, statistics for predicted values, residuals, and standardized versions of predicted values and residuals are displayed in a Residuals Statistics table.
- CASEWISE requests a Casewise Diagnostics table for cases whose absolute value of *ZRESID* is greater than 2. Values of the studentized residual are listed for each case.
- SCATTERPLOT requests a plot of the standardized predicted value and the studentized residual.
- PARTIALPLOT requests partial regression plots for all independent variables.
- SAVE adds the standardized predicted value, standardized residual, Cook's distance, and leverage to the active dataset as new variables.
- All other options are set to their default values.

### Diagnosing multicollinearity

```
REGRESSION
  /MISSING LISTWISE
  /STATISTICS COEFF R ANOVA COLLIN ZPP TOL
  /CRITERIA=PIN(.05) POUT(.10)
  /NOORIGIN
  /DEPENDENT lnsales
  /METHOD=ENTER type price engine_s horsepower wheelbas width length curb_wgt
  fuel_cap mpg .
```

- The procedure creates a model for *lnsales* using *type*, *price*, *engine\_s*, *horsepower*, *wheelbas*, *width*, *length*, *curb\_wgt*, *fuel\_cap*, and *mpg* as predictors.
- STATISTICS requests coefficient estimates, multiple R statistics, an ANOVA table, collinearity diagnostics, part and partial correlations, and tolerance.
- All other options are set to their default values.

### Stepwise regression

```
FACTOR
  /VARIABLES ztype zprice zengine_ zhorsepo zwheelba zwidth zlength zcurb_wg zfuel_ca zmpg
  /MISSING LISTWISE
  /ANALYSIS ztype zprice zengine_ zhorsepo zwheelba zwidth zlength zcurb_wg zfuel_ca zmpg
  /PRINT INITIAL EXTRACTION ROTATION
  /CRITERIA FACTORS(10) ITERATE(25)
  /EXTRACTION PC
  /CRITERIA ITERATE(25)
  /ROTATION VARIMAX
  /SAVE REG(ALL)
  /METHOD=CORRELATION.
```

- The procedure creates a model for *lnsales* using *ztype*, *zprice*, *zengine\_*, *zhorsepo*, *zwheelba*, *zwidth*, *zlength*, *zcurb\_wg*, *zfuel\_ca*, and *zmpg* as predictors.
- METHOD specifies that predictors should be added to the model using stepwise selection.
- RESIDUALS requests a histogram and normal probability plot of the standardized residuals. Cases are labeled by values of *model*. Because residuals processing has been requested, statistics for predicted values, residuals, and standardized versions of predicted values and residuals are displayed in a Residuals Statistics table.



- CASEWISE requests a Casewise Diagnostics table for cases whose absolute value of *ZRESID* is greater than 2. Values of the standardized residual are listed for each case.
- SCATTERPLOT requests a plot of the standardized predicted value and the studentized residual.
- PARTIALPLOT requests partial regression plots for all independent variables.
- SAVE adds the standardized predicted value, standardized residual, Cook's distance, and leverage to the active dataset as new variables.
- All other options are set to their default values.

## VARIABLES Subcommand

VARIABLES names all the variables to be used in the analysis.

- The minimum specification is a list of two variables or the keyword ALL or COLLECT. COLLECT, which must be specified in parentheses, is the default.
- Only one VARIABLES subcommand is allowed, and it must precede any DEPENDENT or METHOD subcommands.
- You can use keyword TO to refer to consecutive variables in the active dataset.
- The order of variables in the correlation matrix constructed by REGRESSION is the same as their order on VARIABLES. If (COLLECT) is used, the order of variables in the correlation matrix is the order in which they are first listed on the DEPENDENT and METHOD subcommands.

**ALL** *Include all user-defined variables in the active dataset.*

**(COLLECT)** *Include all variables named on the DEPENDENT and METHOD subcommands. COLLECT is the default if the VARIABLES subcommand is omitted. COLLECT must be specified in parentheses. If COLLECT is used, the METHOD subcommands must specify variable lists.*

### Example

```
REGRESSION VARIABLES=(COLLECT)
  /DEPENDENT=SAVINGS
  /METHOD=STEP POP15 POP75 INCOME
  /METHOD=ENTER GROWTH.
```

- COLLECT requests that the correlation matrix include *SAVINGS*, *POP15*, *POP75*, *INCOME*, and *GROWTH*. Since COLLECT is the default, the VARIABLES subcommand could have been omitted.
- The DEPENDENT subcommand defines a single equation in which *SAVINGS* is the dependent variable.
- The first METHOD subcommand requests that the block of variables *POP15*, *POP75*, and *INCOME* be considered for inclusion using a stepwise procedure.
- The second METHOD subcommand adds variable *GROWTH* to the equation.

## DEPENDENT Subcommand

DEPENDENT specifies a list of variables and requests that an equation be built for each. DEPENDENT is required.

- The minimum specification is a single variable. There is no default variable list.
- Only one `DEPENDENT` subcommand can be specified. It must be followed by at least one `METHOD` subcommand.
- Keyword `TO` on a `DEPENDENT` subcommand refers to the order in which variables are specified on the `VARIABLES` subcommand. If `VARIABLES= (COLLECT)`, `TO` refers to the order of variables in the active dataset.
- If `DEPENDENT` names more than one variable, an equation is built for each using the same independent variables and methods.

## ***METHOD Subcommand***

`METHOD` specifies a variable selection method and names a block of variables to be evaluated using that method. `METHOD` is required.

- The minimum specification is a method keyword and, for some methods, a list of variables. The actual keyword `METHOD` can be omitted.
- When more than one `METHOD` subcommand is specified, each `METHOD` subcommand is applied to the equation that resulted from the previous `METHOD` subcommands.
- The default variable list for methods `FORWARD`, `BACKWARD`, `STEPWISE`, and `ENTER` consists of all variables named on `VARIABLES` that are not named on the `DEPENDENT` subcommand. If `VARIABLES= (COLLECT)`, the variables must be specified for these methods.
- There is no default variable list for the `REMOVE` and `TEST` methods.
- Keyword `TO` in a variable list on `METHOD` refers to the order in which variables are specified on the `VARIABLES` subcommand. If `VARIABLES= (COLLECT)`, `TO` refers to the order of variables in the active dataset.

The available stepwise methods are as follows:

<b>BACKWARD [varlist]</b>	<i>Backward elimination.</i> Variables in the block are considered for removal. At each step, the variable with the largest probability-of- $F$ value is removed, provided that the value is larger than <code>POUT</code> . <a href="#">For more information, see the topic CRITERIA Subcommand on p. 1711.</a> If no variables are in the equation when <code>BACKWARD</code> is specified, all independent variables in the block are first entered.
<b>FORWARD [varlist]</b>	<i>Forward entry.</i> Variables in the block are added to the equation one at a time. At each step, the variable not in the equation with the smallest probability of $F$ is entered if the value is smaller than <code>PIN</code> . <a href="#">For more information, see the topic CRITERIA Subcommand on p. 1711.</a>
<b>STEPWISE [varlist]</b>	<i>Stepwise selection.</i> If there are independent variables already in the equation, the variable with the largest probability of $F$ is removed if the value is larger than <code>POUT</code> . The equation is recomputed without the variable and the process is repeated until no more independent variables can be removed. Then, the independent variable not in the equation with the smallest probability of $F$ is entered if the value is smaller than <code>PIN</code> . All variables in the equation are again examined for removal. This process continues until no variables in the equation can be removed and no variables not in the equation are eligible for entry, or until the maximum number of steps has been reached. <a href="#">For more information, see the topic CRITERIA Subcommand on p. 1711.</a>

The methods that enter or remove the entire variable block in a single step are as follows:

<b>ENTER [varlist]</b>	<i>Forced entry.</i> All variables specified are entered in a single step in order of decreasing tolerance. You can control the order in which variables are entered by specifying the variables on multiple <code>METHOD=ENTER</code> subcommands.
<b>REMOVE varlist</b>	<i>Forced removal.</i> All variables specified are removed in a single step. <code>REMOVE</code> requires a variable list.
<b>TEST (varlist) (varlist)</b>	<i>R<sup>2</sup> change and its significance for sets of independent variables.</i> This method first adds all variables specified on <code>TEST</code> to the current equation. It then removes in turn each subset from the equation and displays requested statistics. Specify test subsets in parentheses. A variable can be used in more than one subset, and each subset can include any number of variables. Variables named on <code>TEST</code> remain in the equation when the method is completed.

### Example

```
REGRESSION VARIABLES=POP15 TO GROWTH, SAVINGS
  /DEPENDENT=SAVINGS
  /METHOD=STEPWISE
  /METHOD=ENTER.
```

- `STEPWISE` applies the stepwise procedure to variables *POP15* to *GROWTH*.
- All variables not in the equation when the `STEPWISE` method is completed will be forced into the equation with `ENTER`.

### Example

```
REGRESSION VARIABLES=(COLLECT)
  /DEPENDENT=SAVINGS
  /METHOD=TEST(MEASURE3 TO MEASURE9) (MEASURE3, INCOME)
  /METHOD=ENTER GROWTH.
```

- The `VARIABLES=(COLLECT)` specification assembles a correlation matrix that includes all variables named on the `DEPENDENT` and `METHOD` subcommands.
- `REGRESSION` first builds the full equation of all the variables named on the first `METHOD` subcommand: *SAVINGS* regressed on *MEASURE3* to *MEASURE9* and *INCOME*. For each set of test variables (*MEASURE3* to *MEASURE9*, and *MEASURE3* and *INCOME*), the  $R^2$  change,  $F$ , probability, sums of squares, and degrees of freedom are displayed.
- *GROWTH* is added to the equation by the second `METHOD` subcommand. Variables *MEASURE3* to *MEASURE9* and *INCOME* are still in the equation when this subcommand is executed.

## STATISTICS Subcommand

`STATISTICS` controls the display of statistics for the equation and for the independent variables.

- If `STATISTICS` is omitted or if it is specified without keywords, `R`, `ANOVA`, `COEFF`, and `OUTS` are displayed (see below).

- If any statistics are specified on `STATISTICS`, only those statistics specifically requested are displayed.
- `STATISTICS` must be specified before `DEPENDENT` and `METHOD` subcommands. The last specified `STATISTICS` affects all equations.

### Global Statistics

<b>DEFAULTS</b>	<i>R, ANOVA, COEFF, and OUTS.</i> These are displayed if <code>STATISTICS</code> is omitted or if it is specified without keywords.
<b>ALL</b>	<i>All statistics except F.</i>

### Equation Statistics

<b>R</b>	<i>Multiple R.</i> <code>R</code> includes $R^2$ , adjusted $R^2$ , and standard error of the estimate displayed in the Model Summary table.
<b>ANOVA</b>	<i>Analysis of variance table.</i> This option includes regression and residual sums of squares, mean square, $F$ , and probability of $F$ displayed in the ANOVA table.
<b>CHANGE</b>	<i>Change in <math>R^2</math>.</i> This option includes the change in $R^2$ between steps, along with the corresponding $F$ and its probability, in the Model Summary table. For each equation, $F$ and its probability are also displayed.
<b>BCOV</b>	<i>Variance-covariance matrix for unstandardized regression coefficients.</i> The statistics are displayed in the Coefficient Correlations table.
<b>XTX</b>	<i>Swept correlation matrix.</i>
<b>COLLIN</b>	<i>Collinearity diagnostics</i> (Belsley et al., 1980). <code>COLLIN</code> includes the variance-inflation factors (VIF) displayed in the Coefficients table, and the eigenvalues of the scaled and uncentered cross-products matrix, condition indexes, and variance-decomposition proportions displayed in the Collinearity Diagnostics table.
<b>SELECTION</b>	<i>Selection statistics.</i> This option includes Akaike information criterion (AIC), Ameniya's prediction criterion (PC), Mallows conditional mean squared error of prediction criterion ( $C_p$ ), and Schwarz Bayesian criterion (SBC) (Judge, Griffiths, Hill, Lutkepohl, and Lee, 1980). The statistics are displayed in the Model Summary table.

### Statistics for the Independent Variables

<b>COEFF</b>	<i>Regression coefficients.</i> This option includes regression coefficients (B), standard errors of the coefficients, standardized regression coefficients (beta), $t$ , and two-tailed probability of $t$ . The statistics are displayed in the Coefficients table.
<b>OUTS</b>	<i>Statistics for variables not yet in the equation that have been named on METHOD subcommands for the equation.</i> <code>OUTS</code> displays the Excluded Variables table showing beta, $t$ , two-tailed probability of $t$ , and minimum tolerance of the variable if it were the only variable entered next.
<b>ZPP</b>	<i>Zero-order, part, and partial correlation.</i> The statistics are displayed in the Coefficients table.
<b>CI(value)</b>	<i>value% confidence interval for the unstandardized regression coefficients.</i> The statistics are displayed in the Coefficients table. If the optional value is unspecified, it defaults to 95. If specified, the value must be greater than 0 and less than or equal to 100.
<b>SES</b>	<i>Approximate standard error of the standardized regression coefficients.</i> (Meyer and Younger, 1976) The statistics are displayed in the Coefficients table.

---

<b>TOL</b>	<i>Tolerance.</i> This option displays tolerance for variables in the equation in the Coefficients table. For variables not yet entered into the equation, TOL displays in the Excluded Variables table the tolerance each variable would have if it were the only variable entered next.
<b>F</b>	<i>F value for B and its probability.</i> This is displayed instead of the <i>t</i> value in the Coefficients or Excluded Variables table.

## **CRITERIA Subcommand**

CRITERIA controls the statistical criteria used to build the regression equations. The way in which these criteria are used depends on the method specified on METHOD. The default criteria are noted in the description of each CRITERIA keyword below.

- The minimum specification is a criterion keyword and its arguments, if any.
- If CRITERIA is omitted or included without specifications, the default criteria are in effect.
- The CRITERIA subcommand must be specified before DEPENDENT and METHOD subcommands. The last specified CRITERIA affects all equations.

## **Tolerance and Minimum Tolerance Tests**

Variables must pass both tolerance and minimum tolerance tests in order to enter and remain in a regression equation. Tolerance is the proportion of the variance of a variable in the equation that is not accounted for by other independent variables in the equation. The minimum tolerance of a variable not in the equation is the smallest tolerance any variable already in the equation would have if the variable being considered were included in the analysis.

If a variable passes the tolerance criteria, it is eligible for inclusion based on the method in effect.

## **Criteria for Variable Selection**

- The ENTER, REMOVE, and TEST methods use only the TOLERANCE criterion.
- BACKWARD removes variables according to the probability of *F*-to-remove (keyword POUT). Specify FOUT to use *F*-to-remove instead.
- FORWARD enters variables according to the probability of *F*-to-enter (keyword PIN). Specify FIN to use *F*-to-enter instead.
- STEPWISE uses both PIN and POUT (or FIN and FOUT) as criteria. If the criterion for entry (PIN or FIN) is less stringent than the criterion for removal (POUT or FOUT), the same variable can cycle in and out until the maximum number of steps is reached. Therefore, if PIN is larger than POUT or FIN is smaller than FOUT, REGRESSION adjusts POUT or FOUT and issues a warning.

- The values for these criteria are specified in parentheses. If a value is not specified, the default values are used.

<b>DEFAULTS</b>	<i>PIN(0.05), POUT(0.10), and TOLERANCE(0.0001).</i> These are the defaults if <b>CRITERIA</b> is omitted. If criteria have been changed, <b>DEFAULTS</b> restores these defaults.
<b>PIN[(value)]</b>	<i>Probability of F-to-enter.</i> The default value is 0.05. Either <b>PIN</b> or <b>FIN</b> can be specified. If more than one is used, the last one specified is in effect.
<b>FIN[(value)]</b>	<i>F-to-enter.</i> The default value is 3.84. Either <b>PIN</b> or <b>FIN</b> can be specified. If more than one is used, the last one specified is in effect.
<b>POUT[(value)]</b>	<i>Probability of F-to-remove.</i> The default value is 0.10. Either <b>POUT</b> or <b>FOUT</b> can be specified. If more than one is used, the last one specified is in effect.
<b>FOUT[(value)]</b>	<i>F-to-remove.</i> The default value is 2.71. Either <b>POUT</b> or <b>FOUT</b> can be specified. If more than one is used, the last one specified is in effect.
<b>TOLERANCE[(value)]</b>	<i>Tolerance.</i> The default value is 0.0001. If the specified tolerance is very low, <b>REGRESSION</b> issues a warning.
<b>MAXSTEPS[(n)]</b>	<i>Maximum number of steps.</i> The value of <b>MAXSTEPS</b> is the sum of the maximum number of steps for each method for the equation. The default values are, for the <b>BACKWARD</b> or <b>FORWARD</b> methods, the number of variables meeting <b>PIN/POUT</b> or <b>FIN/FOUT</b> criteria, and for the <b>STEPWISE</b> method, twice the number of independent variables.

## Confidence Intervals

<b>CIN[(value)]</b>	<i>Reset the value of the percent for confidence intervals.</i> The default is 95%. The specified value sets the percentage interval used in the computation of temporary variable types <b>MCIN</b> and <b>ICIN</b> .
---------------------	--

### Example

```
REGRESSION VARIABLES=POP15 TO GROWTH, SAVINGS
/CRITERIA=PIN(.1) POUT(.15)
/DEPENDENT=SAVINGS
/METHOD=FORWARD.
```

- The **CRITERIA** subcommand relaxes the default criteria for entry and removal for the **FORWARD** method. Note that the specified **PIN** is less than **POUT**.

## ORIGIN and NOORIGIN Subcommands

**ORIGIN** and **NOORIGIN** control whether or not the constant is suppressed. By default, the constant is included in the model (**NOORIGIN**).

- The specification is either the **ORIGIN** or **NOORIGIN** subcommand.
- **ORIGIN** and **NOORIGIN** must be specified before the **DEPENDENT** and **METHOD** subcommands. The last specified remains in effect for all equations.
- **ORIGIN** requests regression through the origin. The constant term is suppressed.
- If you specify **ORIGIN**, statistics requested on the **DESCRIPTIVES** subcommand are computed as if the mean were 0.

- `ORIGIN` and `NOORIGIN` affect the way the correlation matrix is built. If matrix materials are used as input to `REGRESSION`, the keyword that was in effect when the matrix was written should be in effect when that matrix is read.

### Example

```
REGRESSION VAR=(COL)
/ORIGIN
/DEP=HOMICIDE
/METHOD=ENTER POVPC .
```

- The `REGRESSION` command requests an equation that regresses *HOMICIDE* on *POVPCT* and suppresses the constant (`ORIGIN`).

## REGWGT Subcommand

The only specification on `REGWGT` is the name of the variable containing the weights to be used in estimating a weighted least-squares model. With `REGWGT`, the default display is the usual `REGRESSION` display.

- `REGWGT` is a global subcommand.
- If more than one `REGWGT` subcommand is specified on a `REGRESSION` procedure, only the last one is in effect.
- `REGWGT` can be used with `MATRIX OUT` but not with `MATRIX IN`.
- Residuals saved from equations using the `REGWGT` command are not weighted. To obtain weighted residuals, multiply the residuals created with `SAVE` by the square root of the weighting variable in a `COMPUTE` statement.
- `REGWGT` is in effect for all equations and affects the way the correlation matrix is built. Thus, if `REGWGT` is specified on a `REGRESSION` procedure that writes matrix materials to a matrix data file, subsequent `REGRESSION` procedures using that file will be automatically weighted.

### Example

```
REGRESSION VARIABLES=GRADE GPA STARTLEV TREATMNT
/DEPENDENT=GRADE
/METHOD=ENTER
/SAVE PRED(P) .
COMPUTE WEIGHT=1/(P*(1-P)) .
REGRESSION VAR=GRADE GPA STARTLEV TREATMNT
/REGWGT=WEIGHT
/DEP=GRADE
/METHOD=ENTER .
```

- `VARIABLES` builds a correlation matrix that includes *GRADE*, *GPA*, *STARTLEV*, and *TREATMNT*.
- `DEPENDENT` identifies *GRADE* as the dependent variable.
- `METHOD` regresses *GRADE* on *GPA*, *STARTLEV*, and *TREATMNT*.
- `SAVE` saves the predicted values from the regression equation as variable *P* in the active dataset. For more information, see the topic [SAVE Subcommand on p. 1722](#).
- `COMPUTE` creates the variable *WEIGHT* as a transformation of *P*.

- The second REGRESSION procedure performs a weighted regression analysis on the same set of variables using *WEIGHT* as the weighting variable.

### Example

```
REGRESSION VAR=GRADE GPA STARTLEV TREATMNT
  /REGWGT=WEIGHT
  /DEP=GRADE
  /METHOD=ENTER
  /SAVE RESID(RGRADE) .
COMPUTE WRGRADE=RGRADE * SQRT(WEIGHT) .
```

- This example illustrates the use of COMPUTE with SAVE to weight residuals.
- REGRESSION performs a weighted regression analysis of *GRADE* on *GPA*, *STARTLEV*, and *TREATMNT*, using *WEIGHT* as the weighting variable.
- SAVE saves the residuals as *RGRADE*. These residuals are not weighted.
- COMPUTE creates variable *WRGRADE*, which contains the weighted residuals.

## DESCRIPTIVES Subcommand

DESCRIPTIVES requests the display of correlations and descriptive statistics. By default, descriptive statistics are not displayed.

- The minimum specification is simply the subcommand keyword DESCRIPTIVES, which obtains MEAN, STDDEV, and CORR.
- If DESCRIPTIVES is specified with keywords, only those statistics specifically requested are displayed.
- Descriptive statistics are displayed only once for all variables named or implied on VARIABLES.
- Descriptive statistics are based on all valid cases for each variable if PAIRWISE or MEANSUBSTITUTION has been specified on MISSING. Otherwise, only cases with valid values for all variables named or implied on the VARIABLES subcommand are included in the calculation of descriptive statistics.
- If regression through the origin has been requested (subcommand ORIGIN), statistics are computed as if the mean were 0.

<b>NONE</b>	<i>No descriptive statistics. This is the default if the subcommand is omitted.</i>
<b>DEFAULTS</b>	<i>MEAN, STDDEV, and CORR. This is the same as specifying DESCRIPTIVES without specifications.</i>
<b>MEAN</b>	<i>Display variable means in the Descriptive Statistics table.</i>
<b>STDDEV</b>	<i>Display variable standard deviations in the Descriptive Statistics table.</i>
<b>VARIANCE</b>	<i>Display variable variances in the Descriptive Statistics table.</i>
<b>CORR</b>	<i>Display Pearson correlation coefficients in the Correlations table.</i>
<b>SIG</b>	<i>Display one-tailed probabilities of the correlation coefficients in the Correlations table.</i>
<b>BADCORR</b>	<i>Display the correlation coefficients only if some coefficients cannot be computed.</i>
<b>COV</b>	<i>Display covariance in the Correlations table.</i>



<b>XPROD</b>	<i>Display sum of squares and cross-product deviations from the mean in the Correlations table.</i>
<b>N</b>	<i>Display numbers of cases used to compute correlation coefficients in the Correlations table.</i>
<b>ALL</b>	<i>All descriptive statistics.</i>

**Example**

```
REGRESSION DESCRIPTIVES=DEFAULTS SIG COV
/VARIABLES=AGE, FEMALE, YRS_JOB, STARTPAY, SALARY
/DEPENDENT=SALARY
/METHOD=ENTER STARTPAY
/METHOD=ENTER YRS_JOB.
```

- The variable means, standard deviations, and number of cases are displayed in the Descriptive Statistics table and the correlation coefficients, one-tailed probabilities of the correlation coefficients, and covariance are displayed in the Correlations table.
- Statistics are displayed for all variables named on `VARIABLES`, even though only variables `SALARY`, `STARTPAY`, and `YRS_JOB` are used to build the equations.
- `STARTPAY` is entered into the equation by the first `METHOD` subcommand. `YRS_JOB` is entered by the second `METHOD` subcommand.

**SELECT Subcommand**

By default, all cases in the active dataset are considered for inclusion on `REGRESSION`. Use `SELECT` to include a subset of cases in the correlation matrix and resulting regression statistics.

- The required specification on `SELECT` is a logical expression.
- The syntax for the `SELECT` subcommand is as follows:

```
/SELECT=varname relation value
```

- The variable named on `SELECT` should not be specified on the `VARIABLES` subcommand.
- The relation can be `EQ`, `NE`, `LT`, `LE`, `GT`, or `GE`.
- Only cases for which the logical expression on `SELECT` is true are included in the calculation of the correlation matrix and regression statistics.
- All other cases, including those with missing values for the variable named on `SELECT`, are not included in the computations.
- If `SELECT` is specified, residuals and predicted values are calculated and reported separately for both selected and unselected cases by default. [For more information, see the topic RESIDUALS Subcommand on p. 1718.](#)
- Cases deleted from the active dataset with `SELECT IF`, a temporary `SELECT IF`, or `SAMPLE` are not passed to `REGRESSION` and are not included among either the selected or unselected cases.

**Example**

```
REGRESSION SELECT SEX EQ 'M'
/VARIABLES=AGE, STARTPAY, YRS_JOB, SALARY
```

```

/DEPENDENT=SALARY
/METHOD=STEP
/RESIDUALS=NORMPROB.

```

- Only cases with the value M for *SEX* are included in the correlation matrix calculated by REGRESSION.
- Separate normal P\_P plots are displayed for cases with *SEX* equal to M and for other cases. For more information, see the topic RESIDUALS Subcommand on p. 1718.

## **MATRIX Subcommand**

MATRIX reads and writes matrix data files. It can read matrix data files or datasets written by previous REGRESSION procedures or data files or datasets written by other procedures such as CORRELATIONS. The matrix materials REGRESSION writes also include the mean, standard deviation, and number of cases used to compute each coefficient. This information immediately precedes the correlation matrix in the matrix file.

- Either IN or OUT and a matrix file or previously declared dataset name in parentheses are required on MATRIX.
- When used, MATRIX must be the first subcommand specified in a REGRESSION procedure.
- ORIGIN and NOORIGIN affect the way the correlation matrix is built. If matrix materials are used as input to REGRESSION, the keyword that was in effect when the matrix was written should be in effect when that matrix is read.

**OUT ('savfile' | 'dataset')** *Write a matrix data file or dataset.* Specify either a filename, a previously declared dataset name, or an asterisk, enclosed in parentheses. Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. If you specify an asterisk (\*), the matrix data file replaces the active dataset. If you specify an asterisk or a dataset name, the file is not stored on disk unless you use SAVE or XSAVE.

**IN ('savfile' | 'dataset')** *Read a matrix data file or dataset.* Specify either a filename, dataset name created during the current session, or an asterisk enclosed in parentheses. An asterisk reads the matrix data from the active dataset. Filenames should be enclosed in quotes and are read from the working directory unless a path is included as part of the file specification.

## **Format of the Matrix Data File**

- The file has two special variables created by the program: *ROWTYPE\_* and *VARNAME\_*.
- *ROWTYPE\_* is a short string variable with values MEAN, STDDEV, N, and CORR (for Pearson correlation coefficient).
- *VARNAME\_* is a short string variable whose values are the names of the variables used to form the correlation matrix. When *ROWTYPE\_* is CORR, *VARNAME\_* gives the variable associated with that row of the correlation matrix.
- The remaining variables in the file are the variables used to form the correlation matrix.
- To suppress the constant term when ORIGIN is used in the analysis, value OCORR (rather than value CORR) is written to the matrix system file. OCORR indicates that the regression passes through the origin.

## Split Files

- When split-file processing is in effect, the first variables in the matrix data file are the split variables, followed by *ROWTYPE\_*, the independent variable, *VARNAME\_*, and the dependent variables.
- A full set of matrix materials is written for each subgroup defined by the split variable(s).
- A split variable cannot have the same variable name as any other variable written to the matrix data file.
- If a split file is in effect when a matrix is written, the same split file must be in effect when that matrix is read.

## Missing Values

- With *PAIRWISE* treatment of missing values, the matrix of N's used to compute each coefficient is included with the matrix materials.
- With *LISTWISE* treatment (the default) or *MEANSUBSTITUTION*, a single N used to calculate all coefficients is included.

## Example

```
REGRESSION MATRIX IN('/data/pay_data.sav') OUT(*)
/VARIABLES=AGE, STARTPAY, YRS_JOB, SALARY
/DEPENDENT=SALARY
/METHOD=STEP.
```

- *MATRIX IN* reads the matrix data file *pay\_data.sav*.
- A stepwise regression analysis of *SALARY* is performed using *AGE*, *STARTPAY*, and *YRS\_JOB*.
- *MATRIX OUT* replaces the active dataset with the matrix data file that was previously stored in the *pay\_data.sav* file.

## MISSING Subcommand

*MISSING* controls the treatment of cases with missing values. By default, a case that has a user-missing or system-missing value for any variable named or implied on *VARIABLES* is omitted from the computation of the correlation matrix on which all analyses are based.

- The minimum specification is a keyword specifying a missing-value treatment.

### LISTWISE

*Delete cases with missing values listwise.* Only cases with valid values for all variables named on the current *VARIABLES* subcommand are used. If *INCLUDE* is also specified, only cases with system-missing values are deleted listwise. *LISTWISE* is the default if the *MISSING* subcommand is omitted.

### PAIRWISE

*Delete cases with missing values pairwise.* Each correlation coefficient is computed using cases with complete data for the pair of variables correlated. If *INCLUDE* is also specified, only cases with system-missing values are deleted pairwise.

- MEANSUBSTITUTION** *Replace missing values with the variable mean.* All cases are included and the substitutions are treated as valid observations. If `INCLUDE` is also specified, user-missing values are treated as valid and are included in the computation of the means.
- INCLUDE** *Includes cases with user-missing values.* All user-missing values are treated as valid values. This keyword can be specified along with the methods `LISTWISE`, `PAIRWISE`, or `MEANSUBSTITUTION`.

**Example**

```
REGRESSION VARIABLES=POP15, POP75, INCOME, GROWTH, SAVINGS
  /DEPENDENT=SAVINGS
  /METHOD=STEP
  /MISSING=MEANSUBSTITUTION.
```

- System-missing and user-missing values are replaced with the means of the variables when the correlation matrix is calculated.

**RESIDUALS Subcommand**

`RESIDUALS` controls the display and labeling of summary information on outliers as well as the display of the Durbin-Watson statistic and histograms and normal probability plots for the temporary variables.

- If `RESIDUALS` is specified without keywords, it displays a histogram of residuals, a normal probability plot of residuals, the values of `$CASENUM` and `ZRESID` for the 10 cases with the largest absolute value of `ZRESID`, and the Durbin-Watson test statistic. The histogram and the normal plot are standardized.
- If any keywords are specified on `RESIDUALS`, only the requested information and plots are displayed.

- DEFAULTS** *DURBIN, NORMPROB(ZRESID), HISTOGRAM(ZRESID), OUTLIERS(ZRESID).* These are the defaults if `RESIDUALS` is used without specifications.
- HISTOGRAM(tempvars)** *Histogram of the temporary variable or variables.* The default is `ZRESID`. You can request histograms for `PRED`, `RESID`, `ZPRED`, `DRESID`, `ADJPRED`, `SRESID`, `SDRESID`, `SEMPRED`, `MAHAL`, `COOK`, and `LEVER`. The specification of any other temporary variable will result in an error.
- NORMPROB(tempvars)** *Normal probability (P-P) plot.* The default is `ZRESID`. The other temporary variables for which normal probability plots are available are `PRED`, `RESID`, `ZPRED`, `DRESID`, `SRESID`, and `SDRESID`. The specification of any other temporary variable will result in an error. Normal probability plots are always displayed in standardized form; therefore, when `PRED`, `RESID`, or `DRESID` is requested, the standardized equivalent `ZPRED`, `ZRESID` or `SDRESID` is displayed.
- OUTLIERS(tempvars)** *The 10 cases with the largest absolute values of the specified temporary variables.* The default is `ZRESID`. The output includes the values of `$CASENUM` and of the temporary variables for the 10 cases. The other temporary variables available for `OUTLIERS` are `RESID`, `SRESID`, `SDRESID`, `DRESID`, `MAHAL`, and `COOK`. The specification of any temporary variable other than these will result in an error.
- DURBIN** *Display Durbin-Watson test statistic in the Model Summary table.*

<b>ID(varname)</b>	<i>ID variable providing case labels for use with point selection mode in the Chart Editor. Applicable to scatterplots produced by SCATTERPLOT, PARTIALPLOT, and RESIDUALS. Any variable in the active dataset can be named.</i>
<b>SEPARATE</b>	<i>Separate reporting of residuals statistics and plots for selected and unselected cases. This is the default.</i>
<b>POOLED</b>	<i>Pooled plots and statistics using all cases in the working file when the SELECT subcommand is in effect. This is an alternative to SEPARATE.</i>

**Example**

```
/RESID=DEFAULT ID(SVAR)
```

- **DEFAULT** produces the default residuals statistics: Durbin-Watson statistic, a normal probability plot and histogram of *ZRESID*, and an outlier listing for *ZRESID*.
- Descriptive statistics for *ZRESID*, *RESID*, *PRED*, and *ZPRED* are automatically displayed.
- *SVAR* is specified as the case identifier on the outlier output.

**CASEWISE Subcommand**

**CASEWISE** requests a Casewise Diagnostics table of residuals. You can specify a temporary residual variable for casewise listing (via the **PLOT** keyword). You can also specify variables to be listed in the table for each case.

- If **CASEWISE** is used without any additional specifications, it displays a Casewise Diagnostics table of *ZRESID* for cases whose absolute value of *ZRESID* is at least 3. By default, the values of the case sequence number, *DEPENDENT*, *PRED*, and *RESID* are listed for each case.
- Defaults remain in effect unless specifically altered.

<b>DEFAULTS</b>	<i>OUTLIERS(3), PLOT(ZRESID), DEPENDENT, PRED, and RESID.</i> These are the defaults if the subcommand is used without specifications.
<b>OUTLIERS(value)</b>	<i>List only cases for which the absolute standardized value of the listed variable is at least as large as the specified value. The default value is 3. Keyword OUTLIERS is ignored if keyword ALL is also present.</i>
<b>ALL</b>	<i>Include all cases in the Casewise Diagnostic table. ALL is the alternative to keyword OUTLIERS.</i>
<b>PLOT(tempvar)</b>	<i>List the values of the temporary variable in the Casewise Diagnostics table. The default temporary variable is ZRESID. Other variables that can be listed are RESID, DRESID, SRESID, and SDRESID. The specification of any temporary variable other than these will result in an error. When requested, RESID is standardized and DRESID is Studentized in the output.</i>
<b>tempvars</b>	<i>Display the values of these variables next to the casewise list entry for each case. The default variables are DEPENDENT (the dependent variable), PRED, and RESID. Any of the other temporary variables can be specified. If an ID variable is specified on RESIDUALS, the ID variable is also listed.</i>

**Example**

```
/CASEWISE=DEFAULT ALL SRE MAH COOK SDR
```

- This example requests a Casewise Diagnostics table of the standardized residuals for all cases.
- *ZRESID*, the dependent variable, and the temporary variables *PRED*, *RESID*, *SRESID*, *MAHAL*, *COOK*, and *SDRESID* are for all cases.

## **SCATTERPLOT Subcommand**

SCATTERPLOT names pairs of variables for scatterplots.

- The minimum specification for SCATTERPLOT is a pair of variables in parentheses. There are no default specifications.
- You can specify as many pairs of variables in parentheses as you want.
- The first variable named in each set of parentheses is plotted along the vertical axis, and the second variable is plotted along the horizontal axis.
- Plotting symbols are used to represent multiple points occurring at the same position.
- You can specify any variable named on the VARIABLES subcommand.
- You can specify *PRED*, *RESID*, *ZPRED*, *ZRESID*, *DRESID*, *ADJPRED*, *SRESID*, *SDRESID*, *SEPPRED*, *MAHAL*, *COOK*, and *LEVER*. The specification of any other temporary variables will result in an error.
- Specify an asterisk before temporary variable names to distinguish them from user-defined variables. For example, use *\*PRED* to specify *PRED*.

### **Example**

```
/SCATTERPLOT (*RES, *PRE) (*RES, SAVINGS)
```

- This example specifies two scatterplots: residuals against predicted values and residuals against the values of the variable *SAVINGS*.

## **PARTIALPLOT Subcommand**

PARTIALPLOT requests partial regression plots. Partial regression plots are scatterplots of the residuals of the dependent variable and an independent variable when both of these variables are regressed on the rest of the independent variables.

- If PARTIALPLOT is included without any additional specifications, it produces a partial regression plot for every independent variable in the equation. The plots appear in the order the variables are specified or implied on the VARIABLES subcommand.
- If variables are specified on PARTIALPLOT, only the requested plots are displayed. The plots appear in the order the variables are listed on the PARTIALPLOT subcommand.
- At least two independent variables must be in the equation for partial regression plots to be produced.

**ALL** *Plot all independent variables in the equation.* This is the default.  
**varlist** *Plot the specified variables.* Any variable entered into the equation can be specified.

**Example**

```
REGRESSION VARS=PLOT15 TO SAVINGS
  /DEP=SAVINGS
  /METH=ENTER
  /RESID=DEFAULTS
  /PARTIAL.
```

- A partial regression plot is produced for every independent variable in the equation.

**OUTFILE Subcommand**

OUTFILE saves the parameter covariance or correlation matrix with parameter estimates, standard errors, significance values, and residual degrees of freedom for each term in the final equation in IBM® SPSS® Statistics format data files. It also saves model information in XML format.

- The OUTFILE subcommand must follow the last METHOD subcommand.
- Only one OUTFILE subcommand is allowed. If you specify more than one, only the last one is executed.
- You must specify at least one keyword and a quoted file specification, enclosed in parentheses. For COVB and CORB, you can specify a previously declared dataset (DATASET DECLARE command) instead of a file.
- You cannot save the parameter statistics as the active dataset.
- COVB and CORB are mutually exclusive.
- MODEL cannot be used if split file processing is on (SPLIT FILE command) or if more than one dependent (DEPENDENT subcommand) variable is specified.
- If you specify an external file name, you should include the .sav extension in the specification. There is no default extension.

**COVB ('savfile'|'dataset')**

*Write the parameter covariance matrix with other statistics.*

**CORB ('savfile'|'dataset')**

*Write the parameter correlation matrix with other statistics.*

**MODEL ('file')**

*Write model information to an XML file. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#)*

**PARAMETER('file')**

*Write parameter estimates only to an XML file. You can use this model file to apply the model information to other data files for scoring purposes. [For more information, see the topic Scoring expressions on p. 116.](#)*

**Example**

```
REGRESSION DEPENDENT=Y
  /METHOD=ENTER X1 X2
  /OUTFILE CORB ('/data/covx1x2y.sav').
```

- The OUTFILE subcommand saves the parameter correlation matrix, and the parameter estimates, standard errors, significance values, and residual degrees of freedom for the constant term, *X1* and *X2*.

## SAVE Subcommand

Use SAVE to add one or more residual or fit variables to the active dataset.

- The specification on SAVE is one or more of the temporary variable types, each followed by an optional name in parentheses for the new variable.
- New variable names must be unique.
- If new names are not specified, REGRESSION generates a rootname using a shortened form of the temporary variable name with a suffix to identify its creation sequence.
- If you specify DFBETA or SDBETA on the SAVE subcommand, the number of new variables saved is the total number of variables in the equation.

**FITS**            *Save all influence statistics.* FITS saves *DFFIT*, *SDFIT*, *DFBETA*, *SDBETA*, and *COVRATIO*. You cannot specify new variable names when using this keyword. Default names are generated.

### Example

```
/SAVE=PRED(PREDVAL) RESID(RESIDUAL) COOK(CDISTANC)
```

- This subcommand adds three variables to the end of the active dataset: *PREDVAL*, containing the unstandardized predicted value for each case; *RESIDUAL*, containing the unstandardized residual; and *CDISTANC*, containing Cook's distance.

### Example

```
/SAVE=PRED RESID
```

- This subcommand adds two variables named *PRE\_1* and *RES\_1* to the end of the active dataset.

### Example

```
REGRESSION DEPENDENT=Y
/METHOD=ENTER X1 X2
/SAVE DFBETA(DFBVAR) .
```

- The SAVE subcommand creates and saves three new variables with the names *DFBVAR0*, *DFBVAR1*, and *DFBVAR2*.

### Example

```
REGRESSION VARIABLES=SAVINGS INCOME POP15 POP75 GROWTH
/DEPENDENT=SAVINGS
/METHOD=ENTER INCOME POP15 POP75
/SAVE=PRED(PREDV) SDBETA(BETA) ICIN.
```

- The SAVE subcommand adds seven variables to the end of the file: *PREDV*, containing the unstandardized predicted value for the case; *BETA0*, the standardized *DFBETA* for the intercept; *BETA1*, *BETA2*, and *BETA3*, the standardized *DFBETA*'s for the three independent variables in the model; *LICI\_1*, the lower bound for the prediction interval for an individual case; and *UICI\_1*, the upper bound for the prediction interval for an individual case.



## TEMPLATE Subcommand

TEMPLATE uses an existing file as a template and applies it to any charts generated by the current REGRESSION command. The template overrides the default settings that are used to create any charts. Templates are created in the Chart Editor by saving an existing chart as a template.

### Example

```
REGRESSION VARIABLES=pop15,pop75,income,growth,savings  
/DEPENDENT=savings  
/METHOD=ENTER pop15,pop75,income  
/METHOD=ENTER growth  
/RESIDUALS HISTOGRAM(ZRESID)  
/TEMPLATE=' /templates/mytemplate.sgt'.
```

## References

- Belsley, D. A., E. Kuh, and R. E. Welsch. 1980. *Regression diagnostics: Identifying influential data and sources of collinearity*. New York: John Wiley and Sons.
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- Cook, R. D. 1977. Detection of influential observations in linear regression. *Technometrics*, 19, 15–18.
- Dillon, W. R., and M. Goldstein. 1984. *Multivariate analysis: Methods and applications*. New York: John Wiley and Sons.
- Hoaglin, D. C., and R. E. Welsch. 1978. The hat matrix in regression and ANOVA. *American Statistician*, 32, 17–22.
- Judge, G. G., W. E. Griffiths, R. C. Hill, H. Lutkepohl, and T. C. Lee. 1980. *The theory and practice of econometrics*, 2nd ed. New York: John Wiley and Sons.
- Meyer, L. S., and M. S. Younger. 1976. Estimation of standardized coefficients. *Journal of the American Statistical Association*, 71, 154–157.
- Velleman, P. F., and R. E. Welsch. 1981. Efficient computing of regression diagnostics. *American Statistician*, 35, 234–242.

# RELIABILITY

RELIABILITY is available in the Statistics Base option.

```
RELIABILITY VARIABLES={varlist}
                        {ALL      }

[/SCALE(scalename)=varlist ]

[/MODEL={ALPHA          } ]
        {SPLIT[(n)]     }
        {GUTTMAN        }
        {PARALLEL       }
        {STRICTPARALLEL}

[/STATISTICS=[DESCRIPTIVE] [SCALE]      [{ANOVA          } ] [ALL] ]
             [COVARIANCES] [TUKEY]      {ANOVA FRIEDMAN}
             [CORRELATIONS] [HOTELLING]  {ANOVA COCHRAN }

[/SUMMARY=[MEANS] [VARIANCE] [COV] [CORR] [TOTAL] [ALL] ]

[/ICC=[ {MODEL(ONEWAY)                                } ]
      { [MODEL({MIXED**})] [TYPE({CONSISTENCY**})] ] }
      {RANDOM } {ABSOLUTE } }
      [CIN={95**}] [TESTVAL={0**}] ]
      {n } {p } ]

[/METHOD=COVARIANCE]

[/MISSING={EXCLUDE**}]
          {INCLUDE }

[/MATRIX = [IN({*          })] [OUT({*          })] [NOPRINT]]
           {'savfile'|'dataset'} {'savfile'|'dataset'}
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 17.0

- Increased limits on numbers of variables allowed on the VARIABLES and SCALE lists.

## Example

```
RELIABILITY VARIABLES=SCORE1 TO SCORE10
  /SCALE (OVERALL) = ALL
  /MODEL = ALPHA
  /SUMMARY = MEANS TOTAL.
```

## Overview

RELIABILITY estimates reliability statistics for the components of multiple-item additive scales. It uses any one of five models for reliability analysis and offers a variety of statistical displays. RELIABILITY can also be used to perform a repeated measures analysis of variance, a

two-way factorial analysis of variance with one observation per cell, Tukey's test for additivity, Hotelling's  $T$ -square test for equality of means in repeated measures designs, and Friedman's two-way analysis of variance on ranks. For more complex repeated measures designs, use the GLM procedure (available in the Advanced Models option).

### **Options**

**Model Type.** You can specify any one of five models on the MODEL subcommand.

**Statistical Display.** Statistics available on the STATISTICS subcommand include descriptive statistics, correlation and covariance matrices, a repeated measures analysis of variance table, Hotelling's  $T$ -square, Tukey's test for additivity, Friedman's chi-square for the analysis of ranked data, and Cochran's  $Q$ .

**Computational Method.** You can force RELIABILITY to use the covariance method, even when you are not requesting any output that requires it, by using the METHOD subcommand.

**Matrix Input and Output.** You can read data in the form of correlation matrices and you can write correlation-type matrix materials to a data file using the MATRIX subcommand.

### **Basic Specification**

The basic specification is VARIABLES and a variable list. By default, RELIABILITY displays the number of cases, number of items, and Cronbach's alpha. Whenever possible, it uses an algorithm that does not require the calculation of the covariance matrix.

### **Subcommand Order**

- VARIABLES must be specified first.
- The remaining subcommands can be named in any order.

### **Operations**

- STATISTICS and SUMMARY are cumulative. If you enter them more than once, all requested statistics are produced for each scale.
- If you request output that is not available for your model or for your data, RELIABILITY ignores the request.
- RELIABILITY uses an economical algorithm whenever possible but calculates a covariance matrix when necessary (see [METHOD Subcommand on p. 1729](#)).

### **Limitations**

- Maximum 1 VARIABLES subcommand.
- Maximum 1 SCALE subcommand.
- Maximum 10,000 variables on the VARIABLES subcommand.
- Maximum 10,000 variables on the SCALE subcommand.

## Examples

```
RELIABILITY VARIABLES=SCORE1 TO SCORE10.
```

- This example analyzes a scale (labeled *ALL* in the display output) that includes all 10 items.
- Because there is no `SUMMARY` subcommand, no summary statistics are displayed.

### Reliability Analysis of Survey Items

```
RELIABILITY
/VARIABLES=any bored critics peers writers director cast
/SCALE(ALPHA)=ALL /MODEL=ALPHA
/STATISTICS=DESCRIPTIVE CORR.
```

- The procedure performs a reliability analysis for variables *any*, *bored*, *critics*, *peers*, *writers*, *director*, and *cast*.
- The scale specified on the `SCALE` subcommand is labeled `ALPHA` and includes all variables specified on the `VARIABLES` subcommand.
- The `MODEL` subcommand specifies that the reliability analysis is based on Cronbach's alpha.
- Item means and standard deviations and the inter-item correlation matrix are requested on the `STATISTICS` subcommand.

### Intraclass Correlation Coefficients

```
RELIABILITY
/VARIABLES=judge1 judge2 judge3 judge4 judge5 judge6 judge7
/SCALE(ALPHA)=ALL
/MODEL=ALPHA
/STATISTICS=DESCRIPTIVE
/ICC=MODEL(RANDOM) TYPE(CONSISTENCY) CIN=95 TESTVAL=0 .
```

- The procedure performs a reliability analysis for variables *judge1*, *judge2*, *judge3*, *judge4*, *judge5*, *judge6*, and *judge7*.
- The scale specified on the `SCALE` subcommand is labeled `ALPHA` and includes all variables specified on the `VARIABLES` subcommand.
- The `MODEL` subcommand specifies that the reliability analysis is based on Cronbach's alpha.
- Item means and standard deviations are requested on the `STATISTICS` subcommand.
- The `ICC` subcommand requests a two-way random effect model for the intraclass correlation coefficient, using the consistency definition. 95% confidence intervals will be produced and a 0.05 significance level in testing the ICC against the value 0.

## VARIABLES Subcommand

`VARIABLES` specifies the variables to be used in the analysis. Only numeric variables can be used.

- `VARIABLES` is required and must be specified first.
- You can use keyword `ALL` to refer to all user-defined variables in the active dataset.

## SCALE Subcommand

SCALE defines a scale for analysis, providing a label for the scale and specifying its component variables. If SCALE is omitted, all variables named on VARIABLES are used, and the label for the scale is *ALL*.

- The label is specified in parentheses after SCALE. It can have a maximum of 255 bytes before being truncated.
- RELIABILITY does not add any new variables to the active dataset. The label is used only to identify the output. If the analysis is satisfactory, use COMPUTE to create a new variable containing the sum of the component items.
- Variables named on SCALE must have been named on the VARIABLES subcommand. Use the keyword ALL to refer to all variables named on the VARIABLES subcommand.

### Example

```
RELIABILITY VARIABLES = ITEM1 TO ITEM20
  /SCALE (A) = ITEM1 TO ITEM10.
```

```
RELIABILITY VARIABLES = ITEM1 TO ITEM20
  /SCALE (B) = ITEM1 TO ITEM20.
```

- Analyses for scales A and B both use only cases that have complete data for items 1 through 20.

## MODEL Subcommand

MODEL specifies the type of reliability analysis for the scale named on the SCALE subcommand.

<b>ALPHA</b>	<i>Cronbach's <math>\alpha</math>.</i> Standardized item $\alpha$ is displayed. This is the default.
<b>SPLIT [(n)]</b>	<i>Split-half coefficients.</i> You can specify a number in parentheses to indicate how many items should be in the second half. For example, MODEL SPLIT (6) uses the last six variables for the second half and all others for the first. By default, each half has an equal number of items, with the odd item, if any, going to the first half.
<b>GUTTMAN</b>	<i>Guttman's lower bounds for true reliability.</i>
<b>PARALLEL</b>	<i>Maximum-likelihood reliability estimate under parallel assumptions.</i> This model assumes that items have the same variance but not necessarily the same mean.
<b>STRICTPARALLEL</b>	<i>Maximum-likelihood reliability estimate under strictly parallel assumptions.</i> This model assumes that items have the same means, the same true score variances over a set of objects being measured, and the same error variance over replications.

## STATISTICS Subcommand

STATISTICS displays optional statistics. There are no default statistics.

- STATISTICS is cumulative. If you enter it more than once, all requested statistics are produced for each scale.

<b>DESCRIPTIVES</b>	<i>Item means and standard deviations.</i>
<b>COVARIANCES</b>	<i>Inter-item variance-covariance matrix.</i>
<b>CORRELATIONS</b>	<i>Inter-item correlation matrix.</i>
<b>SCALE</b>	<i>Scale means and scale variances.</i>
<b>TUKEY</b>	<i>Tukey's test for additivity.</i> This helps determine whether a transformation of the items is needed to reduce nonadditivity. The test displays an estimate of the power to which the items should be raised in order to be additive.
<b>HOTELLING</b>	<i>Hotelling's T-square.</i> This is a test for equality of means among the items.
<b>ANOVA</b>	<i>Repeated measures analysis of variance table.</i>
<b>FRIEDMAN</b>	<i>Friedman's chi-square and Kendall's coefficient of concordance.</i> These apply to ranked data. You must request ANOVA in addition to FRIEDMAN; Friedman's chi-square appears in place of the usual $F$ test. If the ANOVA keyword is not specified, the FRIEDMAN keyword is silently ignored.
<b>COCHRAN</b>	<i>Cochran's Q.</i> This applies when all items are dichotomies. You must request ANOVA in addition to COCHRAN; the $Q$ statistic appears in place of the usual $F$ test. If the ANOVA keyword is not specified, the COCHRAN keyword is silently ignored.
<b>ALL</b>	<i>All applicable statistics.</i>

## ICC Subcommand

ICC displays intraclass correlation coefficients for single measure and average measure. Single measure applies to single measurements—for example, the rating of judges, individual item scores, or the body weights of individuals. Average measure, however, applies to average measurements, for example, the average rating of  $k$  judges, or the average score for a  $k$ -item test.

<b>MODEL</b>	<i>Model.</i> You can specify the model for the computation of ICC. There are three keywords for this option. ONEWAY is the one-way random effects model (people effects are random). RANDOM is the two-way random effect model (people effects and the item effects are random). MIXED is the two-way mixed (people effects are random and the item effects are fixed). MIXED is the default. Only one model can be specified.
<b>TYPE</b>	<i>Type of definition.</i> When the model is RANDOM or MIXED, one of the two TYPE keywords may be given. CONSISTENCY is the consistency definition and ABSOLUTE is the absolute agreement definition. For the consistency coefficient, the between measures variance is excluded from the denominator variance, and for absolute agreement, it is not.
<b>CIN</b>	<i>The value of the percent for confidence interval and significance level of the hypothesis testing.</i>
<b>TESTVAL</b>	<i>The value with which an estimate of ICC is compared.</i> The value should be between 0 and 1.

## SUMMARY Subcommand

SUMMARY displays summary statistics for each individual item in the scale.

- **SUMMARY** is cumulative. If you enter it more than once, all requested statistics are produced for the scale.
- You can specify one or more of the following:

<b>MEANS</b>	<i>Statistics on item means.</i> The average, minimum, maximum, range, ratio of maximum to minimum, and variance of the item means.
<b>VARIANCE</b>	<i>Statistics on item variances.</i> This displays the same statistics as for <b>MEANS</b> .
<b>COVARIANCES</b>	<i>Statistics on item covariances.</i> This displays the same statistics as for <b>MEANS</b> .
<b>CORRELATIONS</b>	<i>Statistics on item correlations.</i> This displays the same statistics as for <b>MEANS</b> .
<b>TOTAL</b>	<i>Statistics comparing each individual item to the scale composed of the other items.</i> The output includes the scale mean, variance, and Cronbach's $\alpha$ without the item, and the correlation between the item and the scale without it.
<b>ALL</b>	<i>All applicable summary statistics.</i>

## **METHOD Subcommand**

By default, **RELIABILITY** uses a computational method that does not require the calculation of a covariance matrix wherever possible. **METHOD** forces **RELIABILITY** to calculate the covariance matrix. Only a single specification applies to **METHOD**:

**COVARIANCE**      *Calculate and use the covariance matrix, even if it is not needed.*

If **METHOD** is not specified, **RELIABILITY** computes the covariance matrix for all variables on each **VARIABLES** subcommand only if any of the following is true:

- You specify a model other than **ALPHA** or **SPLIT**.
- You request **COV**, **CORR**, **FRIEDMAN**, or **HOTELLING** on the **STATISTICS** subcommand.
- You request anything other than **TOTAL** on the **SUMMARY** subcommand.
- You write the matrix to a matrix data file, using the **MATRIX** subcommand.

## **MISSING Subcommand**

**MISSING** controls the deletion of cases with user-missing data.

- **RELIABILITY** deletes cases from analysis if they have a missing value for any variable named on the **VARIABLES** subcommand. By default, both system-missing and user-missing values are excluded.

**EXCLUDE**      *Exclude user-missing and system-missing values.* This is the default.

**INCLUDE**      *Treat user-missing values as valid.* Only system-missing values are excluded.

## **MATRIX Subcommand**

**MATRIX** reads and writes IBM® SPSS® Statistics matrix data files.

- Either **IN** or **OUT** and the matrix file in parentheses are required. When both **IN** and **OUT** are used on the same **RELIABILITY** procedure, they can be specified on separate **MATRIX** subcommands or on the same subcommand.

- If both `IN` and `OUT` are used on the same `RELIABILITY` command and there are grouping variables in the matrix input file, these variables are treated as if they were split variables. Values of the grouping variables in the input matrix are passed on to the output matrix (see [Split Files on p. 1731](#)).
- OUT ('savfile']'dataset')**      *Write a matrix data file or dataset.* Specify either a filename, a previously declared dataset name, or an asterisk, enclosed in parentheses. Filenames should be enclosed in quotes and are stored in the working directory unless a path is included as part of the file specification. If you specify an asterisk (\*), the matrix data file replaces the active dataset. If you specify an asterisk or a dataset name, the file is not stored on disk unless you use `SAVE` or `XSAVE`.
- IN ('savfile']'dataset')**      *Read a matrix data file or dataset.* Specify either a filename, dataset name created during the current session, or an asterisk enclosed in parentheses. An asterisk reads the matrix data from the active dataset. Filenames should be enclosed in quotes and are read from the working directory unless a path is included as part of the file specification.

### **Matrix Output**

- `RELIABILITY` writes correlation-type matrices that include the number of cases, means, and standard deviations with the matrix materials (see [Format of the Matrix Data File](#) below for a description of the file). These matrix materials can be used as input to `RELIABILITY` or other procedures.
- Any documents contained in the active dataset are not transferred to the matrix file.
- `RELIABILITY` displays the scale analysis when it writes matrix materials. To suppress the display of scale analysis, specify keyword `NOPRINT` on `MATRIX`.

### **Matrix Input**

- `RELIABILITY` can read a matrix data file created by a previous `RELIABILITY` command or by another procedure. The matrix input file must have records of type `N`, `MEAN`, `STDDEV`, and `CORR` for each split-file group. For more information, see the `Universals` section.
- Variable names, variable and value labels, and print and write formats are read from the dictionary of the matrix data file.
- `MATRIX=IN` cannot be used unless an active dataset has already been defined. To read an existing matrix data file at the beginning of a session, use `GET` to retrieve the matrix file and then specify `IN(*)` on `MATRIX`.

### **Format of the Matrix Data File**

- The matrix data file includes two special variables: `ROWTYPE_` and `VARNAME_`. Variable `ROWTYPE_` is a short string variable having values `N`, `MEAN`, `STDDEV`, and `CORR`. Variable `VARNAME_` is a short string variable whose values are the names of the variables used to form the correlation matrix.
- When `ROWTYPE_` is `CORR`, `VARNAME_` gives the variable associated with that row of the correlation matrix.
- The remaining variables in the matrix file are the variables used to form the correlation matrix.



## Split Files

- When split-file processing is in effect, the first variables in the matrix data file will be the split variables, followed by *ROWTYPE\_*, *VARNAME\_*, and the dependent variable(s).
- If grouping variables are in the matrix input file, their values are between *ROWTYPE\_* and *VARNAME\_*. The grouping variables are treated like split-file variables.
- A full set of matrix materials is written for each split-file group defined by the split variables.
- A split variable cannot have the same variable name as any other variable written to the matrix data file.
- If split-file processing is in effect when a matrix is written, the same split file must be in effect when that matrix is read by any procedure.

## Missing Values

Missing-value treatment affects the values written to a matrix data file. When reading a matrix data file, be sure to specify a missing-value treatment on *RELIABILITY* that is compatible with the treatment that was in effect when the matrix materials were generated.

### Example: Matrix Output to External File

```
DATA LIST / TIME1 TO TIME5 1-10.
BEGIN DATA
  0 0 0 0 0
  0 0 1 1 0
  0 0 1 1 1
  0 1 1 1 1
  0 0 0 0 1
  0 1 0 1 1
  0 0 1 1 1
  1 0 0 1 1
  1 1 1 1 1
  1 1 1 1 1
END DATA.
RELIABILITY VARIABLES=TIME1 TO TIME5
/MATRIX=OUT('/data/relmtx.sav').
LIST.
```

- *RELIABILITY* reads data from the active dataset and writes one set of matrix materials to file *relmtx.sav*.
- The active dataset is still the file defined by *DATA LIST*. Subsequent commands are executed in this file.

### Example: Matrix Output to Active Dataset

```
DATA LIST / TIME1 TO TIME5 1-10.
BEGIN DATA
  0 0 0 0 0
  0 0 1 1 0
  0 0 1 1 1
  0 1 1 1 1
  0 0 0 0 1
  0 1 0 1 1
  0 0 1 1 1
```

## RELIABILITY

```

1 0 0 1 1
1 1 1 1 1
1 1 1 1 1
END DATA.
RELIABILITY VARIABLES=TIME1 TO TIME5
/MATRIX=OUT(*) NOPRINT.
LIST.

```

- RELIABILITY writes the same matrix as in the previous example. However, the matrix data file replaces the active dataset. The LIST command is executed in the matrix file, not in the file defined by DATA LIST.
- Because NOPRINT is specified on MATRIX, scale analyses are not displayed.

**Example: Matrix Output to Active Dataset**

```

GET FILE='/data/relmtx.sav'.
RELIABILITY VARIABLES=ALL
/MATRIX=IN(*).

```

- This example assumes that you are starting a new session and want to read an existing matrix data file. GET retrieves the matrix data file *relmtx.sav*.
- MATRIX=IN specifies an asterisk because the matrix data file is the active dataset. If MATRIX=IN('/data/relmtx.sav') is specified, an error message is issued.
- If the GET command is omitted, an error message is issued.

**Example: Matrix Input from External File**

```

GET FILE='/data/personnel.sav'.
FREQUENCIES VARIABLE=AGE.

RELIABILITY VARIABLES=ALL
/MATRIX=IN('/data/relmtx.sav').

```

- This example performs a frequencies analysis on file *personnel.sav* and then uses a different file containing matrix data for RELIABILITY. The file is an existing matrix data file. In order for this to work, the analysis variables named in *relmtx.sav* must also exist in *personnel.sav*.
- *relmtx.sav* must have records of type N, MEAN, STDDEV, and CORR for each split-file group.
- *relmtx.sav* does not replace *personnel.sav* as the active dataset.

**Example: Matrix Input from Working File**

```

GET FILE='/data/personnel.sav'.
CORRELATIONS VARIABLES=V1 TO V5
/MATRIX=OUT(*).
RELIABILITY VARIABLES=V1 TO V5
/MATRIX=IN(*).

```

- RELIABILITY uses matrix input from procedure CORRELATIONS. An asterisk is used to specify the active dataset for both the matrix output from CORRELATIONS and the matrix input for RELIABILITY.

# RENAME VARIABLES

```
RENAME VARIABLES {(varname=varname) [(varname ...)]}  
                 {(varnames=varnames) }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
RENAME VARIABLES (JOB CAT=TITLE) .
```

## Overview

RENAME VARIABLES changes the names of variables in the active dataset while preserving their original order, values, variable labels, value labels, missing values, and print and write formats.

### Basic Specification

- The basic specification is an old variable name, an equals sign, and the new variable name. The equals sign is required.

### Syntax Rules

- Multiple sets of variable specifications are allowed. Each set can be enclosed in parentheses.
- You can specify a list of old variable names followed by an equals sign and a list of new variable names. The same number of variables must be specified on both lists. A single set of parentheses enclosing the entire specification is required for this method.
- Keyword `TO` can be used on the left side of the equals sign to refer to variables in the active dataset, and on the right side of the equals sign to generate new variable names.
- Old variable names do not need to be specified according to their order in the active dataset.
- Name changes take place in one operation. Therefore, variable names can be exchanged between two variables (see the Examples on p. 1733).
- Multiple `RENAME VARIABLES` commands are allowed.
- `RENAME VARIABLES` cannot follow either a `TEMPORARY` or a `MODEL PROGRAM` command.

## Examples

### Renaming Multiple Variables

```
RENAME VARIABLES (MOHIRED=MOSTART) (YRHIRED=YRSTART) .
```

- `MOHIRED` is renamed to `MOSTART` and `YRHIRED` to `YRSTART`. The parentheses are optional.

```
RENAME VARIABLES (MOHIRED YRHIRED=MOSTART YRSTART) .
```

- The same name changes are specified as in the previous example. The parentheses are required, since variable lists are used.

### ***Exchanging Variable Names***

```
RENAME VARIABLES (A=B) (B=A) .
```

- Variable names are exchanged between two variables: A is renamed to B, and B is renamed to A.

## ***Mixed Case Variable Names***

You can use the `RENAME VARIABLES` command to change the case of any characters in a variable name.

### ***Example***

```
RENAME VARIABLES (newvariable = NewVariable) .
```

- For the existing variable name specification, case is ignored. Any combination of upper and lower case will work.
- For the new variable name, case will be preserved as entered for display purposes.

# REPEATING DATA

```
REPEATING DATA [FILE=file]

[/ENCODING='encoding specification']

/STARTS=beg col[-end col] /OCCURS={value }
                        {varname}

[/LENGTH={value } ] [/CONTINUED[=beg col[-end col]]]
                        {varname}

[/ID={col loc}=varname] [/{TABLE } ]
                        {format }      {NOTABLE}

/DATA=variable specifications
```

## Release History

### Release 16.0

- ENCODING subcommand added for Unicode support.

## Example

```
INPUT PROGRAM.
DATA LIST / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
REPEATING DATA STARTS=12 /OCCURS=NUMVEH
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.
```

## Overview

REPEATING DATA reads input cases whose records contain repeating groups of data. For each repeating group, REPEATING DATA builds one output case in the active dataset. All of the repeating groups in the data must contain the same type of information, although the number of groups for each input case may vary. Information common to the repeating groups for each input case can be recorded once for that case and then spread to each resulting output case. In this respect, a file with a repeating data structure is like a hierarchical file with both levels of information recorded on a single record rather than on separate record types. For information on reading hierarchical files, see FILE TYPE-END FILE TYPE.

REPEATING DATA must be used within an INPUT PROGRAM structure or within a FILE TYPE structure with mixed or nested data. In an INPUT PROGRAM structure, REPEATING DATA must be preceded by a DATA LIST command. In a FILE TYPE structure, DATA LIST is needed only if there are variables to be spread to each resulting output case.

## Options

**Length of Repeating Groups.** If the length of the repeating groups varies across input cases, you can specify a variable that indicates the length on the LENGTH subcommand. You can also use LENGTH if you do not want to read all the data in each repeating group.

**Continuation Records.** You can use the `CONTINUED` subcommand to indicate that the repeating groups for each input case are contained on more than one record. You can check the value of an identification variable across records for the same input case using the `ID` subcommand.

**Summary Tables.** You can suppress the display of the table that summarizes the names, locations, and formats of the variables specified on the `DATA` subcommand using the `NOTABLE` subcommand.

### **Basic Specification**

The basic specification requires three subcommands: `STARTS`, `OCCURS`, and `DATA`.

- `STARTS` specifies the beginning column of the repeating data segments. When there are continuation records, `STARTS` can specify the ending column of the last repeating group on the first record of each input case.
- `OCCURS` specifies the number of repeating groups on each input case. `OCCURS` can specify a number if the number of repeating groups is the same for all input cases. Otherwise, `OCCURS` should specify the name of a variable whose value for each input case indicates the number of repeating groups for that case.
- `DATA` specifies names, location within the repeating segment, and format for each variable to be read from the repeated groups.

### **Subcommand Order**

- `DATA` must be the last subcommand specified on `REPEATING DATA`.
- The remaining subcommands can be named in any order.

### **Syntax Rules**

- `REPEATING DATA` can be specified only within an `INPUT PROGRAM` structure, or within a `FILE TYPE` structure with mixed or nested data. `DATA LIST`, `REPEATING DATA`, and any transformation commands used to build the output cases must be placed within the `INPUT PROGRAM` or `FILE TYPE` structure. Transformations that apply to the output cases should be specified after the `END INPUT PROGRAM` or `END FILE TYPE` command.
- `LENGTH` must be used if the last variable specified on the `DATA` subcommand is not read from the last position of each repeating group or if the length of the repeating groups varies across input cases.
- `CONTINUED` must be used if repeating groups for each input case are continued on successive records.
- The `DATA LIST` command used with `REPEATING DATA` must define all fixed-format data for the records.
- Repeating groups are usually recorded at the end of the fixed-format records, but fixed-format data may follow the repeating data in data structures such as IBM SMF and RMF records. Use the following sequence in such cases.

```
DATA LIST .../* Read the fixed-format data before repeating data
REREAD COLUMNS= .../* Skip repeating data
DATA LIST .../* Read the fixed-format data after repeating data
REPEATING DATA .../*Read repeating data
```

## Operations

- Fixed-location data specified on the `DATA LIST` are spread to each output case.
- If `LENGTH` is not specified, the program uses the default length for repeating data groups, which is determined from specifications on the `DATA` subcommand. For more information on the default length, see the `LENGTH` subcommand.

## Cases Generated

- The number of output cases generated is the number specified on the `OCCURS` subcommand. Physical record length or whether fields are non-blank does not affect the number of cases generated.
- If the number specified for `OCCURS` is nonpositive or missing, no cases are generated.

## Records Read

- If `CONTINUED` is not specified, all repeating groups are read from the first record of each input case.
- If `CONTINUED` is specified, the first continuation record is read when the first record for the input case is exhausted, that is, when the next repeating group would extend past the end of the record. The ending column for the first record is defined on `STARTS`. If the ending column is not specified on `STARTS`, the logical record length is used.
- Subsequent continuation records are read when the current continuation record is exhausted. Exhaustion of the current continuation record is detected when the next repeating group would extend past the end of the record. The ending column for continuation records is defined on `CONTINUED`. If the ending column is not specified on `CONTINUED`, the logical record length is used.
- For inline data, the record length is always 80. For data stored in a file, the record length is generally whatever was specified on the `FILE HANDLE` command or the default of 1024. Shorter records are extended with blanks when they are read. For IBM implementations, the physical record length is available and is used.

## Reading Past End of Record

If one or more fields extend past the end of the actual record, or if `CONTINUED` is specified and the ending column specified on either `STARTS` or `CONTINUED` is beyond the end of the actual record, the program takes the following action:

- For string data with format `A`, the data record is considered to be extended logically with blanks. If the entire field lies past the end of the record, the resulting value will be all blanks.
- For numeric data, a warning is issued and the resulting value is system-missing.

## Examples

### Basic Example

\* Build a file with each case representing one vehicle and

## REPEATING DATA

spread information about the household to each case.

```

INPUT PROGRAM.
DATA LIST / SEQNUM 2-4 Numpers 6-7 Numveh 9-10.
REPEATING DATA STARTS=12 /OCCURS=NUMVEH
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.

BEGIN DATA
1001 02 02 FORD      T8PONTIAC C6
1002 04 01 CHEVY    C4
1003 02 03 CADILAC C8FORD      T6VW      C4
END DATA.
LIST.

```

- Data are extracted from a file representing household records. Each input case is recorded on a single record; there are no continuation records.
- The total number of persons living in the house and number of vehicles owned by the household is recorded on each record. The first field of numbers (columns 1–4) for each record is an identification number unique to each record. The next two fields of numbers are number of persons in household and number of vehicles. The remainder of the record contains repeating groups of information about each vehicle: the make of vehicle, model, and number of cylinders.
- INPUT PROGRAM indicates the beginning of the input program and END INPUT PROGRAM indicates the end of the input program.
- DATA LIST reads the variables from the household portion of the record. All fixed-format variables are defined on DATA LIST.
- REPEATING DATA reads the information from the repeating groups and builds the output cases. Repeating groups start in column 12. The number of repeating groups for each input case is given by the value of variable NUMVEH. Three variables are defined for each repeating group: MAKE, MODEL, and NUMCYL.
- The first input record contains two repeating groups, producing two output cases in the active dataset. One output case is built from the second input record which contains information on one vehicle, and three output cases are built from the third record. The values of the fixed-format variables defined on DATA LIST are spread to every new case built in the active dataset. Six cases result, as shown below.

SEQNUM	Numpers	Numveh	MAKE	MODEL	NUMCYL
1	2	2	FORD	T	8
1	2	2	PONTIAC	C	6
2	4	1	CHEVY	C	4
3	2	3	CADILAC	C	8
3	2	3	FORD	T	6
3	2	3	VW	C	4

NUMBER OF CASES READ = 6      NUMBER OF CASES LISTED = 6

### Using REPEATING DATA With Mixed File Types

\* Use REPEATING DATA with FILE TYPE MIXED: read only type 3 records.

```

FILE TYPE MIXED RECORD=#SEQNUM 2-4.
RECORD TYPE 003.
REPEATING DATA STARTS=12 /OCCURS=3
  /DATA=MAKE 1-8(A) MODEL 9(A) NUMCYL 10.

```



```

END FILE.
END FILE TYPE.

BEGIN DATA
1001 02 02 FORD      T8PONTIAC C6
1002 04 01 CHEVY    C4
1003 02 03 CADILAC C8FORD      T6VW      C4
END DATA.
LIST.

```

- The task in this example is to read only the repeating data for records with value 003 for variable *#SEQNUM*.
- REPEATING DATA is used within a FILE TYPE structure, which specifies a mixed file type. The record identification variable *#SEQNUM* is located in columns 2–4.
- RECORD TYPE specifies that only records with value 003 for *#SEQNUM* are copied into the active dataset. All other records are skipped.
- REPEATING DATA indicates that the repeating groups start in column 12. The OCCURS subcommand indicates there are three repeating groups on each input case, and the DATA subcommand specifies names, locations, and formats for the variables in the repeating groups.
- The DATA LIST command is not required in this example, since none of the information on the input case is being spread to the output cases. However, if there were multiple input cases with value 003 for *#SEQNUM* and they did not all have three repeating groups, DATA LIST would be required to define a variable whose value for each input case indicated the number of repeating groups for that case. This variable would then be specified on the OCCURS subcommand.

### Using Transformations With REPEATING DATA

```

INPUT PROGRAM.
DATA LIST / PARENTID 1 DATE 3-6 NCHILD 8.
REPEATING DATA STARTS=9 /OCCURS=NCHILD
  /DATA=BIRTHDAY 2-5 VACDATE 7-10.
END INPUT PROGRAM.

COMPUTE AGE=DATE - BIRTHDAY.
COMPUTE VACAGE=VACDATE - BIRTHDAY.

DO IF PARENTID NE LAG(PARENTID,1) OR $CASENUM EQ 1.
COMPUTE CHILD=1.
ELSE.
COMPUTE CHILD=LAG(CHILD,1)+1.
END IF.
FORMAT AGE VACAGE CHILD (F2).

BEGIN DATA
1 1987 2 1981 1983 1982 1984
2 1988 1 1979 1984
3 1988 3 1978 1981 1981 1986 1983 1986
4 1988 1 1984 1987
END DATA.
LIST.

```

- Data are from a file that contains information on parents within a school district. Each input case is recorded on a single record; there are no continuation records.
- Each record identifies the parents by a number and indicates how many children they have. The repeating groups give the year of birth and year of vaccination for each child.

## REPEATING DATA

- REPEATING DATA indicates that the repeating groups begin in column 9. The value of *NCHILD* indicates how many repeating groups there are for each record.
- The first two COMPUTE commands compute the age for each child and age at vaccination. These transformation commands are specified outside the input program.
- Because the repeating groups do not have descriptive values, the DO IF structure computes variable *CHILD* to distinguish between the first-born child, second-born child, and so forth. The value for *CHILD* will be 1 for the first-born, 2 for the second-born, and so forth. The LIST output is shown below.

PARENTID	DATE	NCHILD	BIRTHDAY	VACDATE	AGE	VACAGE	CHILD
1	1987	2	1981	1983	6	2	1
1	1987	2	1982	1984	5	2	2
2	1988	1	1979	1984	9	5	1
3	1988	3	1978	1981	10	3	1
3	1988	3	1981	1986	7	5	2
3	1988	3	1983	1986	5	3	3
4	1988	1	1984	1987	4	3	1

NUMBER OF CASES READ = 7 NUMBER OF CASES LISTED = 7

## STARTS Subcommand

STARTS indicates the beginning location of the repeating data segment on the first record of each input case. STARTS is required and can specify either a number or a variable name.

- If the repeating groups on the first record of each input case begin in the same column, STARTS specifies a column number.
- If the repeating groups on the first record of each input case do not begin in the same column, STARTS specifies the name of a variable whose value for each input case indicates the beginning location of the repeating groups on the first record. The variable can be defined on DATA LIST or created by transformation commands that precede REPEATING DATA.
- When repeating groups are continued on multiple records for each input case, STARTS must also specify an ending location if there is room on the logical record length for more repeating groups than are contained on the first record of each input case. The ending column applies only to the first record of each input case. See the CONTINUED subcommand for an example.
- The ending column can be specified as a number or a variable name. Specifications for the beginning column and the ending column are separated by a hyphen. The values of the variable used to define the ending column must be valid values and must be larger than the starting value.
- If the variable specified for the ending column is undefined or missing for an input case, the program displays a warning message and builds no output cases from that input case. If the variable specified for the ending column on STARTS has a value that is less than the value specified for the starting column, the program issues a warning and builds output cases only from the continuation records of that input case; it does not build cases from the first record of the case.
- If the ending location is required but not supplied, the program generates output cases with system-missing values for the variables specified on the DATA subcommand and may misread all data after the first or second record in the data file (see the CONTINUED subcommand).

**Repeating Groups in the Same Location**

```

INPUT PROGRAM.
DATA LIST FILE=VEHICLE / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
REPEATING DATA STARTS=12 /OCCURS=NUMVEH
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.

```

- `STARTS` specifies column number 12. The repeating groups must therefore start in column 12 of the first record of each input case.

**Repeating Groups in Varying Locations**

```

INPUT PROGRAM.
DATA LIST FILE=VEHICLE / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
+ DO IF (SEQNUM LE 100).
+ COMPUTE FIRST=12.
+ ELSE.
+ COMPUTE FIRST=15.
+ END IF.
REPEATING DATA STARTS=FIRST /OCCURS=NUMVEH
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.

```

- This example assumes that each input case is recorded on a single record and that there are no continuation records. Repeating groups begin in column 12 for all records with sequence numbers 1 through 100 and in column 15 for all records with sequence numbers greater than 100.
- The sequence number for each record is defined as variable *SEQNUM* on the `DATA LIST` command. The `DO IF-END IF` structure creates the variable *FIRST* with value 12 for records with sequence numbers through 100 and value 15 for records with sequence numbers greater than 100.
- Variable *FIRST* is specified on the `STARTS` subcommand.

**OCCURS Subcommand**

`OCCURS` specifies the number of repeating groups for each input case. `OCCURS` is required and specifies a number if the number of groups is the same for all input cases or a variable if the number of groups varies across input cases. The variable must be defined on a `DATA LIST` command or created with transformation commands.

**Specifying the Number of Repeating Groups Using a Data Field**

```

INPUT PROGRAM.
DATA LIST / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
REPEATING DATA STARTS=12 /OCCURS=NUMVEH
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.

BEGIN DATA
1001 02 02 FORD      T8PONTIAC C6
1002 04 01 CHEVY    C4
1003 02 03 CADILAC  C8FORD      T6VW      C4
END DATA.
LIST.

```

- Data for each input case are recorded on a single record; there are no continuation records.
- The value for variable *NUMVEH* in columns 9 and 10 indicates the number of repeating groups on each record. One output case is built in the active dataset for each occurrence of a repeating group.
- In the data, *NUMVEH* has the value 2 for the first case, 1 for the second, and 3 for the third. Thus, six cases are built from these records. If the value of *NUMVEH* is 0, no cases are built from that record.

### **Specifying a Fixed Number of Repeating Groups**

\* Read only the first repeating group from each record.

```
INPUT PROGRAM.
DATA LIST FILE=VEHICLE / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
REPEATING DATA STARTS=12 /OCCURS=1
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.
LIST.
```

- Since *OCCURS* specifies that there is only one repeating group for each input case, only one output case is built from each input case regardless of the actual number of repeating groups.

## **DATA Subcommand**

*DATA* specifies a name, location within each repeating segment, and format for each variable to be read from the repeating groups. *DATA* is required and must be the last subcommand on *REPEATING DATA*.

- The specifications for *DATA* are the same as for the *DATA LIST* command.
- The specified location of the variables on *DATA* is their location within each repeating group—*not* their location within the record.
- Any input format available on the *DATA LIST* command can be specified on the *DATA* subcommand. Both FORTRAN-like and the column-style specifications can be used.

### **Example**

```
INPUT PROGRAM.
DATA LIST FILE=VEHICLE / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
REPEATING DATA STARTS=12 /OCCURS=NUMVEH
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.
LIST.
```

- Variable *MAKE* is a string variable read from positions 1 through 8 of each repeating group; *MODEL* is a single-character string variable read from position 9; and *NUMCYL* is a one-digit numeric variable read from position 10.
- The *DATA LIST* command defines variables *SEQNUM*, *NUMPERS*, and *NUMVEH*. These variables are spread to each output case built from the repeating groups.

## FILE Subcommand

REPEATING DATA always reads the file specified on its associated DATA LIST or FILE TYPE command. The FILE subcommand on REPEATING DATA explicitly specifies the name of the file.

- FILE must specify the same file as its associated DATA LIST or FILE TYPE command.

### Example

```
INPUT PROGRAM.
DATA LIST FILE=VEHICLE / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
REPEATING DATA FILE=VEHICLE /STARTS=12 /OCCURS=NUMVEH
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.
END INPUT PROGRAM.
LIST.
```

- FILE on REPEATING DATA specifically identifies the *VEHICLE* file, which is also specified on the DATA LIST command.

## ENCODING Subcommand

ENCODING specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is UTF8. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: `Locale` (the current locale setting), UTF8, UTF16, UTF16BE (big endian), UTF16LE (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- In Unicode mode, the defined width of string variables is tripled for code page and UTF-16 text data files. Use [ALTER TYPE](#) to automatically adjust the defined width of string variables.

## LENGTH Subcommand

LENGTH specifies the length of each repeating data group. The default length is the number of columns between the beginning column of the repeating data groups and the ending position of the last variable specified on DATA. (For the first record of each input case, STARTS specifies the beginning column of the repeating groups. For continuation records, repeating groups are read from column 1 by default or from the column specified on CONTINUED.)

- The specification on LENGTH can be a number or the name of a variable.
- LENGTH must be used if the last variable specified on the DATA subcommand is not read from the last position of each repeating group, or if the length of the repeating groups varies across input cases.
- If the length of the repeating groups varies across input cases, the specification must be a variable whose value for each input case is the length of the repeating groups for that case. The variable can be defined on DATA LIST or created with transformation commands.

- If the value of the variable specified on `LENGTH` is undefined or missing for an input case, the program displays a warning message and builds only one output case for that input case.

### **Example**

- \* Read only the variable `MAKE` for each vehicle.
- \* The data contain two values that are not specified on the `DATA` subcommand. The first is in position 9 of the repeating groups, and the second is in position 10.

```
INPUT PROGRAM.
DATA LIST FILE=VEHICLE / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.
REPEATING DATA STARTS=12 /OCCURS=NUMVEH /LENGTH=10
  /DATA=MAKE 1-8 (A).
END INPUT PROGRAM.
```

- `LENGTH` indicates that each repeating group is 10 columns long. `LENGTH` is required because `MAKE` is not read from the last position of each repeating group. As illustrated in previous examples, each repeating group also includes variable `MODEL` (position 9) and `NUMCYL` (position 10).
- `DATA` specifies that `MAKE` is in positions 1 through 8 of each repeating group. Positions 9 and 10 of each repeating group are skipped.

## **CONTINUED Subcommand**

`CONTINUED` indicates that the repeating groups are contained on more than one record for each input case.

- Each repeating group must be fully recorded on a single record: a repeating group cannot be split across records.
- The repeating groups must begin in the same column on all continuation records.
- If `CONTINUED` is specified without beginning and ending columns, the program assumes that the repeating groups begin in column 1 of continuation records and searches for repeating groups by scanning to the end of the record or to the value specified by `OCCURS`. [For more information, see the topic Operations on p. 1737.](#)
- If the repeating groups on continuation records do not begin in column 1, `CONTINUED` must specify the column in which the repeating groups begin.
- If there is room on the logical record length for more repeating groups than are contained on the first record of each input case, the `STARTS` subcommand must indicate an ending column for the records. The ending column on `STARTS` applies only to the first record of each input case.
- If there is room on the logical record length for more repeating groups than are contained on the continuation records of each input case, the `CONTINUED` subcommand must indicate an ending column. The ending column on `CONTINUED` applies to all continuation records.

### **Basic Example**

- \* This example assumes the logical record length is 80.

```
INPUT PROGRAM.
DATA LIST / ORDERID 1-5 NITEMS 7-8.
```

```
REPEATING DATA STARTS=10 /OCCURS=NITEMS /CONTINUED=7
  /DATA=ITEM 1-9 (A) QUANTITY 11-13 PRICE (DOLLAR7.2,1X).
END INPUT PROGRAM.
```

```
BEGIN DATA
10020 07 01-923-89 001 25.99 02-899-56 100 101.99 03-574-54 064 61.29
10020 04-780-32 025 13.95 05-756-90 005 56.75 06-323-47 003 23.74
10020 07-350-95 014 11.46
20030 04 01-781-43 010 10.97 02-236-54 075 105.95 03-655-83 054 22.99
20030 04-569-38 015 75.00
END DATA.
LIST.
```

- Data are extracted from a mail-order file. Each input case represents one complete order. The data show two complete orders recorded on a total of five records.
- The order number is recorded in columns 1 through 5 of each record. The first three records contain information for order 10020; the next two records contain information for order 20030. The second field of numbers on the first record of each order indicates the total number of items ordered. The repeating groups begin in column 10 on the first record and in column 7 on continuation records. Each repeating data group represents one item ordered and contains three variables—the item inventory number, the quantity ordered, and the price.
- DATA LIST defines variables *ORDERID* and *NITEMS* on the first record of each input case.
- STARTS on REPEATING DATA indicates that the repeating groups on the first record of each input case begin in column 10.
- OCCURS indicates that the total number of repeating groups for each input case is the value of *NITEMS*.
- CONTINUED must be used because the repeating groups are continued on more than one record for each input case. CONTINUED specifies a beginning column because the repeating groups begin in column 7 rather than in column 1 on the continuation records.
- DATA defines variables *ITEM*, *QUANTITY*, and *PRICE* for each repeating data group. *ITEM* is in positions 1–9, *QUANTITY* is in positions 11–13, and *PRICE* is in positions 14–20 and is followed by one blank column. The length of the repeating groups is therefore 21 columns. The LIST output is shown below.

ORDERID	NITEMS	ITEM	QUANTITY	PRICE
10020	7	01-923-89	1	\$25.99
10020	7	02-899-56	100	\$101.99
10020	7	03-574-54	64	\$61.29
10020	7	04-780-32	25	\$13.95
10020	7	05-756-90	5	\$56.75
10020	7	06-323-47	3	\$23.74
10020	7	07-350-95	14	\$11.46
20030	4	01-781-43	10	\$10.97
20030	4	02-236-54	75	\$105.95
20030	4	03-655-83	54	\$22.99
20030	4	04-569-38	15	\$75.00

```
NUMBER OF CASES READ =      11      NUMBER OF CASES LISTED =      11
```

### **Specifying an Ending Column on the STARTS Subcommand**

\* This example assumes the logical record length is 80.

```
INPUT PROGRAM.
DATA LIST / ORDERID 1-5 NITEMS 7-8.
REPEATING DATA STARTS=10-55 /OCCURS=NITEMS /CONTINUED=7
```

## REPEATING DATA

```

/ DATA=ITEM 1-9 (A) QUANTITY 11-13 PRICE (DOLLAR7.2,1X).
END INPUT PROGRAM.

```

```

BEGIN DATA
10020 07 01-923-89 001 25.99 02-899-56 100 101.99
10020 03-574-54 064 61.29 04-780-32 025 13.95 05-756-90 005 56.75
10020 06-323-47 003 23.74 07-350-95 014 11.46
20030 04 01-781-43 010 10.97 02-236-54 075 105.95
20030 03-655-83 054 22.99 04-569-38 015 75.00
END DATA.
LIST.

```

- Data are the same as in the previous example; however, records are entered differently. The first record for each input case contains only two repeating groups.
- DATA LIST defines variables *ORDERID* and *NITEMS* in columns 1–8 on the first record of each input case. Column 9 is blank. DATA defines variables *ITEM*, *QUANTITY*, and *PRICE* in positions 1–20 of each repeating group, followed by a blank. Thus, each repeating group is 21 columns wide. The length of the first record of each input case is therefore 51 columns: 21 columns for each of two repeating groups, plus the eight columns defined on DATA LIST, plus column 9, which is blank. The operating system's logical record length is 80, which allows room for one more repeating group on the first record of each input case. STARTS must therefore specify an ending column that does not provide enough columns for another repeating group; otherwise, the program creates an output case with missing values for the variables specified on DATA.
- STARTS specifies that the program is to scan only columns 10–55 of the first record of each input case looking for repeating data groups. It will scan continuation records beginning in column 7 until the value specified on the OCCURS subcommand is reached.

### Specifying an Ending Column on the CONTINUED Subcommand

\* This example assumes the logical record length is 80.

```

INPUT PROGRAM.
DATA LIST / ORDERID 1-5 NITEMS 7-8.
REPEATING DATA STARTS=10-55 /OCCURS=NITEMS /CONTINUED=7-55
/ DATA=ITEM 1-9 (A) QUANTITY 11-13 PRICE (DOLLAR7.2,1X).
END INPUT PROGRAM.

```

```

BEGIN DATA
10020 07 01-923-89 001 25.99 02-899-56 100 101.99
10020 03-574-54 064 61.29 04-780-32 025 13.95
10020 05-756-90 005 56.75 06-323-47 003 23.74
10020 07-350-95 014 11.46
20030 04 01-781-43 010 10.97 89-236-54 075 105.95
20030 03-655-83 054 22.99 04-569-38 015 75.00
END DATA.
LIST.

```

- The data are the same as in the previous two examples, but records are entered differently. The first record and the continuation records for each input case store only two repeating groups each.
- The operating system's logical record length is 80, which allows room for more repeating groups on all records.



- `STARTS` specifies that the program is to scan only columns 10-55 of the first record of each input case looking for repeating data groups.
- `CONTINUED` specifies that the program is to scan only columns 7-55 of all continuation records.

## ***ID Subcommand***

`ID` compares the value of an identification variable across records of the same input case. `ID` can be used only when `CONTINUED` is specified. The identification variable must be defined on a `DATA LIST` command and must be recorded on all records in the file.

- The `ID` subcommand has two specifications: the location of the variable on the continuation records and the name of the variable (as specified on the `DATA LIST` command). The specifications must be separated from each other by an equals sign.
- The format specified on the `ID` subcommand must be the same as the format specified for the variable on `DATA LIST`. However, the location can be different on the continuation records.
- If the values of the identification variable are not the same on all records for a single input case, the program displays an error message and stops reading data.

### ***Example***

```
INPUT PROGRAM.
DATA LIST / ORDERID 1-5 NITEMS 7-8.
REPEATING DATA STARTS=10-50 /OCCURS=NITEMS
 /CONTINUED=7 /ID=1-5=ORDERID
 /DATA=ITEM 1-9 (A) QUANTITY 11-13 PRICE 15-20 (2).
END INPUT PROGRAM.
```

```
BEGIN DATA
10020 04 45-923-89 001 25.9923-899-56 100 101.99
10020 63-780-32 025 13.9554-756-90 005 56.75
20030 03 45-781-43 010 10.9789-236-54 075 105.95
20030 32-569-38 015 75.00
END DATA.
LIST.
```

- The order number in the data is recorded in columns 1-5 of each record.
- `ORDERID` is defined on the `DATA LIST` command as a five-column integer variable. The first specification on the `ID` subcommand must therefore specify a five-column integer variable. The location of the variable can be different on continuation records.

## ***TABLE and NOTABLE Subcommands***

`TABLE` displays a table summarizing all variables defined on the `DATA` subcommand. The summary table lists the names, locations, and formats of the variables and is identical in format to the summary table displayed by the `DATA LIST` command. `NOTABLE` suppresses the table. `TABLE` is the default.

### ***Example***

```
INPUT PROGRAM.
```

---

*REPEATING DATA*

```
DATA LIST FILE=VEHICLE / SEQNUM 2-4 NUMPERS 6-7 NUMVEH 9-10.  
REPEATING DATA STARTS=12 /OCCURS=NUMVEH /NOTABLE  
  /DATA=MAKE 1-8 (A) MODEL 9 (A) NUMCYL 10.  
END INPUT PROGRAM.
```

- NOTABLE suppresses the display of the summary table.

# REPORT

REPORT is available in the Statistics Base option.

```
REPORT [/FORMAT={MANUAL } ] [NOLIST ] [ALIGN({LEFT } )] [TSPACE({1})]
      {AUTOMATIC} {LIST[(n)]} {CENTER} {n}
      {RIGHT }
      [CHDSPACE({1})] [FTSPACE({1})] [SUMSPACE({1})] [COLSPACE({4})]
      {n} {n} {n} {n}
      [BRKSPACE({ 1 } )] [LENGTH({1,length})] [MARGINS({1,width})]
      { n } {t,b } {l,r }
      {-1†} {*,* } {*,* }
      [CHALIGN({TOP } )] [UNDERSCORE({OFF})] [PAGE1({1})] [MISSING {'. '}]
      {BOTTOM†} {ON†} {n} {s' }
      [ONEBREAKCOLUMN {OFF**}] [INDENT {2**}] [CHWRAP {OFF**}] [PREVIEW {OFF**}]
      {ON } {n } {ON } {ON }

[/OUTFILE=file]
[/STRING=stringname (varname[(width)] [(BLANK)] ['literal'])]
/VARIABLES=varname ({VALUE}) [+ varname({VALUE})] ['col head'] [option list]
      {LABEL} {LABEL}
      {DUMMY} {DUMMY}
```

where option list can contain any of the following:

```
(width) (OFFSET({0 } )) ((LEFT } )
      {n } {CENTER†}
      {CENTER†} {RIGHT }

[/MISSING={VAR } ]
      {NONE }
      {LIST[([varlist] [(1)])]}
      {n}

[/TITLE={(LEFT } ) 'line1' 'line2'... ] [FOOTNOTE={(LEFT } ) 'line1' 'line2'... ]
      {CENTER} {CENTER}
      {RIGHT } {RIGHT }
      [ ] [PAGE] [ ] [DATE] [ ] [var]

[/BREAK=varlist ['col head'] [option list]]
```

where option list can contain any of the following:

```
(width) ({VALUE } ) ((NOTOTAL}) ((SKIP({1 } ))
      {LABEL†} {TOTAL } {n}
      {PAGE[(RESET)]}
      (OFFSET({0 } )) (UNDERSCORE([varlist])) ((LEFT } ) ({NONAME})
      {n } {CENTER†} {CENTER†} {NAME }
      {CENTER†} {RIGHT }

[/SUMMARY=function...['summary title'][(break col #)] [SKIP({0})]
      {n}
      or
[/SUMMARY=PREVIOUS[({1})]
      {n}
```

where function is

```
aggregate [(varname[({PLAIN } )])[(d)][varname...]]
      {format††}
      or
composite(argument) [(report col[({PLAIN } )])[(d)]]
      {format††}
```

\*\*Default if the keyword is omitted.

†Default if `FORMAT=AUTOMATIC`.

††Any printable output format is valid. See `FORMATS`.

*Aggregate functions:*

VALIDN	VARIANCE	PLT(n)
SUM	KURTOSIS	PIN(min,max)
MIN	SKEWNESS	FREQUENCY(min,max)
MAX	MEDIAN(min,max)	PERCENT(min,max)
MEAN	MODE(min,max)	
STDDEV	PGT(n)	

*Composite functions:*

```

DIVIDE(arg1 arg2 [factor])
MULTIPLY(arg1...argn)
PCT(arg1 arg2)
SUBTRACT(arg1 arg2)
ADD(arg1...argn)
GREAT(arg1...argn)
LEAST(arg1...argn)
AVERAGE(arg1...argn)

```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

**Example**

```

REPORT FORMAT=LIST
  /VARIABLES=PRODUCT (LABEL) ' ' 'Retail' 'Products'
    SALES 'Annual' 'Sales' '1981'
  /BREAK=DEPT 'Department' (LABEL)
  /SUMMARY=VALIDN (PRODUCT) MEAN (SALES).

```

## Overview

REPORT produces case listings and summary statistics and gives you considerable control over the appearance of the output. REPORT calculates all the univariate statistics available in `DESCRIPTIVES` and the statistics and subpopulation means available in `MEANS`. In addition, REPORT calculates statistics not directly available in any other procedure, such as computations involving aggregated statistics.

REPORT provides complete report format defaults but also lets you customize a variety of table elements, including column widths, titles, footnotes, and spacing. Because REPORT is so flexible and the output has so many components, it is often efficient to preview report output using a small number of cases until you find the format that best suits your needs.

### **Basic specification**

The basic specification depends on whether you want a listing report or a summary report. A listing report without subgroup classification requires `FORMAT` and `VARIABLES`. A listing report with subgroup classification requires `FORMAT`, `VARIABLES`, and `BREAK`. A summary report requires `VARIABLES`, `BREAK`, and `SUMMARY`.

**Listing reports.** `FORMAT=LIST` and `VARIABLES` with a variable list are required. Case listings are displayed for each variable named on `VARIABLES`. There are no break groups or summary statistics unless `BREAK` or `SUMMARY` is specified.

**Summary reports.** `VARIABLES`, `BREAK`, and `SUMMARY` are required. The report is organized according to the values of the variable named on `BREAK`. The variable named on `BREAK` must be named on a preceding `SORT CASES` command. Specified statistics are computed for the variables specified on `VARIABLES` for each subgroup defined by the break variables.

### **Subcommand order**

The following order must be observed among subcommands when they are used:

- `FORMAT` must precede all other subcommands.
- `VARIABLES` must precede `BREAK`.
- `OUTFILE` must precede `BREAK`.
- Each `SUMMARY` subcommand must immediately follow its associated `BREAK`. Multiple `SUMMARY` subcommands associated with the same `BREAK` must be specified consecutively.
- `TITLE` and `FOOTNOTE` can appear anywhere after `FORMAT` except between `BREAK` and `SUMMARY`.
- `MISSING` must follow `VARIABLES` and precede the first `BREAK`.
- `STRING` must precede `VARIABLES`.

### **Syntax rules**

- Only one each of the `FORMAT`, `STRING`, `VARIABLES`, and `MISSING` subcommands is allowed.
- To obtain multiple break groups, use multiple `BREAK` subcommands.
- To obtain multiple summaries for a break level, specify multiple `SUMMARY` subcommands for the associated `BREAK`.
- Keywords on `REPORT` subcommands have default specifications that are in effect if the keyword is not specified. Specify keywords only when you wish to change a default.
- Keywords are enclosed in parentheses if the subcommand takes variable names as arguments.

### **Operations**

- `REPORT` processes cases sequentially. When the value of a break variable changes, `REPORT` displays a statistical summary for cases processed since the last set of summary statistics was displayed. Thus, the file must be sorted in order on the break variable or variables.
- The maximum width and page length of the report are determined by the `SET` command.

- If a column is not wide enough to display numeric values, `REPORT` first rounds decimal digits, then converts to scientific notation if possible, and then displays asterisks. String variables that are wider than the column are truncated.
- The format used to display values in case listings is controlled by the dictionary format of the variable. Each statistical function in `REPORT` has a default format.

### Limitations

- Maximum 500 variables per `VARIABLES` subcommand. You can specify more than 500 variables if you stack them. [For more information, see the topic VARIABLES subcommand on p. 1758.](#)
- Maximum 10 dummy variables per `VARIABLES` subcommand.
- Maximum 20 `MODE` and `MEDIAN` requests per `SUMMARY` subcommand.
- Maximum 20 `PGT`, `PLT`, and `PIN` requests per `SUMMARY` subcommand.
- Maximum 50 strings per `STRING` subcommand.
- The length of titles and footnotes cannot exceed the report width.
- The length of string variables created on `STRING` cannot exceed the page width.
- There is no fixed limit on the number of `BREAK` and `SUMMARY` subcommands. However, the page width limits the number of variables that can be displayed and thereby limits the number of break variables.
- The maximum width of a report is 255 characters.
- The number of report variables that can be specified depends upon the width of the report, the width of the variable columns, and the number of `BREAK` subcommands.
- Maximum 50 variables for the `FREQUENCY` or `PERCENT` functions.
- Memory requirements significantly increase if `FREQUENCY`, `PERCENT`, `MEDIAN`, or `MODE` is requested for variables with a wide range of values. The amount of workspace required is  $20 + 8 * (\text{max} - \text{min} + 1)$  bytes per variable per function per break. If the same range is used for different statistics for the same variable, only one set of cells is collected. For example, `FREQUENCY(1,100)(VARA) PERCENT(1,100)(VARA)` requires only 820 bytes.
- If `TOTAL` is in effect, workspace requirements are almost doubled.
- Memory requirements also increase if value labels are displayed for variables with many value labels. The amount of workspace required is  $4 + 24 * n$  bytes per variable, where  $n$  is the number of value labels specified for the variable.

## Examples

```

SORT CASES BY DEPT.
REPORT FORMAT=LIST
  /VARIABLES=PRODUCT (LABEL) ' ' 'Retail' 'Products'
    SALES 'Annual' 'Sales' '1981'
  /BREAK=DEPT 'Department' (LABEL)
  /SUMMARY=VALIDN (PRODUCT) MEAN (SALES) 'No.Sold,Mean Sales'.

```

- This report is a listing of products and sales by department. A summary of the total number of products sold and the average sales by department is also produced.

- Cases are first sorted by *DEPT* so that cases are grouped by department for the case listing and for the calculation of statistics.
- *FORMAT* requests a report that lists individual cases within each break group.
- *VARIABLES* specifies *PRODUCT* and *SALES* as the report variables. Keyword *LABEL* requests that the case listings for *PRODUCT* display value labels instead of values. Three-line column headings are provided for each report column. The first line of the column heading is blank for the variable *PRODUCT*.
- *BREAK* identifies *DEPT* as the break variable and provides a one-line column title for the break column. *LABEL* displays the value label instead of the value itself.
- *SUMMARY* calculates the valid number of cases for *PRODUCT* and the mean of *SALES* for each value of *DEPT*. A title is provided for the summary line to override the default title, *VALIDN*.

## Defaults

**Column heads.** *REPORT* uses variable labels as default column heads; if no variable labels have been specified, variable names are used. If *ONEBREAKCOLUMN* is *ON*, the default head for the first *BREAK* subcommand is used.

**Column widths.** Default column widths are determined by *REPORT*, using the maximum of the following for each column:

- The widest print format in the column, whether it is a variable print format or a summary print format.
- The width of any temporary variable defined with the *STRING* subcommand on *REPORT*.
- If a column heading is assigned, the length of the longest title line in the heading when *CHWRAP* is off, and the longest word in the title when *CHWRAP* is on. Underscores, which are removed on printing, can be used to create longer words in the title.
- When no column heading is specified, the length of the longest word in the variable label, or the length of the variable name.
- If you specify *LABEL* on *VARIABLES* or *BREAK*, the length of the variable's longest value label. If *FORMAT=MANUAL* is in effect, 20 is the maximum value used for this criterion.
- The minimum column width is 8 when *FORMAT=MANUAL*; it can be less when *FORMAT=AUTOMATIC*.

**Automatic fit.** When the above criteria for column width result in a report that is too wide for the report margins, *FORMAT=AUTOMATIC* shrinks the report. *AUTOMATIC* performs the following two steps sequentially, stopping as soon as the report fits within the margins:

1. *AUTOMATIC* reduces intercolumn spacing incrementally until it reaches a minimum intercolumn space of 1. It will never reduce it to 0.
2. *AUTOMATIC* shortens widths for strings specified on the *STRING* subcommand or for value label strings when the *LABEL* option is specified. It begins with the longest string if that string is at least 15 characters wide and shortens the column width as much as needed (up to 40% of its length), wrapping the string within the new width. If necessary, it repeats the step, using different defined strings. It will not shorten the column width of the same string twice.

REPORT does *not* implement the automatic fit unless AUTOMATIC is specified on the FORMAT subcommand.

**AUTOMATIC versus MANUAL defaults.** Many default settings depend on whether you specify AUTOMATIC or MANUAL on FORMAT. The following table shows the defaults according to either of the specifications.

Table 207-1  
Keyword default settings

Subcommand	Keyword	Default for AUTOMATIC	Default for MANUAL
FORMAT	ALIGN	left	left
	BRKSPACE		
	summary report	1	1
	listing report	-1	1
	CHALIGN	bottom	top
	CHDSPACE	1	1
	COLSPACE	4	4
	FTSPACE	1	1
	LENGTH	1,system length	1,system length
	LIST NOLIST	NOLIST	NOLIST
	MARGINS	1,system width	1,system width
	MISSING	.	.
	PAGE1	1	1
	SUMSPACE	1	1
	TSPACE	1	1
	UNDERSCORE	on	off
	ONEBREAKCOLUMN	off	off
	INDENT <sup>1</sup>	2	2
	CHWRAP	off	off
	PREVIEW	off	off
VARIABLES	LABEL VALUE DUMMY	VALUE	VALUE
	LEFT CENTER RIGHT	CENTER <sup>2</sup>	RIGHT for numbers LEFT for strings
	OFFSET	CENTER	0
BREAK	LABEL VALUE	LABEL	VALUE
	LEFT CENTER RIGHT	CENTER <sup>2</sup>	RIGHT for numbers LEFT for strings
	NAME NONAME	NONAME	NONAME
	OFFSET	CENTER <sup>3</sup>	0
	PAGE	off	off
	SKIP	1	1
	TOTAL NOTOTAL	NOTOTAL	NOTOTAL
UNDERSCORE	off	off	
SUMMARY	PREVIOUS	1	1



Subcommand	Keyword	Default for AUTOMATIC	Default for MANUAL
	SKIP	0	0

<sup>1</sup> No effect when ONEBREAKCOLUMN is on.

<sup>2</sup> LEFT when ONEBREAKCOLUMN is on.

<sup>3</sup> 0 when ONEBREAKCOLUMN is on.

## Options

**Format.** REPORT provides full format defaults and offers you optional control over page length, vertical spacing, margin and column widths, page titles, footnotes, and labels for statistics. The maximum width and length of the report are controlled by specifications on the SET command. The FORMAT subcommand on REPORT controls how the report is laid out on a page and whether case listings are displayed. The VARIABLES subcommand specifies the variables that are listed or summarized in the report (**report variables**) and controls the titles, width, and contents of report columns. The BREAK subcommand specifies the variables that define groups (**break variables**) and controls the titles, width, and contents of break columns. SUMMARY specifies statistics and controls the titles and spacing of summary lines. The TITLE and FOOTNOTE subcommands control the specification and placement of multiple-line titles and footnotes. STRING concatenates variables to create temporary variables that can be specified on VARIABLES or BREAK.

**Output file.** You can direct reports to a file separate from the file used for the rest of the output from your session using the OUTFILE subcommand.

**Statistical display.** The statistical display is controlled by the SUMMARY subcommand. Statistics can be calculated for each category of a break variable and for the group as a whole. Available statistics include mean, variance, standard deviation, skewness, kurtosis, sum, minimum, maximum, mode, median, and percentages. Composite functions perform arithmetic operations using two or more summary statistics calculated on single variables.

**Missing values.** You can override the default to include user-missing values in report statistics and listings with the MISSING subcommand. You can also use FORMAT to define a missing-value symbol to represent missing data.

## FORMAT subcommand

FORMAT controls the overall width and length of the report and vertical spacing. Keywords and their arguments can be specified in any order.

- `MANUAL` and `AUTOMATIC` are alternatives. The default is `MANUAL`.
- `LIST` and `NOLIST` are alternatives. The default is `NOLIST`.

<b>MANUAL</b>	<i>Default settings for manual format.</i> <code>MANUAL</code> displays values for break variables, right-justifies numeric values and their column headings, left-justifies value labels and string values and their column headings, top-aligns and does not underscore column headings, extends column widths to accommodate the variable's longest value label (but not the longest word in the variable label) up to a width of 20, and generates an error message when a report is too wide for its margins. <code>MANUAL</code> is the default.
<b>AUTOMATIC</b>	<i>Default settings for automatic format.</i> <code>AUTOMATIC</code> displays labels for break variables, centers all data, centers column headings but left-justifies column headings if value labels or string values exceed the width of the longest word in the heading, bottom-aligns and underscores column headings, extends column widths to accommodate the longest word in a variable label or the variable's longest value label, and shrinks a report that is too wide for its margins.
<b>LIST(n)</b>	<i>Individual case listing.</i> The values of all variables named on <code>VARIABLES</code> are displayed for each case. The optional <i>n</i> inserts a blank line after each <i>n</i> cases. By default, no blank lines are inserted. Values for cases are listed using the default formats for the variables.
<b>NOLIST</b>	<i>No case listing.</i> This is the default.
<b>PAGE(n)</b>	<i>Page number for the first page of the report.</i> The default is 1.
<b>LENGTH(t,b)</b>	<i>Top and bottom line numbers of the report.</i> You can specify any numbers to define the report page length. By default, the top of the report begins at line 1, and the bottom of the report is the last line of the system page. You can use an asterisk for <i>t</i> or <i>b</i> to indicate a default value. If the specified length does not allow even one complete line of information to be displayed, <code>REPORT</code> extends the length specification and displays a warning.
<b>MARGINS(l,r)</b>	<i>Columns for the left and right margins.</i> The right column cannot exceed 255. By default, the left margin is display column 1 and the right margin is the rightmost display column of the system page. You can use an asterisk for <i>l</i> or <i>r</i> to indicate a default value.
<b>ALIGN</b>	<i>Placement of the report relative to its margins.</i> <code>LEFT</code> , <code>CENTER</code> , or <code>RIGHT</code> can be specified in the parentheses following the keyword. <code>LEFT</code> left-justifies the report. <code>CENTER</code> centers the report between its margins. <code>RIGHT</code> right-justifies the report. The default is <code>LEFT</code> .
<b>COLSPACE(n)</b>	<i>Number of spaces between each column.</i> The default is 4 or the average number of spaces that will fit within report margins, whichever is less. When <code>AUTOMATIC</code> is in effect, <code>REPORT</code> overrides the specified column spacing if necessary to fit the report between its margins.
<b>CHALIGN</b>	<i>Alignment of column headings.</i> Either <code>TOP</code> or <code>BOTTOM</code> can be specified in the parentheses following the keyword. <code>TOP</code> aligns all column headings with the first, or top, line of multiple-line headings. <code>BOTTOM</code> aligns headings with the last, or bottom, line of multiple-line headings. When <code>AUTOMATIC</code> is in effect, the default is <code>BOTTOM</code> ; when <code>MANUAL</code> is in effect, the default is <code>TOP</code> .
<b>UNDERSCORE</b>	<i>Underscores for column headings.</i> Either <code>ON</code> or <code>OFF</code> can be specified in the parentheses following the keyword. <code>ON</code> underscores the bottom line of each column heading for the full width of the column. <code>OFF</code> does not underscore column headings. The default is <code>ON</code> when <code>AUTOMATIC</code> is in effect and <code>OFF</code> when <code>MANUAL</code> is in effect.
<b>TSPACE(n)</b>	<i>Number of blank lines between the report title and the column heads.</i> The default is 1.
<b>CHDSPACE(n)</b>	<i>Number of blank lines beneath the longest column head.</i> The default is 1.

<b>BRKSPACE(n)</b>	<i>Number of blank lines between the break head and the next line. The next line is a case if LIST is in effect or the first summary line if NOLIST is in effect. BRKSPACE (-1) places the first summary statistic or the first case listing on the same line as the break value. When a summary line is placed on the same line as the break value, the summary title is suppressed. When AUTOMATIC is in effect, the default is -1; when MANUAL is in effect, it is 1.</i>
<b>SUMSPACE(n)</b>	<i>Number of blank lines between the last summary line at the lower break level and the first summary line at the higher break level when they break simultaneously. SUMSPACE also controls spacing between the last listed case and the first summary line if LIST is in effect. The default is 1.</i>
<b>FTSPACE(n)</b>	<i>Minimum number of blank lines between the last listing on the page and the footnote. The default is 1.</i>
<b>MISSING 's'</b>	<i>Missing-value symbol. The symbol can be only one character and represents both system- and user-missing values. The default is a period.</i>
<b>ONEBREAKCOLUMN</b>	<i>Display subgroups defined on multiple BREAK subcommands in a single column. You can specify OFF or ON in parentheses after the keyword. The default is OFF. When ONEBREAKCOLUMN is ON, it applies to all BREAK subcommands. <a href="#">For more information, see the topic BREAK subcommand on p. 1762.</a></i>
<b>INDENT(n)</b>	<i>Indentation of break values and summary titles of each successive subgroup defined by one BREAK subcommand in a single break column. INDENT is effective only when ONEBREAKCOLUMN is on. Multiple variables specified on one BREAK subcommand are indented as a block. The default specification is 2. When ONEBREAKCOLUMN is OFF, specification on INDENT is ignored.</i>
<b>CHWRAP</b>	<i>Automatically wrap user-specified column heads. You can specify OFF or ON in parentheses after the keyword. The default is OFF. When CHWRAP is ON, user-specified heads for either break or variable columns are wrapped. If multiple lines are specified for a head, each line is wrapped, if necessary, independent of other lines. To prevent wrapping at blanks, use the underscore character ( _ ) to signify a hard blank in your head specification. The underscore serves as a hard blank only in user-specified heads and only when CHWRAP is ON. The underscore does not appear in the printed heading.</i>
<b>PREVIEW</b>	<i>Display the first page of output only. You can specify OFF or ON either in parentheses or with one blank space separating the specification from the keyword. The default is OFF. When PREVIEW is ON, the program stops processing after the first page for you to quickly check the format of your report.</i>

## ***OUTFILE subcommand***

OUTFILE directs the report to a file separate from the file used for the rest of the output from your session. This allows you to print the report without having to delete the extraneous material that would be present in the output.

- OUTFILE must follow FORMAT and must precede BREAK.
- You can append multiple reports to the same file by naming the same file on the OUTFILE subcommand for each REPORT command.

Note: The first time a particular file is named on the OUTFILE subcommand within a job or session, the file is created or, if it already exists, it is erased. It remains open until the job or session finishes. Later uses of the same file in REPORT append their output to it. Attempts to erase the file may fail.

**Example**

```
REPORT FORMAT=AUTOMATIC LIST
  /OUTFILE=PRSNLRPT
  /VARIABLES=LNAME AGE TENURE JTENURE SALARY
  /BREAK=DIVISION
  /SUMMARY=MEAN.
```

```
REPORT FORMAT=AUTOMATIC
  /OUTFILE=PRSNLRPT
  /VARIABLES=LNAME AGE TENURE JTENURE SALARY
  /BREAK=DIVISION
  /SUMMARY=MEAN
  /SUMMARY=MIN
  /SUMMARY=MAX.
```

- Both a listing report and a summary report are written to file *PRSNLRPT*.

**VARIABLES subcommand**

The required **VARIABLES** subcommand names the variables to be listed and summarized in the report. You can also use **VARIABLES** to control column titles, column widths, and the contents of report columns.

- The minimum specification on **VARIABLES** is a list of report variables. The number of variables that can be specified is limited by the system page width.
- Each report variable defines a report column. The value of the variable or an aggregate statistic calculated for the variable is displayed in that variable's report column.
- Variables are assigned to columns in the order in which they are named on **VARIABLES**.
- Variables named on **BREAK** can also be named on **VARIABLES**.
- When **FORMAT=LIST**, variables can be stacked in a single column by linking them with plus signs (+) on the **VARIABLES** subcommand. If no column heading is specified, **REPORT** uses the default heading from the first variable on the list. Only values from the first variable in the column are used to calculate summaries.
- Optional specifications apply only to the immediately preceding variable or list of variables implied by the **TO** keyword. Options can be specified in any order.
- All optional specifications except column headings must be enclosed in parentheses; column headings must be enclosed in quotes.

**Column contents**

The following options can be used to specify the contents of the report column for each variable:

<b>(VALUE)</b>	<i>Display the values of the variable.</i> This is the default.
<b>(LABEL)</b>	<i>Display value labels.</i> If value labels are not defined, values are displayed.
<b>(DUMMY)</b>	<i>Display blank spaces.</i> <b>DUMMY</b> defines a report column for a variable that does not exist in the active dataset. Dummy variables are used to control spacing or to reserve space for statistics computed for other variables. Do not name an existing variable as a dummy variable.

- `VALUE` and `LABEL` have no effect unless `LIST` has been specified on the `FORMAT` subcommand.
- When `AUTOMATIC` is in effect, value labels or string values are centered in the column based on the length of the longest string or label; numeric values are centered based on the width of the widest value or summary format. When `MANUAL` is in effect, value labels or string values are left-justified in the column and numeric values are right-justified. (See the `OFFSET` keyword.)

## ***Column heading***

The following option can be used to specify a heading for the report column:

**'column heading'**      *Column heading for the preceding variable.* The heading must be enclosed in quotes. If no column heading is specified, the default is the variable label or, if no variable label has been specified, the variable name.

- To specify multiple-line headings, enclose each line in quotes, using the conventions for strings. The specifications for title lines should be separated by at least one blank.
- Default column headings wrap for as many lines as are required to display the entire label. If `AUTOMATIC` is in effect, user-specified column headings appear exactly as specified, even if the column width must be extended. If `MANUAL` is in effect, user-specified titles wrap to fit within the column width.

## ***Column heading alignment***

The following options can be used to specify how column headings are aligned:

**(LEFT)**                      *Left-aligned column heading.*  
**(CENTER)**                    *Centered column heading.*  
**(RIGHT)**                     *Right-aligned column heading.*

- If `AUTOMATIC` is in effect, column headings are centered within their columns by default. If value labels or string values exceed the width of the longest word in the heading, the heading is left-justified.
- If `MANUAL` is in effect, column headings are left-justified for value labels or string values and right-justified for numeric values by default.

## Column format

The following options can be used to specify column width and adjust the position of the column contents:

- |                 |   |
|-----------------|---|
| <b>(width)</b>  | <i>Width for the report column.</i> If no width is specified for a variable, REPORT determines a default width using the criteria described under <a href="#">Defaults</a> . If you specify a width that is not wide enough to display numeric values, REPORT first rounds decimal digits, then converts to scientific notation if possible, and then displays asterisks. Value labels or string values that exceed the width are wrapped.  |
| <b>(OFFSET)</b> | <i>Position of the report column contents.</i> The specification is either <i>n</i> or CENTER specified in parentheses. OFFSET( <i>n</i> ) indicates the number of spaces to offset the contents from the left for value labels or string values, and from the right for numeric values. OFFSET(CENTER) centers contents within the center of the column. If AUTOMATIC is in effect, the default is CENTER. If MANUAL is in effect, the default is 0. Value labels and string values are left-justified and numeric values are right-justified. |

### Example

```
/VARIABLES=V1 TO V3(LABEL) (15)
V4 V5 (LABEL)(OFFSET (2))(10)
SEPI (DUMMY) (2) ' '
V6 'Results using' "Lieben's Method" 'of Calculation'
```

- The width of the columns for variables *V1* through *V3* is 15 each. Value labels are displayed for these variables in the case listing.
- The column for variable *V4* uses the default width. Values are listed in the case listing.
- Value labels are displayed for variable *V5*. The column width is 10. Column contents are offset two spaces from the left.
- *SEPI* is a dummy variable. The column width is 2, and there is at least one space on each side of *SEPI*. Thus, there are at least four blanks between the columns for *V5* and *V6*. *SEPI* is given a null title to override the default column title *SEPI*.
- *V6* has a three-line title. Its column uses the default width, and values are listed in the case listing.

## STRING subcommand

STRING creates a temporary string variable by concatenating variables and user-specified strings. These variables exist only within the REPORT procedure.

- The minimum specification is a name for the string variable followed by a variable name or a user-specified string enclosed in parentheses.
- The name assigned to the string variable must be unique.
- Any combination of string variables, numeric variables, and user-specified strings can be used in the parentheses to define the string.
- Keyword TO cannot be used within the parentheses to imply a variable list.
- More than one string variable can be defined on STRING.

- If a case has a missing value for a variable within the parentheses, the variable passes the missing value to the temporary variable without affecting other elements specified.
- A string variable defined in REPORT cannot exceed the system page width.
- String variables defined on STRING can be used on VARIABLES or BREAK.

The following options can be used to specify how components are to be concatenated:

<b>(width)</b>	<i>Width of the preceding variable within the string.</i> The default is the dictionary width of the variable. The maximum width for numeric variables within the string definition is 16. The maximum width for a string variable is the system page width. If the width specified is less than that required by the value, numeric values are displayed as asterisks and string values are truncated. If the width exceeds the width of a value, numeric values are padded with zeros on the left and string values are padded with blanks on the right.
<b>(BLANK)</b>	<i>Left-pad values of the preceding numeric variable with blanks.</i> The default is to left-pad values of numeric variables with zeros. If a numeric variable has a dollar or comma format, it is automatically left-padded with blanks.
<b>'literal'</b>	<i>User-specified string.</i> Any combination of characters can be specified within quotes.

### Example

```
/STRING=JOB1 (AVAR NVAR)
      JOB2 (AVAR (2) NVAR (3))
      JOB3 (AVAR (2) NVAR (BLANK) (4))
```

- STRING defines three string variables to be used within the report.
- Assume that *AVAR* is a string variable read from a four-column field using keyword FIXED on DATA LIST and that *NVAR* is a computed numeric variable with the default format of eight columns with two implied decimal places.
- If a case has value KJ for *AVAR* and value 241 for *NVAR*, *JOB1* displays the value 'KJ 00241.00', *JOB2* the value 'KJ241', and *JOB3* the value 'KJ 241'. If *NVAR* has the system-missing value for a case, *JOB1* displays the value 'KJ'.

### Example

```
/STRING=SOCSEC (S1 '-' S2 '-' S3)
```

- STRING concatenates the three variables *S1*, *S2*, and *S3*, each of which contains a segment of the social security number.
- Hyphens are inserted between the segments when the values of *SOCSEC* are displayed.
- This example assumes that the variables *S1*, *S2*, and *S3* were read from three-column, two-column, and four-column fields respectively, using the keyword FIXED on DATA LIST. These variables would then have default format widths of 3, 2, and 4 columns and would not be left-padded with zeros.

## ***BREAK subcommand***

**BREAK** specifies the variables that define the subgroups for the report, or it specifies summary totals for reports with no subgroups. **BREAK** also allows you to control the titles, width, and contents of break columns and to begin a new page for each level of the break variable.

- A break occurs when any one of the variables named on **BREAK** changes value. Cases must be sorted by the values of all **BREAK** variables on all **BREAK** subcommands.
- The **BREAK** subcommand must precede the **SUMMARY** subcommand that defines the summary line for the break.
- A break column is reserved for each **BREAK** subcommand if **ONEBREAKCOLUMN** is **OFF** (the default).
- To obtain multiple break levels, specify multiple break variables on a **BREAK** subcommand.
- If more than one variable is specified on a **BREAK** subcommand, a single break column is used. The value or value label for each variable is displayed on a separate line in the order in which the variables are specified on **BREAK**. The first variable specified changes most slowly. The default column width is the longest of the default widths for any of the break variables.
- To obtain summary totals without any break levels, use keyword **TOTAL** in parentheses on **BREAK** without listing any variables. **TOTAL** must be specified on the first **BREAK** subcommand.
- When **MISSING=VAR** is specified, user-missing values are displayed in case listings but are not included in summary statistics. When **NONE** is specified, user-missing values are ignored. System-missing values are displayed as missing in case and break listings.
- Optional specifications apply to all variables in the break column and to the break column as a whole. Options can be specified in any order following the last variable named.
- All optional specifications except column headings must be enclosed in parentheses; column headings must be enclosed in quotes.

### ***Column contents***

The following can be used to specify the contents of the break column:

- |                |  |
|----------------|--|
| <b>(VALUE)</b> | <i>Display values of the break variables.</i>  |
| <b>(LABEL)</b> | <i>Display value labels.</i> If no value labels have been defined, values are displayed. |

- The value or label is displayed only once for each break change but it is repeated at the top of the page in a multiple-page break group.
- When **AUTOMATIC** is in effect, the default is **LABEL**; when **MANUAL** is in effect, the default is **VALUE**.
- When **AUTOMATIC** is in effect, the value or label is centered in the column. When **MANUAL** is in effect, value labels and string values are left-justified and numeric values are right-justified. Keywords **OFFSET**, **ONEBREAKCOLUMN**, and **INDENT** can also affect positioning.



## Column heading

The following option specifies headings used for the break column.

**'column heading'**      *Column heading for the break column.* The heading must be included in quotes. The default heading is the variable label of the break variable or, if no label has been defined, the variable name. If the break column is defined by more than one variable, the label or name of the first variable is used. If `ONEBREAKCOLUMN` is ON, the specified or implied column heading for the first `BREAK` subcommand is used.

- To specify multiple-line headings, enclose each line in a set of quotes, following the conventions for strings. Separate the specifications for heading lines with at least one blank.
- Default column headings wrap for as many lines as are required to display the entire label.
- User-specified column headings appear exactly as specified if `CHWRAP` is OFF (the default). If `CHWRAP` is ON, any user-defined line longer than the specified or default column width is automatically wrapped.

## Column heading alignment

The following options can be used to specify how column headings are aligned:

**(LEFT)**                      *Left-aligned column heading.*  
**(CENTER)**                    *Centered column heading.*  
**(RIGHT)**                     *Right-aligned column heading.*

- When `AUTOMATIC` is in effect, column headings are centered within their columns by default. If value labels or string values exceed the width of the longest word in the heading, the heading is left-justified.
- When `MANUAL` is in effect, column headings are left-justified for value labels or string values and right-justified for numeric values.
- When `ONEBREAKCOLUMN` is ON, all column contents are left aligned. Specifications of `CENTER` and `RIGHT` on `BREAK` are ignored.

## Column format

The following options can be used to format break columns:

**(width)**                      *Column width for the break column.* If no width is specified for a variable, `REPORT` determines a default width using the criteria described under [Defaults](#). If `ONEBREAKCOLUMN` is ON, the column width specified or implied by the first `BREAK` subcommand is used. If you specify a width that is not wide enough to display numeric values, `REPORT` first rounds decimal digits, then converts them to scientific notation if possible, and then displays asterisks. Value labels or string values that exceed the width are wrapped.

**(OFFSET)**                     *Position of the break column contents.* The specification is either `n` or `CENTER` specified in parentheses. `OFFSET (n)` indicates the number of spaces to offset the contents from the left for value labels or string values, and from the right for numeric values. `OFFSET (CENTER)` centers contents within the column. If `AUTOMATIC` is in effect, the default is `CENTER`. If `MANUAL` is in effect, the

default is 0: value labels and string values are left-justified and numeric values are right-justified. If ONEBREAKCOLUMN is ON, the offset is applied along with the indentation specified on INDENT, always from the left. The specification of CENTER on OFFSET is ignored.

- (UNDERSCORE)** *Use underscores below case listings.* Case listing columns produced by FORMAT LIST are underscored before summary statistics are displayed. You can optionally specify the names of one or more report variables after UNDERSCORE; only the specified columns are underscored.
- (TOTAL)** *Display the summary statistics requested on the next SUMMARY subcommand for all the cases in the report.* TOTAL must be specified on the first BREAK subcommand and applies only to the next SUMMARY subcommand specified.
- (NOTOTAL)** *Display summary statistics only for each break.* This is the default.
- (SKIP(n))** *Skip n lines after the last summary line for a break before beginning the next break.* The default for n is 1.
- (PAGE)** *Begin each break on a new page.* If RESET is specified on PAGE, the page counter resets to the PAGE1 setting on the FORMAT subcommand every time the break value changes for the specified variable. PAGE cannot be specified for listing reports with no break levels.
- (NAME)** *Display the name of the break variable next to each value or value label of the break variable.* NAME requires enough space for the length of the variable name plus two additional characters (for a colon and a blank space) in addition to the space needed to display break values or value labels. NAME is ignored if the break column width is insufficient.
- (NONAME)** *Suppress the display of break variable names.* This is the default.

### Example

```

SORT DIVISION BRANCH DEPT.
REPORT FORMAT=AUTOMATIC MARGINS (1,70) BRKSPACE(-1)

/VARIABLES=SPACE(DUMMY) ' ' (4)
          SALES 'Annual' 'Sales' '1981' (15) (OFFSET(2))
          EXPENSES 'Annual' 'Expenses' '1981' (15) (OFFSET(2))

/BREAK=DIVISION
          BRANCH (10) (TOTAL) (OFFSET(1))
/SUMMARY=MEAN

/BREAK=DEPT 'Department' (10)
/SUMMARY=MEAN.

```

- This example creates a report with three break variables. *BRANCH* breaks within values of *DIVISION*, and *DEPT* breaks within values of *BRANCH*.
- FORMAT sets margins to a maximum of 70 columns and requests that summary lines be displayed on the same line as break values. Because LIST is not specified on FORMAT, only summary statistics are displayed.
- VARIABLES defines three report columns, each occupied by a report variable: *SPACE*, *SALES*, and *EXPENSES*.
- The variable *SPACE* is a dummy variable that exists only within REPORT. It has a null heading and a width of 4. It is used as a space holder to separate the break columns from the report columns.
- *SALES* has a three-line heading and a width of 15. The values of *SALES* are offset two spaces from the right.

- *EXPENSES* is the third report variable and has the same width and offset specifications as *SALES*.
- The leftmost column in the report is reserved for the first two break variables, *DIVISION* and *BRANCH*. Value labels are displayed, since this is the default for *AUTOMATIC*. The break column has a width of 10 and the value labels are offset one space from the left. Value labels more than nine characters long are wrapped. The default column heading is used. *TOTAL* requests a summary line at the end of the report showing the mean of all cases in the report.
- The first *SUMMARY* subcommand displays the mean of each report variable in its report column. This line is displayed each time the value of *DIVISION* or *BRANCH* changes.
- The third break variable, *DEPT*, occupies the second column from the left in the report. The break column has a width of 10 and has a one-line heading. Value labels are displayed in the break column, and those exceeding 10 characters are wrapped.
- The second *SUMMARY* subcommand displays the mean for each report variable when the value of *DEPT* changes.

## ***SUMMARY subcommand***

*SUMMARY* calculates a wide range of aggregate and composite statistics.

- *SUMMARY* must be specified if *LIST* is not specified on *FORMAT*.
- The minimum specification is an aggregate or a composite function and its arguments. This must be the first specification on *SUMMARY*.
- Each *SUMMARY* subcommand following a *BREAK* subcommand specifies a new summary line.
- The default location of the summary title is the column of the break variable to which the summary applies. When more than one function is named on *SUMMARY*, the default summary title is that of the function named first. Both the title and its default column location can be altered. [For more information, see the topic Summary titles on p. 1769.](#)
- The default format can be altered for any function. ([For more information, see the topic Summary print formats on p. 1770.](#))
- *SUMMARY* subcommands apply only to the preceding *BREAK* subcommand. If there is no *SUMMARY* subcommand after a *BREAK* subcommand, no statistics are displayed for that break level.
- To use the summary specifications from a previous *BREAK* subcommand for the current *BREAK* subcommand, specify keyword *PREVIOUS* on *SUMMARY*. [For more information, see the topic Other summary keywords on p. 1771.](#)
- Summary statistics are displayed in report columns. With aggregate functions, you can compute summary statistics for all report variables or for a subset. [For more information, see the topic Aggregate functions on p. 1766.](#) With composite functions, you can compute summaries for all or a subset of report variables and you have additional control over the placement of summary statistics in particular report columns. [For more information, see the topic Composite functions on p. 1768.](#)
- Multiple summary statistics requested on one *SUMMARY* subcommand are all displayed on the same line. More than one function can be specified on *SUMMARY* as long as you do not attempt to place two results in the same report column (*REPORT* will not be executed if you

do). To place results of more than one function in the same report column, use multiple SUMMARY subcommands.

- Any composite and aggregate functions except FREQUENCY and PERCENT can be specified on the same summary line.
- To insert blank lines between summaries when more than one summary line is requested for a break, use keyword SKIP followed by the number of lines to skip in parentheses. The default is 0. [For more information, see the topic Other summary keywords on p. 1771.](#)

## Aggregate functions

Use the aggregate functions to request descriptive statistics for report variables.

- If no variable names are specified as arguments to an aggregate function, the statistic is calculated for all variables named on VARIABLES (all report variables).
- To request an aggregate function for a subset of report variables, specify the variables in parentheses after the function keyword.
- All variables specified for an aggregate function must have been named on VARIABLES.
- Keyword TO cannot be used to specify a list of variables for an aggregate function.
- The result of an aggregate function is always displayed in the report column reserved for the variable for which the function was calculated.
- To use several aggregate functions for the same report variable, specify multiple SUMMARY subcommands. The results are displayed on different summary lines.
- The aggregate functions FREQUENCY and PERCENT have special display formats and cannot be placed on the same summary line with other aggregate or composite functions. They can be specified only once per SUMMARY subcommand.
- Aggregate functions use only cases with valid values.

<b>VALIDN</b>	<i>Valid number of cases.</i> This is the only function available for string variables.
<b>SUM</b>	<i>Sum of values.</i>
<b>MIN</b>	<i>Minimum value.</i>
<b>MAX</b>	<i>Maximum value.</i>
<b>MEAN</b>	<i>Mean.</i>
<b>STDDEV</b>	<i>Standard deviation.</i> Aliases are SD and STDEV.
<b>VARIANCE</b>	<i>Variance.</i>
<b>KURTOSIS</b>	<i>Kurtosis.</i>
<b>SKEWNESS</b>	<i>Skewness.</i>
<b>MEDIAN(min,max)</b>	<i>Median value for values within the range.</i> MEDIAN sets up integer-valued bins for counting all values in the specified range. Noninteger values are truncated when the median is calculated.
<b>MODE(min,max)</b>	<i>Modal value for values within the range.</i> MODE sets up integer-valued bins for counting all values in the specified range. Noninteger values are truncated when the mode is calculated.
<b>PGT(n)</b>	<i>Percentage of cases with values greater than n.</i> Alias PCGT.
<b>PLT(n)</b>	<i>Percentage of cases with values less than n.</i> Alias PCLT.

<b>PIN(min,max)</b>	<i>Percentage of cases within the inclusive value range specified. Alias PCIN.</i>
<b>FREQUENCY(min,max)</b>	<i>Frequency counts for values within the inclusive range. FREQUENCY sets up integer-valued bins for counting all values in the specified range. Noninteger values are truncated when the frequency is computed. FREQUENCY cannot be mixed with other aggregate statistics on a summary line.</i>
<b>PERCENT(min,max)</b>	<i>Percentages for values within the inclusive range. PERCENT sets up integer-valued bins for counting all values in the specified range. Noninteger values are truncated when the percentages are computed. PERCENT cannot be mixed with other aggregate statistics on a summary line.</i>

### Example

```

SORT CASES BY BVAR AVAR.
REPORT FORMAT=AUTOMATIC LIST /VARIABLES=XVAR YVAR ZVAR

  /BREAK=BVAR
  /SUMMARY=SUM
  /SUMMARY=MEAN (XVAR YVAR ZVAR)
  /SUMMARY=VALIDN (XVAR)

  /BREAK=AVAR
  /SUMMARY=PREVIOUS .

```

- `FORMAT` requests a case listing, and `VARIABLES` establishes a report column for variables `XVAR`, `YVAR`, and `ZVAR`. The report columns have default widths and titles.
- Both break variables, `BVAR` and `AVAR`, have default widths and headings.
- Every time the value of `BVAR` changes, three summary lines are displayed. The first line contains the sums for variables `XVAR`, `YVAR`, and `ZVAR`. The second line contains the means of all three variables. The third line displays the number of valid cases for `XVAR` in the report column for `XVAR`.
- Every time the value of `AVAR` changes within each value of `BVAR`, the three summary lines requested for `BVAR` are displayed. These summary lines are based on cases with the current values of `BVAR` and `AVAR`.

### Example

```

SORT CASES BY DEPT.
REPORT FORMAT=AUTOMATIC
  /VARIABLES=WAGE BONUS TENURE
  /BREAK=DEPT (23)
  /SUMMARY=SUM(WAGE BONUS) MEAN(TENURE) 'Sum Income: Mean Tenure'.

```

- `SUMMARY` defines a summary line consisting of the sums of `WAGE` and `BONUS` and the mean of `TENURE`. The result of each aggregate function is displayed in the report column of the variable for which the function is calculated.
- A title is assigned to the summary line. A width of 23 is defined for the break column to accommodate the title for the summary line.

## Composite functions

Use composite functions to obtain statistics based on aggregated statistics, to place a summary statistic in a column other than that of the report variable for which it was calculated, or to manipulate variables not named on VARIABLES.

- Composite functions can be computed for the following aggregate functions: VALIDN, SUM, MIN, MAX, MEAN, STDEV, VARIANCE, KURTOSIS, SKEWNESS, PGT, PLT, and PIN. Constants can also be arguments to composite functions.
- When used within composite functions, aggregate functions can have only one variable as an argument.
- A composite function and its arguments cannot be separated by other SUMMARY specifications.
- The result of a composite function can be placed in any report column, including columns of dummy or string variables, by specifying a target column. To specify a target column, enclose the variable name of the column in parentheses after the composite function and its arguments. By default, the results of a composite function are placed in the report column of the first variable specified on the composite function that is also specified on VARIABLES.
- The format for the result of a composite function can be specified in parentheses after the name of the column location, within the parentheses that enclose the column-location specification.

<b>DIVIDE</b> (arg arg	<i>Divide the first argument by the second and then multiply the result by the factor if it is specified.</i>
<b>MULTIPLY</b> (arg	<i>Multiply the arguments.</i>
<b>PCT</b> (arg	<i>The percentage of the first argument over the second.</i>
<b>SUBTRACT</b> (arg	<i>Subtract the second argument from the first.</i>
<b>ADD</b> (arg	<i>Add the arguments.</i>
<b>GREAT</b> (arg	<i>The maximum of the arguments.</i>
<b>LEAST</b> (arg	<i>The minimum of the arguments.</i>
<b>AVERAGE</b> (arg	<i>The average of the arguments.</i>

### Example

```

SORT CASES BY DEPT.
REPORT FORMAT=AUTOMATIC BRKSPACE(-1)
  /VARIABLES=WAGE BONUS SPACE1 (DUMMY) ' ' BNFT1 BNFT2 SPACE2 (DUMMY) ' '
  /BREAK=DEPT

  /SUMMARY=MEAN(WAGE BONUS BNFT1 BNFT2)
    ADD(VALIDN(WAGE)) (SPACE2)

  /SUMMARY=ADD(SUM(WAGE) SUM(BONUS))
    ADD(SUM(BNFT1) SUM(BNFT2)) 'Totals' SKIP(1)

  /SUMMARY=DIVIDE(MEAN(WAGE) MEAN(BONUS)) (SPACE1 (COMMA)(2))
    DIVIDE(MEAN(BNFT1) MEAN(BNFT2)) (SPACE2 (COMMA)(2)) 'Ratios'
    SKIP(1) .

```

- VARIABLES defines six report columns. The columns for WAGE, BONUS, BNFT1, and BNFT2 contain aggregate statistics based on those variables. The variables SPACE1 and SPACE2 are dummy variables that are created for use as space holders; each is given a blank heading to suppress the default column heading.

- The first SUMMARY computes the means of the variables *WAGE*, *BONUS*, *BNFT1*, and *BNFT2*. Because *BRKSPACE=-1*, this summary line will be placed on the same line as the break value and will have no summary title. The means are displayed in the report column for each variable. SUMMARY also computes the valid number of cases for *WAGE* and places the result in the *SPACE2* column.
- The second SUMMARY adds the sum of *WAGE* to the sum of *BONUS*. Since no location is specified, the result is displayed in the *WAGE* column. In addition, the sum of *BNFT1* is added to the sum of *BNFT2* and the result is placed in the *BNFT1* column. The title for the summary line is *Totals*. One line is skipped before the summary line requested by this SUMMARY subcommand is displayed.
- The third summary line divides the mean of *WAGE* by the mean of *BONUS* and places the result in *SPACE1*. The ratio of the mean of *BNFT1* to the mean of *BNFT2* is displayed in the *SPACE2* column. The results are displayed with commas and two decimal places. The title for the summary line is *Ratios*. One line is skipped before the summary line requested by this SUMMARY subcommand is displayed.

### Summary titles

- You can specify a summary title enclosed in quotes, following the conventions for strings.
- The summary title must be specified after the first function and its arguments. It cannot separate any function from its arguments.
- A summary title can be only one line long.
- A summary title wider than the break column extends into the next break column to the right. If the title is wider than all of the available break columns, it is truncated.
- Only one summary title can be specified per summary line. If more than one is specified, the last is used.
- The summary title is left- or right-justified depending upon whether the break title is left- or right-justified.
- The default location for the summary title is the column of the *BREAK* variable to which the summary applies. With multiple breaks, you can override the default placement of the title by specifying, in parentheses following the title, the number of the break column in which you want the summary title to be displayed.
- In a report with no break levels, REPORT displays the summary title above the summary line at the left margin.

Table 207-2  
Default title for summary lines

Function	Title
VALIDN	N
VARIANCE	Variance
SUM	Sum
MEAN	Mean
STDDEV	StdDev
MIN	Minimum
MAX	Maximum

<b>Function</b>	<b>Title</b>
SKEWNESS	Skewness
KURTOSIS	Kurtosis
PGT(n)	>n
PLT(n)	<n
PIN(min,max)	In $n_1$ to $n_2$
FREQUENCY(min,max)	Total
PERCENT(min,max)	Total
MEDIAN(min,max)	Median
MODE(min,max)	Mode

### **Summary print formats**

All functions have default formats that are used to display results. You can override these defaults by specifying a format keyword and/or the number of decimal places.

- Any printable formats or the `PLAIN` keyword can be specified. Format specifications must be enclosed in parentheses.
- For aggregate functions, the format and/or number of decimal places is specified after the variable name, within the parentheses that enclose the variable name. The variable must be explicitly named as an argument.
- For composite functions, the format and/or number of decimal places is specified after the variable name of the column location, within the parentheses that enclose the variable name. The column location must be explicitly specified.
- If the report column is wide enough, `SUM`, `MEAN`, `STDDEV`, `MIN`, `MAX`, `MEDIAN`, `MODE`, and `VARIANCE` use `DOLLAR` or `COMMA` format if a `DOLLAR` or `COMMA` format has been declared for the variable on either the `FORMATS` or `PRINT FORMATS` command.
- If the column is not wide enough to display the decimal digits for a given function, `REPORT` displays fewer decimal places. If the column is not wide enough to display the integer portion of the number, `REPORT` uses scientific notation if possible, or, if not, displays asterisks.
- An exact value of 0 is displayed with one 0 to the left of the decimal point and as many 0 digits to the right as specified by the format. A number less than 1 in absolute value is displayed without a 0 to the left of the decimal point, except with `DOLLAR` and `COMMA` formats.

**(PLAIN)** *Uses the setting on SET DECIMAL for the thousands separator and decimal delimiter. PLAIN overrides dictionary formats. This is the default for all functions except SUM, MEAN, STDDEV, MIN, MAX, MEDIAN, MODE, and VARIANCE. For these functions, the default is the dictionary format of the variable for which the function is computed.*

**(d)** *Number of decimal places.*

### **Example**

```
/SUMMARY=MEAN(INCOME (DOLLAR) (2))
ADD(SUM(INCOME) SUM(WEALTH)) (WEALTH (DOLLAR) (2))
```



- SUMMARY displays the mean of *INCOME* with dollar format and two decimal places. The result is displayed in the *INCOME* column.
- The sums of *INCOME* and *WEALTH* are added, and the result is displayed in the *WEALTH* column with dollar format and two decimal places.

Table 207-3

*Default print formats for functions*

<b>Function</b>	<b>Format type</b>	<b>Width</b>	<b>Decimal places</b>
VALIDN	F	5	0
SUM	Dictionary	Dictionary + 2	Dictionary
MEAN	Dictionary	Dictionary	Dictionary
STDDEV	Dictionary	Dictionary	Dictionary
VARIANCE	Dictionary	Dictionary	Dictionary
MIN	Dictionary	Dictionary	Dictionary
MAX	Dictionary	Dictionary	Dictionary
SKEWNESS	F	5	2
KURTOSIS	F	5	2
PGT	PCT	6	1
PLT	PCT	6	1
PIN	PCT	6	1
MEDIAN	Dictionary	Dictionary	Dictionary
MODE	Dictionary	Dictionary	Dictionary
PERCENT	F	6	1
FREQUENCY	F	5	0
DIVIDE	F	Dictionary	0
PCT	PCT	6	2
SUBTRACT	F	Dictionary	0
ADD	F	Dictionary	0
GREAT	F	Dictionary	0
LEAST	F	Dictionary	0
AVERAGE	F	Dictionary	0
MULTIPLY	F	Dictionary	0

Where DATE formats are specified, functions with the dictionary format type display the DATE formats, using the column width as the display width.

### ***Other summary keywords***

The following additional keywords can be specified on SUMMARY. These keywords are not enclosed in parentheses.

- SKIP(n)** *Blank lines before the summary line.* The default is 0. If SKIP is specified for the first SUMMARY subcommand for a BREAK, it skips the specified lines after skipping the number of lines specified for BRKSPACE on FORMAT. Similarly, with case listings SKIP skips *n* lines after the blank line at the end of the listing.
- PREVIOUS(n)** *Use the SUMMARY subcommands specified for the nth BREAK.* If *n* is not specified, PREVIOUS refers to the set of SUMMARY subcommands for the previous BREAK. If an integer is specified, the SUMMARY subcommands from the

*n*th BREAK are used. If PREVIOUS is specified, no other specification can be used on that SUMMARY subcommand.

## **TITLE and FOOTNOTE subcommands**

TITLE and FOOTNOTE provide titles and footnotes for the report.

- TITLE and FOOTNOTE are optional and can be placed anywhere after FORMAT except between the BREAK and SUMMARY subcommands.
- The specification on TITLE or FOOTNOTE is the title or footnote in quotes. To specify a multiple-line title or footnote, enclose each line in quotes and separate the specifications for each line by at least one blank.
- The default REPORT title is the title specified on the TITLE command. If there is no TITLE command specified in your session, the default REPORT title is the first line of the header.
- Titles begin on the first line of the report page. Footnotes end on the last line of the report page.
- Titles and footnotes are repeated on each page of a multiple-page report.
- The positional keywords LEFT, CENTER, and RIGHT can each be specified once. The default is CENTER.
- If the total width needed for the combined titles or footnotes for a line exceeds the page width, REPORT generates an error message.

<b>LEFT</b>	<i>Left-justify titles or footnotes within the report page margins.</i>
<b>RIGHT</b>	<i>Right-justify titles or footnotes within the report page margins.</i>
<b>CENTER</b>	<i>Center titles and footnotes within the report page width.</i>

The following can be specified as part of the title or footnote.

<b>)PAGE</b>	<i>Display the page number right-justified in a five-character field.</i>
<b>)DATE</b>	<i>Display the current date in the form dd/mmm/yy, right-justified in a nine-character field.</i>
<b>)var</b>	<i>Display this variable's value label at this position.</i> If you specify a variable that has no value label, the value is displayed, formatted according to its print format. You cannot specify a scratch or system variable or a variable created with the STRING subcommand. If you want to use a variable named DATE or PAGE in the file, change the variable's name with the RENAME VARIABLES command before you use it on the TITLE or FOOTNOTE subcommands, to avoid confusion with the )PAGE and )DATE keywords.

- )PAGE, )DATE, and )var are specified within quotes and can be mixed with string segments within the quotes.
- A variable specified on TITLE or FOOTNOTE must be defined in the active dataset, but does not need to be included as a column on the report.
- One label or value from each variable specified on TITLE or FOOTNOTE is displayed on every page of the report. If a new page starts with a case listing, REPORT takes the value label from the first case listed. If a new page starts with a BREAK line, REPORT takes the value label from the first case of the new break group. If a new page starts with a summary line, REPORT takes the value label from the last case of the break group being summarized.

**Example**

```
/TITLE=LEFT 'Personnel Report' 'Prepared on )DATE'
      RIGHT 'Page: )PAGE'
```

- `TITLE` specifies two lines for a left-justified title and one line for a right-justified title. These titles are displayed at the top of each page of the report.
- The second line of the left-justified title contains the date on which the report was processed.
- The right-justified title displays the page number following the string *Page:* on the same line as the first line of the left-justified title.

**MISSING subcommand**

`MISSING` controls the treatment of cases with missing values.

- `MISSING` specifications apply to variables named on `VARIABLES` and `SUMMARY` and to strings created with the `STRING` subcommand.
- Missing-value specifications are ignored for variables named on `BREAK` when `MISSING=VAR` or `NONE`. There is one break category for system-missing values and one for each user-missing value.
- The character used to indicate missing values is controlled by the `FORMAT` subcommand.

**VAR**

*Missing values are treated separately for each variable.* Missing values are displayed in case listings but are not included in the calculation of summary statistics on a function-by-function basis. This is the default.

**NONE**

*User-missing values are treated as valid values.* This applies to all variables named on `VARIABLES`.

**LIST([varlist][n])**

*Cases with the specified number of missing values across the specified list of variables are not used.* The variable list and *n* are specified in parentheses. If *n* is not specified, the default is 1. If no variables are specified, all variables named on `VARIABLES` are assumed.

**Example**

```
/MISSING= LIST (XVAR, YVAR, ZVAR 2)
```

- Any case with two or more missing values across the variables *XVAR*, *YVAR*, and *ZVAR* is omitted from the report.

# REPOSITORY ATTRIBUTES

```
REPOSITORY ATTRIBUTES FILE='file specification'  
  [DESCRIPTION='description']  
  [KEYWORDS='keywords']  
  [AUTHOR='author']  
  [VERSIONLABEL='label']  
  [EXPIRATION=days]  
  [TOPICS='topics']  
  [/SECURITY ID='id' [PERMISSION= [READ**] [WRITE] [DELETE] [MODIFY] [OWNER] ] ]
```

\*\* Default if the keyword is omitted.

## **Release History**

Release 16.0

- Command introduced as PER ATTRIBUTES.

Release 17.0

- VERSIONLABEL keyword extended to support multiple labels.

Release 18.0

- PER ATTRIBUTES deprecated. Command name changed to REPOSITORY ATTRIBUTES.

## **Example**

```
REPOSITORY COPY  
  FILE='/myscripts/cust_value.py'  
  OUTFILE='SPSSCR://scripts/cust_value.py'.  
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
  DESCRIPTION='Customer Value Calculation'  
  KEYWORDS='customer;value'.
```

## **Overview**

The REPOSITORY ATTRIBUTES command allows you to set attributes—such as a version label and security settings—for an object in a IBM® SPSS® Collaboration and Deployment Services Repository.

- When copying to a IBM SPSS Collaboration and Deployment Services Repository with the REPOSITORY COPY command, use the REPOSITORY ATTRIBUTES command (following REPOSITORY COPY) to specify attributes of the object.

## **Basic Specification**

The basic specification is the FILE keyword, which specifies the IBM SPSS Collaboration and Deployment Services Repository object whose attributes are to be set. All other keywords and subcommands are optional.

**Syntax Rules**

- The `SECURITY` subcommand can be specified multiple times.
- Each keyword can only be specified once.
- Keywords and subcommands can be used in any order.
- Keywords and subcommand names must be spelled in full.
- Equals signs (=) shown in the syntax chart are required.

**Operations**

- Use of the `REPOSITORY ATTRIBUTES` command requires a connection to a IBM SPSS Collaboration and Deployment Services Repository. Connections are established with the `REPOSITORY CONNECT` command.
- `REPOSITORY ATTRIBUTES` overwrites any existing values of specified attributes.

**FILE Keyword**

The `FILE` keyword is required and specifies the IBM® SPSS® Collaboration and Deployment Services Repository object whose attributes are to be set.

- The form of the file specification for an object in a IBM SPSS Collaboration and Deployment Services Repository is the scheme name `SPSSCR`, followed by a colon, either one or two slashes (forward or backward), and a file path, all enclosed in quotes. For example:

```
'SPSSCR://scripts/myscript.py'
```

- Paths can be specified with forward slashes (/) or backslashes (\).
- You can define a file handle to a file or a directory in a IBM SPSS Collaboration and Deployment Services Repository and use that handle in file specifications for IBM SPSS Collaboration and Deployment Services Repository objects.
- You can use the `CD` command to set the working directory to a directory in the currently connected IBM SPSS Collaboration and Deployment Services Repository, allowing you to use relative paths in file specifications for IBM SPSS Collaboration and Deployment Services Repository objects.
- File specifications for IBM SPSS Collaboration and Deployment Services Repository objects must specify the filename exactly as provided when the file was stored. If the file was stored with an extension, then you must supply the extension. If the file was stored without an extension then do not include one.

For examples of file specifications, see the examples for the `REPOSITORY COPY` command on p. 1783.

**DESCRIPTION Keyword**

The `DESCRIPTION` keyword specifies a description for an object in a IBM® SPSS® Collaboration and Deployment Services Repository and replaces any existing description for the object. Specify the value as a quoted string.

**Example**

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
DESCRIPTION='Customer Value Calculation'.
```

**KEYWORDS Keyword**

KEYWORDS specifies one or more keywords to associate with an object in a IBM® SPSS® Collaboration and Deployment Services Repository to aid in searching. Specify the value as a quoted string.

- Multiple keywords should be separated by semicolons.
- Blank spaces at the beginning and end of each keyword are ignored, but blank spaces within keywords are honored.
- The specified keywords replace any existing ones for the object.

**Example**

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
KEYWORDS='customer;value'.
```

**AUTHOR Keyword**

The AUTHOR keyword specifies the author of an object in a IBM® SPSS® Collaboration and Deployment Services Repository. By default, the author is set to the login name of the user who created the object. Specify the value as a quoted string.

**Example**

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
AUTHOR='GSWEET'.
```

**VERSIONLABEL Keyword**

The VERSIONLABEL keyword specifies one or more version labels for an object in a IBM® SPSS® Collaboration and Deployment Services Repository—for example, “production” or “development”. Two versions of an object cannot have any of the same labels. If you specify a label associated with another version, it will be removed from that version and associated with the one you’re modifying. For a single label, specify the value as a quoted string. For multiple labels, specify the value as a blank-separated list of quoted strings—for example, 'label1' 'label2'.

- By default, the specified labels will be applied to the latest version of the object that is accessible to the current user. You can apply labels to a version other than the latest one by specifying the version in the file specification on the FILE keyword. [For more information, see the topic File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects on p. 2230.](#)

**Example**

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
  VERSIONLABEL='development'.
```

**EXPIRATION Keyword**

The `EXPIRATION` keyword specifies an expiration date for an object in a IBM® SPSS® Collaboration and Deployment Services Repository. This provides a mechanism to make sure dated information is not displayed after a certain date. Expired documents are not deleted but are automatically removed to a special category where they can be accessed only by the site administrator, who can delete, archive, or republish them.

- Specify the value as an integer representing the number of days from the current day (inclusive) to the last day (inclusive) that the document will be active.

**Example**

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
  EXPIRATION = 366.
```

**TOPICS Keyword**

The `TOPICS` keyword allows you to associate an object in a IBM® SPSS® Collaboration and Deployment Services Repository with one or more topics. Topics allow you to organize documents by subject matter and have a hierarchical structure.

- Topics are specified as a quoted path that includes each level of the hierarchy for that topic, with successive levels separated by a forward slash. A forward slash at the beginning of the path is optional.
- Use a semicolon as the delimiter when specifying multiple topics.
- Objects can only be associated with existing topics. `REPOSITORY ATTRIBUTES` cannot be used to create new topics.
- The specified topics replace any existing ones for the object.

**Example**

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
  TOPICS = '/engineering/scripts;/marketing/analyses'.
```

**SECURITY Subcommand**

The `SECURITY` subcommand allows you to specify security settings for an object in a IBM® SPSS® Collaboration and Deployment Services Repository. Specify the identifier of the user or group with the `ID` keyword, and specify the access level with the `PERMISSION` keyword. You can specify settings for multiple users or groups by including multiple instances of the `SECURITY` subcommand.

**ID Keyword**

The ID keyword specifies the ID of the user or group for which access is being granted. Specify the value in quotes.

**PERMISSION Keyword**

The PERMISSION keyword is optional and specifies the access level granted to the specified user or group. One or more of the following access levels can be specified: READ, WRITE, MODIFY (grants ability to modify permissions), DELETE, and OWNER.

- If PERMISSION is omitted, the access level is READ.
- READ access is always granted, whether or not it is specified on the PERMISSION keyword.
- If OWNER is specified, all other values are ignored (an owner has all permissions).

**Example**

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'  
  /SECURITY ID='admin' PERMISSION=OWNER  
  /SECURITY ID='--everyone--' PERMISSION=READ WRITE.
```



# REPOSITORY CONNECT

```
REPOSITORY CONNECT
  /SERVER HOST='host[:{8080**}]' [SSL={NO**}]
           {port }           {YES }
  /LOGIN  USER='userid'
          PASSWORD='password'
          [PROVIDER='security provider']
          [ENCRYPTEDPWD={YES**}]
           {NO }
```

\*\* Default if the keyword or value is omitted.

## **Release History**

Release 15.0

- Command introduced as PER CONNECT.

Release 17.0

- DOMAIN keyword deprecated on the LOGIN subcommand.
- PROVIDER keyword introduced on the LOGIN subcommand.

Release 18.0

- PER CONNECT deprecated. Command name changed to REPOSITORY CONNECT.

## **Example**

```
REPOSITORY CONNECT
  /SERVER HOST='PER1'
  /LOGIN  USER='MyUserID'
          PASSWORD='abc12345'
          ENCRYPTEDPWD=NO.
```

## **Overview**

The REPOSITORY CONNECT command establishes a connection to a IBM® SPSS® Collaboration and Deployment Services Repository and logs in the user. A connection enables you to store objects to, and retrieve objects from, a IBM SPSS Collaboration and Deployment Services Repository.

### **Options**

**Server.** You can specify a connection port and whether to connect to the specified server using Secure Socket Layer (SSL) technology, if it is enabled on the server.

**Login.** You can specify whether the password is provided as encrypted or unencrypted (plain text).

**Single Sign-On.** Single sign-on allows users to connect to a IBM SPSS Collaboration and Deployment Services Repository without explicitly providing a user ID and password. The necessary authentication is performed with the user's existing credentials on the current machine—obtained, for example, from Windows Active Directory. If your IBM SPSS Collaboration and Deployment Services Repository supports single sign-on (contact your local system administrator for details), you can omit the `LOGIN` subcommand.

### **Basic Specification**

The basic specification for `REPOSITORY CONNECT` is the host server, user name, and password. By default, server port 8080 is used, the connection is established without SSL, and the specified password is assumed to be encrypted. To create an encrypted password, generate (paste) the `REPOSITORY CONNECT` command syntax from the IBM® SPSS® Collaboration and Deployment Services Repository Connect dialog box.

### **Syntax Rules**

- Each subcommand can be specified only once.
- Subcommands can be used in any order.
- An error occurs if a keyword or attribute is specified more than once within a subcommand.
- Equals signs (=) and forward slashes (/) shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.

### **Operations**

- `REPOSITORY CONNECT` establishes a connection to an IBM SPSS Collaboration and Deployment Services Repository and logs in the specified user. Any existing connection terminates when the new one is established.
- You can specify whether a connection is established when the `REPOSITORY CONNECT` command is encountered or whether the connection is deferred until it is required. The behavior is determined by the `REPDEFER` subcommand of the `SET` command.
- The connection terminates if the IBM® SPSS® Statistics session ends.
- An error occurs if a connection cannot be established to the specified host server.
- An error occurs if the connection cannot be authenticated—for example, if the password is invalid for the specified user.

### **Example**

```
REPOSITORY CONNECT
  /SERVER HOST='PER1:80'
  /LOGIN  USER='MyUserID'
          PASSWORD='abc12345'
          ENCRYPTEDPWD=NO.
```

- The `SERVER` subcommand specifies a connection to host 'PER1' on port 80.
- `ENCRYPTEDPWD=NO` indicates that the password is not encrypted.

## ***SERVER Subcommand***

The `SERVER` subcommand specifies the host server and whether to establish a secure connection.

<b>HOST</b>	<i>Server that hosts the IBM® SPSS® Collaboration and Deployment Services Repository.</i> Specify the name of the server in quotes. The default port is 8080. To connect to another port, specify the port number after the host name; for example, 'PER1:80'. A colon must separate the host name and port.
<b>SSL</b>	<i>Use Secure Socket Layer technology.</i> Specifies whether to establish a secure connection to the host server. The default is <code>NO</code> . SSL is available only if supported on the host server.

## ***LOGIN Subcommand***

The `LOGIN` subcommand specifies login information, including user name and password.

<b>USER</b>	<i>User name.</i> Specify the user name in quotes.
<b>PASSWORD</b>	<i>Password.</i> Specify the password in quotes.
<b>PROVIDER</b>	<i>Security provider.</i> You can optionally specify the security provider against which to validate the user/password combination. Contact your local administrator for details.
<b>ENCRYPTED-PWD</b>	<i>Password encryption.</i> By default, the specified password is treated as encrypted. To indicate that the password is entered as plain text, specify <code>ENCRYPTEDPWD=NO</code> .

# REPOSITORY COPY

```
REPOSITORY COPY
  FILE='file specification'
  OUTFILE='file specification'.
```

## **Release History**

Release 16.0

- Command introduced as `PER COPY`.

Release 18.0

- `PER COPY` deprecated. Command name changed to `REPOSITORY COPY`.

## **Example**

```
REPOSITORY COPY
  FILE='/myscripts/demo.py'
  OUTFILE='SPSSCR://scripts/demo.py'.
```

## **Overview**

The `REPOSITORY COPY` command allows you to copy an arbitrary file from the local file system to a IBM® SPSS® Collaboration and Deployment Services Repository or to copy a file from a IBM SPSS Collaboration and Deployment Services Repository to the local file system.

- When copying to a IBM SPSS Collaboration and Deployment Services Repository, use the `REPOSITORY ATTRIBUTES` command—following `REPOSITORY COPY`—to specify properties such as a description, keywords, and security settings for the object.

## **Basic Specification**

The only specification is the `FILE` keyword, which specifies the source file, and the `OUTFILE` keyword, which specifies the target location. Each keyword can specify a location in the local file system or a location in the current IBM SPSS Collaboration and Deployment Services Repository.

## **File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects**

- The form of the file specification for an object in a IBM SPSS Collaboration and Deployment Services Repository is the scheme name `SPSSCR`, followed by a colon, either one or two slashes (forward or backward), and a file path, all enclosed in quotes. For example:

```
'SPSSCR://scripts/myscript.py'
```

- Paths can be specified with forward slashes (/) or backslashes (\).

- You can define a file handle to a file or a directory in a IBM SPSS Collaboration and Deployment Services Repository and use that handle in file specifications for IBM SPSS Collaboration and Deployment Services Repository objects.
- You can use the `CD` command to set the working directory to a directory in the currently connected IBM SPSS Collaboration and Deployment Services Repository, allowing you to use relative paths in file specifications for IBM SPSS Collaboration and Deployment Services Repository objects.
- File extensions are not added to files stored to a IBM SPSS Collaboration and Deployment Services Repository. Files are stored to a IBM SPSS Collaboration and Deployment Services Repository with an automatically determined MIME type that describes the type of file. Including a file extension is not necessary but is recommended.
- When copying a file from a IBM SPSS Collaboration and Deployment Services Repository you must specify the filename exactly as provided when the file was stored. If the file was stored with an extension, then you must supply the extension. If the file was stored without an extension then do not include one.
- When copying from a IBM SPSS Collaboration and Deployment Services Repository, the latest version of the specified object that is accessible to the current user is retrieved. To specify a version by label, use `SPSSCR://<path>#L.<label>`. To specify a version by the version marker (consists of an integer version number followed by a time stamp that uniquely specifies a version), use `SPSSCR://<path>#M.<version marker>`. [For more information, see the topic File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects on p. 2230.](#)

### **Operations**

- `REPOSITORY COPY` will fail if a IBM SPSS Collaboration and Deployment Services Repository location is specified (for either `FILE` or `OUTFILE`) and there is no connection to a IBM SPSS Collaboration and Deployment Services Repository. Connections are established with the `REPOSITORY CONNECT` command.
- Specifying an existing IBM SPSS Collaboration and Deployment Services Repository file on the `OUTFILE` keyword results in a new version of the file, as opposed to overwriting the latest version.
- The `PERMISSIONS`, `SAVE`, and `XSAVE` commands will generate errors if you attempt to set permissions for IBM SPSS Collaboration and Deployment Services Repository objects. Permissions for IBM SPSS Collaboration and Deployment Services Repository objects are set with the `REPOSITORY ATTRIBUTES` command or using administration tools included with IBM® SPSS® Collaboration and Deployment Services.
- The `ERASE FILE` command will generate an error if you attempt to delete a IBM SPSS Collaboration and Deployment Services Repository object.

## **Examples**

### **Copying from the Local File System to a IBM SPSS Collaboration and Deployment Services Repository**

```
REPOSITORY COPY
FILE='/myscripts/cust_value.py'
OUTFILE='SPSSCR://scripts/cust_value.py' .
```

## REPOSITORY COPY

```
REPOSITORY ATTRIBUTES FILE='SPSSCR://scripts/cust_value.py'
KEYWORDS='customer;value'
/SECURITY ID='rprime' PERMISSION=OWNER.
```

- The local file `/myscripts/cust_value.py` is copied to the location `/scripts/cust_value.py` in the current IBM® SPSS® Collaboration and Deployment Services Repository.
- The `REPOSITORY ATTRIBUTES` command specifies keywords and security settings for this IBM SPSS Collaboration and Deployment Services Repository object.

**Copying from a IBM SPSS Collaboration and Deployment Services Repository to the Local File System**

```
REPOSITORY COPY
FILE='SPSSCR://scripts/demo.py'
OUTFILE='/myscripts/demo.py'.
```

- The IBM SPSS Collaboration and Deployment Services Repository object `/scripts/demo.py` is copied to the local file `/myscripts/demo.py`.

**Copying from a Labelled Version in a IBM SPSS Collaboration and Deployment Services Repository to the Local File System**

```
REPOSITORY COPY
FILE='SPSSCR://scripts/demo.py#L.production'
OUTFILE='/myscripts/demo.py'.
```

- The version of the IBM SPSS Collaboration and Deployment Services Repository object `/scripts/demo.py` with the label `production` is copied to the local file `/myscripts/demo.py`.

**Using a File Handle to a IBM SPSS Collaboration and Deployment Services Repository Location**

```
FILE HANDLE custval /NAME='SPSSCR://CustomerValue'.
REPOSITORY COPY
FILE='custval/cust_value.py'
OUTFILE='/myscripts/cust_value.py'.
```

- The handle `custval` is associated with the IBM SPSS Collaboration and Deployment Services Repository directory `/CustomerValue`, and the `REPOSITORY COPY` command retrieves the latest version of the file `cust_value.py` from this directory in the current IBM SPSS Collaboration and Deployment Services Repository.

**Setting the Working Directory to a IBM SPSS Collaboration and Deployment Services Repository Location**

```
CD 'SPSSCR://CustomerValue'.
REPOSITORY COPY
FILE='cust_value.py'
OUTFILE='/myscripts/cust_value.py'.
```

- The working directory is set to the IBM SPSS Collaboration and Deployment Services Repository directory `/CustomerValue`, and the `REPOSITORY COPY` command retrieves the latest version of the file `cust_value.py` from this directory in the current IBM SPSS Collaboration and Deployment Services Repository.

# REREAD

```
REREAD [FILE=file]
       [COLUMN=expression]
```

## **Example**

```
INPUT PROGRAM.
DATA LIST /KIND 10-14 (A).

DO IF (KIND EQ 'FORD').
REREAD.
DATA LIST /PARTNO 1-2 PRICE 3-6 (DOLLAR,2) QUANTITY 7-9.
END CASE.

ELSE IF (KIND EQ 'CHEVY').
REREAD.
DATA LIST /PARTNO 1-2 PRICE 15-18 (DOLLAR,2) QUANTITY 19-21.
END CASE.
END IF.

END INPUT PROGRAM.

BEGIN DATA
111295100FORD
11      CHEVY 295015
END DATA.
```

## **Overview**

REREAD instructs the program to reread a record in the data. It is available only within an INPUT PROGRAM structure and is generally used to define data using information obtained from a previous reading of the record. REREAD is usually specified within a conditional structure, such as DO IF-END IF, and is followed by a DATA LIST command. When it receives control for a case, REREAD places the pointer back to the column specified for the current case and begins reading data as defined by the DATA LIST command that follows.

### **Options**

**Data Source.** You can use inline data or data from an external file specified on the FILE subcommand. Using external files allows you to open multiple files and merge data.

**Beginning Column.** You can specify a beginning column other than column 1 using the COLUMN subcommand.

### **Basic Specification**

The basic specification is the command keyword REREAD. The program rereads the current case according to the data definitions specified on the following DATA LIST.

**Subcommand Order**

Subcommands can be specified in any order.

**Syntax Rules**

- REREAD is available only within an INPUT PROGRAM structure.
- Multiple REREAD commands can be used within the input program. Each must be followed by an associated DATA LIST command.

**Operations**

- REREAD causes the next DATA LIST command to reread the most recently processed record in the specified file.
- When it receives control for a case, REREAD places the pointer back to column 1 for the current case and begins reading data as defined by the DATA LIST that follows. If the COLUMN subcommand is specified, the pointer begins reading in the specified column and uses it as column 1 for data definition.
- REREAD can be used to read part of a record in FIXED format and the remainder in LIST format. Mixing FIXED and FREE formats yields unpredictable results.
- Multiple REREAD commands specified without an intervening DATA LIST do not have a cumulative effect. All but the last are ignored.

**Examples****Using REREAD to Process Different Record Formats**

```

INPUT PROGRAM.
DATA LIST /PARTNO 1-2 KIND 10-14 (A).

DO IF (KIND EQ 'FORD').
REREAD.
DATA LIST /PRICE 3-6 (DOLLAR,2) QUANTITY 7-9.
END CASE.

ELSE IF (KIND EQ 'CHEVY').
REREAD.
DATA LIST /PRICE 15-18 (DOLLAR,2) QUANTITY 19-21.
END CASE.
END IF.
END INPUT PROGRAM.

BEGIN DATA
111295100FORD          CHAPMAN AUTO SALES
121199005VW           MIDWEST VOLKSWAGEN SALES
11 395025FORD          BETTER USED CARS
11      CHEVY 195005      HUFFMAN SALES & SERVICE
11      VW 595020         MIDWEST VOLKSWAGEN SALES
11      CHEVY 295015      SAM'S AUTO REPAIR
11      CHEVY 210 20      LONGFELLOW CHEVROLET
9555032 VW             HYDE PARK IMPORTS
END DATA.
LIST.
```



- Data are extracted from an inventory of automobile parts. The automobile part number always appears in columns 1 and 2, and the automobile type always appears in columns 10 through 14. The location of other information such as price and quantity depends on both the part number and the type of automobile.
- The first `DATA LIST` extracts the part number and type of automobile.
- Depending on the information from the first `DATA LIST`, the records are reread using one of two `DATA LIST` commands, pulling the price and quantity from different places.
- The two `END CASE` commands limit the active dataset to only those cases with part 11 and automobile type Ford or Chevrolet. Without the `END CASE` commands, cases would be created for other part numbers and automobile types, with missing values for price, quantity, and buyer.

The output from the `LIST` command is as follows:

```
PARTNO KIND    PRICE QUANTITY
    11  FORD  $12.95    100
    11  FORD  $3.95     25
    11  CHEVY $1.95     5
    11  CHEVY $2.95    15
    11  CHEVY $2.10    20
```

### **Multiple REREAD Commands for the Same Record**

```
INPUT PROGRAM.
DATA LIST      NOTABLE/ CDIMAGE 1-20(A).
REREAD        COLUMN = 6. /* A, C, and E are in column 6
REREAD        COLUMN = 11. /* B, D, and F are in column 11
DATA LIST     NOTABLE/ INFO 1(A).
END INPUT PROGRAM.
BEGIN DATA
1   A   B
2   C   D
3   E   F
END DATA.
LIST.
```

- Multiple `REREAD` commands are used without an intervening `DATA LIST`. Only the last one is used. Thus, the starting column comes from the last `REREAD` specified and the pointer is reset to column 11.

The output from the `LIST` command is as follows:

```
CDIMAGE          INFO
1   A   B        B
2   C   D        D
3   E   F        F
```

## **FILE Subcommand**

`FILE` specifies an external raw data file from which the next `DATA LIST` command reads data.

- The default file is the file specified on the immediately preceding `DATA LIST` command.

## REREAD

- If the file specified on `FILE` is not the default file, the same file must be specified on the next `DATA LIST`. Otherwise, the `FILE` subcommand is ignored and the `DATA LIST` command reads the next record from the file specified on it or, if no file is specified, from the file specified on the previous `DATA LIST` command.

**Example**

```

INPUT PROGRAM.
DATA LIST FILE=UPDATE END=#EOF NOTABLE
  /#ID 1-3.                /*Get rep ID in new sales file.
DATA LIST FILE = SALESREP NOTABLE
  /ID 1-3 SALES 4-11(F,2)
  NEWSALE 12-19(F,2).     /*Get rep record from master file.

LOOP IF #EOF OR (#ID GT ID). /*If UPDATE ends or no new sales made.
+ COMPUTE NEWSALE = 0.      /*Set NEWSALE to 0
+ END CASE.                /*Build a case.
+ DATA LIST FILE = SALESREP NOTABLE
  /ID 1-3 SALES 4-11(F,2)
  NEWSALE 12-19(F,2).     /*Continue reading masterfile.
END LOOP

DO IF NOT #EOF.           /*If new sales made.
+ REREAD FILE=UPDATE COLUMN = 4. /*Read new sales from UPDATE.
+ DATA LIST FILE=UPDATE
  /NEWSALE 1-8(F,2).
+ COMPUTE SALES=SALES+NEWSALE. /*Update master file.
END IF.
END CASE.                 /*Build a case.
END INPUT PROGRAM.

LIST.

```

- This example uses `REREAD` to merge two raw data files (`SALESREP` and `UPDATE`).
- Both files are sorted by sales representative ID number. The `UPDATE` file contains only records for sales representatives who have made new sales, with variables `ID` and `NEWSALE`. The master file `SALESREP` contains records for all sales representatives, with variables `SALES` (which contains year-to-date sales) and `NEWSALE` (which contains the update values each time the file is updated).
- If a sales representative has made no new sales, there is no matching ID in the `UPDATE` file. When `UPDATE` is exhausted or when the ID's in the two files do not match, the loop structure causes the program to build a case with `NEWSALE` equal to 0 and then continue reading the master file.
- When the ID's match (and the `UPDATE` file is not yet exhausted), the `REREAD` command is executed. The following `DATA LIST` rereads the record in `UPDATE` that matches the `ID` variable. `NEWSALE` is read from the `UPDATE` file starting from column 4 and `SALES` is updated. Note that the following `DATA LIST` specifies the same file.
- When the updated base is built, the program returns to the first `DATA LIST` command in the input program and reads the next ID from the `UPDATE` file. If the `UPDATE` file is exhausted (`#EOF=1`), the loop keeps reading records from the master file until it reaches the end of the file.
- The same task can be accomplished using `MATCH FILES`. With `MATCH FILES`, the raw data must be read and saved as IBM® SPSS® Statistics data files first.

## COLUMN Subcommand

COLUMN specifies a beginning column for the REREAD command to read data. The default is column 1. You can specify a numeric expression for the column.

### Specifying a Numeric Expression for the Column

```
INPUT PROGRAM.
DATA LIST /KIND 10-14 (A).
COMPUTE #COL=1.
IF (KIND EQ 'CHEVY') #COL=13.

DO IF (KIND EQ 'CHEVY' OR KIND EQ 'FORD').
REREAD COLUMN #COL.
DATA LIST /PRICE 3-6 (DOLLAR,2) QUANTITY 7-9.
END CASE.
END IF.
END INPUT PROGRAM.
BEGIN DATA
111295100FORD          CHAPMAN AUTO SALES
121199005VW           MIDWEST VOLKSWAGEN SALES
11 395025FORD         BETTER USED CARS
11      CHEVY 195005   HUFFMAN SALES & SERVICE
11      VW 595020     MIDWEST VOLKSWAGEN SALES
11      CHEVY 295015   SAM'S AUTO REPAIR
12      CHEVY 210 20   LONGFELLOW CHEVROLET
9555032 VW           HYDE PARK IMPORTS
END DATA.
LIST.
```

- The task in this example is to read *PRICE* and *QUANTITY* for Chevrolets and Fords only. A scratch variable is created to indicate the starting column positions for *PRICE* and *QUANTITY*, and a single `DATA LIST` command is used to read data for both types of automobiles.
- Scratch variable `#COL` is set to 13 for Chevrolets and 1 for all other automobiles. For Fords, the data begin in column 1. Variable *PRICE* is read from columns 3–6 and *QUANTITY* is read from columns 7–9. When the record is a Chevrolet, the data begins in column 13. Variable *PRICE* is read from columns 15–18 (15 is 3, 16 is 4, and so forth), and *QUANTITY* is read from columns 19–21.

### Reading Both FIXED and LIST Input With REREAD

```
INPUT PROGRAM.
DATA LIST     NOTABLE FIXED/ A 1-14(A).  /*Read the FIXED portion
REREAD       COLUMN = 15.
DATA LIST     LIST/ X Y Z.              /*Read the LIST portion
END INPUT PROGRAM.
```

\* The value 1 on the first record is in column 15.

```
BEGIN DATA
FIRST RECORD  1 2 3 -1 -2 -3
NUMBER 2     4 5
THE THIRD    6 7 8
#4
FIFTH AND LAST9 10 11
END DATA.
LIST.
```

- Columns 1–14 are read in `FIXED` format. `REREAD` then resets the pointer to column 15. Thus, beginning in column 15, values are read in `LIST` format.
- The second `DATA LIST` specifies only three variables. Thus, the values `-1`, `-2`, and `-3` on the first record are not read.
- The program generates a warning for the missing value on record 2 and a second warning for the three missing values on record 4.
- On the fifth and last record there is no delimiter between value `LAST` and value `9`. `REREAD` can still read the `9` in `LIST` format.

# RESPONSE RATE

RESPONSE RATE is available in the Direct Marketing option.

```
RESPONSE RATE
/VARIABLES
  NUMERATOR=varname
  DENOMINATOR=varname
  GROUP=varname
[/MINRATE {RATE=NONE**|value    }]
          {FORMULA=value1, value1}
[/MAXCOUNT {COUNT=ALL**|value}]
           {PERCENT=value    }
```

\*\*Default if subcommand omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release history

Release 18.0

- Command introduced.

## Example

```
RESPONSE RATE
/VARIABLES
  NUMERATOR=Responses
  DENOMINATOR=Customers
  GROUP=Rank
/MINRATE RATE=5
/MAXCOUNT COUNT=6000.
```

## Overview

RESPONSE RATE calculates group response rates and cumulative group response rates, expressed as a percentage, where response rate is calculated as NUMERATOR divided by DENOMINATOR for each category of GROUP, and cumulative response rate is calculated as the cumulative sum of NUMERATOR divided by the cumulative sum of DENOMINATOR for the current and all preceding categories of GROUP.

- Output includes a table of response rates, sum of the DENOMINATOR variable, cumulative response rates, and cumulative sum of the DENOMINATOR variable in ascending order of the GROUP variable, and a line chart of response rates and cumulative sum of the DENOMINATOR variable.
- If MINRATE or MAXCOUNT are specified, the table is color-coded based on the specified cumulative value(s), and the line charts include reference lines at the specified value(s).

The `RESPONSE RATE` command is intended primarily for use as part of the Postal Code Response Rate dialog interface, and some of the results refer to a new, aggregated dataset that is created by the dialog interface. When you use the `RESPONSE RATE` command outside of the Postal Code Response Rate dialog interface, this new, aggregated dataset is not created.

### **Basic specification**

- The basic specification is the command name `RESPONSE RATE` and a `VARIABLES` subcommand.
- The `MINRATE` and `MAXCOUNT` subcommands are optional.

### **Operations**

- Cases with missing values for any variables specified on the `VARIABLES` subcommand are excluded from the analysis (listwise deletion).
- If the sum of `DENOMINATOR` is 0 or negative for a value of `GROUP`, that group is excluded from the analysis.

### **Subcommand order**

Subcommands can be specified in any order.

### **Syntax rules**

- The `VARIABLES` subcommand is required.
- The `VARIABLES` subcommand must include `NUMERATOR`, `DENOMINATOR` and `GROUP` variables.
- Equals signs (=) shown in the syntax chart and in examples are required.
- Command, subcommand, and keyword names must be spelled out fully. Abbreviation is not allowed.

## **Examples**

### **Minimum Specification**

```
RESPONSE RATE
/VARIABLES
  NUMERATOR=Responses
  DENOMINATOR=Customers
  GROUP=PostalRank.
```

- Response rates are displayed for each group defined by *PostalRank*.
- Results are displayed in ascending order of the value of *PostalRank*.

### **MINRATE and MAXCOUNT**

```
RESPONSE RATE
/VARIABLES
  NUMERATOR=Responses
  DENOMINATOR=Customers
```

```
GROUP=PostalRank
/MINRATE=5
/MAXCOUNT=6000.
```

- Rows with a cumulative response rate above 5% and a cumulative customer count below 6,000 are displayed with a green background in the table; those with cumulative response rates below 5% or a cumulative count above 6,000 are displayed with a red background.
- The upper panel of the chart includes a reference line at the 5% response rate level, and the lower panel includes a reference line at a cumulative count of 6,000 customers.

## VARIABLES subcommand

The **VARIABLES** subcommand specifies the variables to use. This subcommand is required and includes three required keywords:

**NUMERATOR.** *The variable used as the numerator to calculate response rates.* The variable must be numeric.

**DENOMINATOR.** *The variable used as the denominator to calculate response rates.* The variable must be numeric.

**GROUP.** *The grouping variable.* Response rates are calculated for each category of the grouping variable. The variable can be string or numeric.

Each keyword must be followed by an equals sign (=) and a variable name.

Response rate is calculated as the **NUMERATOR** divided by **DENOMINATOR** for each category of **GROUP**. Cumulative response rate is calculated as the cumulative sum of **NUMERATOR** divided by the cumulative sum of **DENOMINATOR**, where cumulative sum is the sum for the current and all preceding (lower) values of **GROUP**. Response rates are expressed as percentages.

### Example

```
RESPONSE RATE
/VARIABLES
  NUMERATOR=Responses
  DENOMINATOR=Customers
  GROUP=PostalRank.
```

## MINRATE subcommand

The optional **MINRATE** subcommand specifies a minimum cumulative response rate to use to color-code the output table and to use as a reference line on the response rate line chart. The following alternatives are available for specifying minimum response rate:

**RATE=NONE.** *No minimum response rate.* This is the default if the subcommand is omitted.

**RATE=value.** *Use the specified percentage value as the minimum cumulative response rate.* Specify a value greater than 0 and less than 100.

**FORMULA=value1,value2.** Calculate minimum cumulative response rate based on the formula  $(value1/value2)*100$ . Both values must be positive numbers. The result should be a value greater than 0 and less than 100. In direct marketing, for example, *value1* might be the cost of mailing a package, and *value2* might be net revenue per response. If the cost of mailing a package is \$0.75 and the net revenue per response is \$56, then the minimum response rate is:  $(0.75/56)*100=1.34\%$

Each keyword must be followed by an equals sign (=) and a variable name.

- If **MINRATE** is specified without **MAXCOUNT**, all rows with a cumulative response rate equal to or greater than the specified minimum are displayed with a green background, and all rows with a cumulative response rate less than the specified minimum are displayed with a red background.
- If **MAXCOUNT** is also specified, only rows with a cumulative response rate above the specified minimum *and* rows with a cumulative count below the specified maximum are displayed with a green background. For example, if the response rate is higher than the specified minimum, but the cumulative count is greater than the specified maximum, then the row will have a red background.

### Example

```
RESPONSE RATE
/VARIABLES
  NUMERATOR=Responses
  DENOMINATOR=Customers
  GROUP=PostalRank
/MINRATE=5.
```

## MAXCOUNT subcommand

The optional **MAXCOUNT** subcommand specifies the maximum cumulative total of the **DENOMINATOR** variable to use to color-code the output table and to use as a reference line on the cumulative count line chart. For each category of the **GROUP** total, the cumulative total is the sum of the **DENOMINATOR** values for all preceding categories plus the current category. For example, if **DENOMINATOR** represents number of customers, then the cumulative total for any **GROUP** category is the total number of customers for all groups up to and including that group.

The following alternatives are available for specifying maximum count:

**COUNT=ALL.** *No maximum.* This is the default if the subcommand is omitted.

**COUNT=value.** *Use the specified value as the maximum cumulative total.* The value must be a positive integer.

**PERCENT=value.** *Determine the maximum cumulative total based on the specified percentage of the overall total.* The value must be greater than 0 and less than 100.

- If **MAXCOUNT** is specified without **MINRATE**, all rows with a cumulative total for the **DENOMINATOR** variable less than or equal to the specified maximum are displayed with a green background, and all subsequent rows are displayed with a red background.



- If `MINRATE` is also specified, rows are only displayed with a green background if the maximum cumulative total is not exceeded *and* the cumulative response rate is higher than the specified minimum. For example, if the response rate is less than the specified minimum, the row will be displayed with a red background regardless of the cumulative total.

***Example***

```
RESPONSE RATE
/VARIABLES
  NUMERATOR=Responses
  DENOMINATOR=Customers
  GROUP=PostalRank
/MAXCOUNT=6000.
```

# RESTORE

RESTORE

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Overview

RESTORE restores SET specifications that were stored by a previous PRESERVE command. RESTORE and PRESERVE are especially useful when using the macro facility. PRESERVE–RESTORE sequences can be nested up to five levels.

### Basic Specification

The only specification is the command keyword. RESTORE has no additional specifications.

### Limitations

RESTORE does not restore SET THREADS and SET MCACHE settings.

## Example

```
PRESERVE.  
SET TNUMBERS=BOTH.  
FREQUENCIES VARIABLES=DIVISION  
RESTORE.  
FREQUENCIES VARIABLES=DIVISION.
```

- PRESERVE stores all current SET specifications.
- SET TNUMBERS=BOTH changes the TNUMBERS setting to display both values and value labels in pivot tables.
- The pivot table results from the first FREQUENCIES command will display both category values and labels.
- RESTORE reestablishes the SET specifications that were in effect prior to when PRESERVE was specified.
- The pivot table results from the second FREQUENCIES command will use the TNUMBERS setting in effect prior to PRESERVE (the default is to display only labels).

# RMV

```
RMV new variables={LINT (varlist)      }
                  {MEAN (varlist [{,2 }]) }
                  {      {,n }      }
                  {      {ALL}      }
                  {MEDIAN (varlist [{,2 }])}
                  {      {,n }      }
                  {      {ALL}      }
                  {SMEAN (varlist)     }
                  {TREND (varlist)     }

[/new variables=function (varlist [,span])]
```

**Table 214-1**

*Function keywords*

LINT	Linear interpolation
MEAN	Mean of surrounding values
MEDIAN	Median of surrounding values
SMEAN	Variable mean
TREND	Linear trend at that point

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Example**

```
RMV NEWVAR1=LINT(OLDVAR1) .
```

## **Overview**

RMV produces new variables by copying existing variables and replacing any system- or user-missing values with estimates computed by one of several methods. You can also use RMV to replace the values of existing variables. The estimated values are computed from valid data in the existing variables.

### **Basic Specification**

The basic specification is one or more new variable names, an equals sign, a function, and an equal number of existing variables. RMV displays a list of the new variables, the number of missing values replaced, the case numbers of the first and last nonmissing cases, the number of valid cases, and the function used to produce the variables.

### **Syntax Rules**

- The existing variables (and span, if specified) must be enclosed in parentheses.
- The equals sign is required.
- You can specify more than one equation on RMV.

- Equations are separated by slashes.
- You can specify only one function per equation.
- You can create more than one new variable per equation by specifying more than one new variable name on the left and an equal number of existing variables on the right.

### **Operations**

- Each new variable is added to the active dataset.
- If the new variable named already exists, its values are replaced.
- If the new variable named does not already exist, it is created.
- If the same variable is named on both sides of the equation, the new variable will replace the existing variable. Valid values from the existing variable are copied into the new variable, and missing values are replaced with estimates.
- Variables are created in the order in which they are specified on the RMV command.
- If multiple variables are created on a single equation, the first new variable is based on the first existing variable, the second new variable is based on the second existing variable, and so forth.
- RMV automatically generates a variable label for each new variable describing the function and variable used to create it and the date and time of creation.
- The format of a new variable depends on the function specified and on the format of the existing variable.
- RMV honors the TSET MISSING setting that is currently in effect.
- RMV does not honor the USE command.

### **Limitations**

- Maximum 1 function per equation.
- There is no limit on the number of variables created by an equation.
- There is no limit on the number of equations per RMV command.

## **LINT Function**

LINT replaces missing values using linear interpolation. The last valid value before the missing value and the first valid value after the missing value are used for the interpolation.

- The only specification on LINT is a variable or variable list in parentheses.
- LINT will not replace missing values at the endpoints of variables.

### **Example**

```
RMV NEWVAR1=LINT(OLDVAR1) .
```

- This example produces a new variable called *NEWVAR1*.
- *NEWVAR1* will have the same values as *OLDVAR1* but with missing values replaced by linear interpolation.

## MEAN Function

MEAN replaces missing values with the mean of valid surrounding values. The number of surrounding values used to compute the mean depends on the span.

- The specification on MEAN is a variable or variable list and a span, in parentheses.
- The span specification is optional and can be any positive integer or keyword ALL.
- A span of  $n$  uses  $n$  valid cases before and after the missing value.
- If span is not specified, it defaults to 2.
- Keyword ALL computes the mean of all valid values.
- MEAN will not replace missing values if there are not enough valid surrounding cases to satisfy the span specification.

### Example

```
RMV B=MEAN(A, 3) .
```

- This example produces a new variable called  $B$ .
- $B$  will have the same values as variable  $A$  but with missing values replaced by means of valid surrounding values.
- Each mean is based on six values—that is, the three nearest valid values on each side of the missing value.

## MEDIAN Function

MEDIAN replaces missing values with the median of valid surrounding values. The number of surrounding values used to compute the median depends on the span.

- The specification on MEDIAN is a variable or variable list and a span, in parentheses.
- The span specification is optional and can be any positive integer or keyword ALL.
- A span of  $n$  uses  $n$  valid cases before and after the missing value.
- If span is not specified, it defaults to 2.
- Keyword ALL computes the median of all valid values.
- MEDIAN will not replace missing values if there are not enough valid surrounding cases to satisfy the span specification.

### Example

```
RMV B=MEDIAN(A, 3) .
```

- This example produces a new variable called  $B$ .
- $B$  will have the same values as  $A$  but with missing values replaced by medians of valid surrounding values.
- Each median is based on six values—that is, the three nearest valid values on each side of the missing value.

## ***SMEAN Function***

SMEAN replaces missing values in the new variable with the variable mean.

- The only specification on SMEAN is a variable or variable list in parentheses.
- The SMEAN function is equivalent to the MEAN function with a span specification of ALL.

### ***Example***

```
RMV VAR1 TO VAR4=SMEAN(VARA VARB VARC VARD) .
```

- Four new variables (*VAR1*, *VAR2*, *VAR3*, and *VAR4*) are created.
- *VAR1* copies the values of *VARA*, *VAR2* copies *VARB*, *VAR3* copies *VARC*, and *VAR4* copies *VARD*. Any missing values in an existing variable are replaced with the mean of that variable.

## ***TREND Function***

TREND replaces missing values in the new variable with the linear trend for that point. The existing variable is regressed on an index variable scaled 1 to *n*. Missing values are replaced with their predicted values.

- The only specification on TREND is a variable or variable list in parentheses.

### ***Example***

```
RMV YHAT=TREND(VARY) .
```

- This example creates a new variable called *YHAT*.
- *YHAT* has the same values as *VARY* but with missing values replaced by predicted values.

# ROC

ROC is available in the Statistics Base option.

```
ROC varlist BY varname({varvalue })
                        {'varvalue'}

[/MISSING = {EXCLUDE**}]
           {INCLUDE  }

[/CRITERIA = [CUTOFF({INCLUDE**})] [TESTPOS({LARGE**}) [CI({95** })]]
            {EXCLUDE  }           {SMALL  }           {value}

           [DISTRIBUTION({FREE** })]]
           {NEGEXPO}

[/PLOT = [CURVE**[(REFERENCE)]] [NONE]]

[/PRINT = [SE][MODELQUALITY] [COORDINATES]].
```

\*\* Default if subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 18.0

- MODELQUALITY keyword introduced.

## Example

```
ROC
  pred BY default (1).
```

## Overview

ROC produces a receiver operating characteristic (ROC) curve and an estimate of the area under the curve.

### Options

**Distributional Assumptions.** In the CRITERIA subcommand, the user can choose the nonparametric or parametric method to estimate the standard error of the area under the curve. Currently, the bi-negative exponential distribution is the only parametric option.

**Optional Output.** In addition to an estimate of the area under the ROC curve, the user may request its standard error, a confidence interval, and a *p* value under the null hypothesis that the area under the curve equals 0.5. A table of cutoff values and coordinates used to plot the ROC curve may also be displayed.

### **Basic Specification**

The basic specification is one variable as the test result variable and one variable as the actual state variable with one of its values. ROC uses the nonparametric (distribution-free) method to calculate the area under the ROC curve. The default and minimum output is a chart of the ROC curve and a table of the area under the curve.

### **Syntax Rules**

- Minimum syntax: You always need a test result variable and one actual state variable with one of its values in the ROC command line.
- The test result variable must be numeric, but the state variable can be any type with any format.
- Subcommands can be specified in any order.
- When a subcommand is duplicated, only the last one is honored given that all duplicates have no syntax errors. A syntax warning is issued.
- Within a subcommand, if two or more exclusive or contradictory keywords are given, the latter keywords override the earlier ones. A syntax warning is issued.
- If a keyword is duplicated within a subcommand, it is silently ignored.

### **Limitations**

- Only the nonparametric method and one parametric method are available as the computational options for the standard error of the area under the curve at this moment. In the future, binormal and bilogistic distributions may be added to the parametric option.

## **Examples**

```
ROC
  assay BY actual (1)
  /PLOT = CURVE(REFERENCE)
  /PRINT = SE COORDINATES
  /CRITERIA = CUTOFF(INCLUDE) TESTPOS(LARGE) DISTRIBUTION(FREE) CI(95)
  /MISSING = EXCLUDE .
```

- *assay* is the test result variable, and *actual* is the actual state variable. The “positive” group is identified by the value 1.
- The PLOT subcommand specifies that the ROC curve chart should be displayed with the diagonal reference line.
- The PRINT subcommand specifies that the standard error of the area estimate should be computed and that the coordinates of the ROC curve should be displayed in a separate table.
- All other options are set to their default values.

### **Comparing Multiple Curves**

```
ROC
  preddef1 preddef2 preddef3 BY default (1)
  /PLOT = CURVE(REFERENCE)
  /PRINT = SE
  /CRITERIA = CUTOFF(INCLUDE) TESTPOS(LARGE) DISTRIBUTION(FREE) CI(95)
  /MISSING = EXCLUDE .
```



- *preddef1*, *preddef2*, and *preddef3* are the test result variables, and *default* is the actual state variable. The “positive” group is identified by the value 1.
- The PLOT subcommand specifies that the ROC curve chart should be displayed with the diagonal reference line.
- The PRINT subcommand specifies that the standard error of the area estimate should be computed and that the coordinates of the ROC curve should be displayed in a separate table.
- All other options are set to their default values.

### ***varlist BY varname(varvalue)***

The *varlist* specifies one or more test result variables. They must be of numeric type.

The *varname* separated from the *varlist* by the word *BY* specifies the actual state variable. It can be of any type and any format. The user must also specify a *varvalue* in brackets after the second *varname* to define the “positive” group of the actual state variable. All other valid state values are assumed to indicate the negative group.

### ***MISSING Subcommand***

Cases with system-missing values in the test result variable and the actual state variable are always excluded from the analysis. However, the *MISSING* subcommand allows the user to redefine the user-missing values to be valid.

- |                |   |
|----------------|---|
| <b>EXCLUDE</b> | <i>Exclude both user-missing and system-missing values. Cases with a system-missing value or a user-missing value in either the test result variable or the actual state variable are excluded from the analysis. This is the active default.</i> |
| <b>INCLUDE</b> | <i>User-missing values are treated as valid. Cases with a system-missing value in either the test result variable or the actual state variable are excluded from the analysis.</i>  |

### ***CRITERIA Subcommand***

The *CRITERIA* subcommand allows the user to decide whether or not the cutoff value is included as the positive test result value, whether or not the larger or smaller value direction of the test result variable indicates the positive test result, the confidence level of the asymptotic confidence interval produced by */PRINT = SE*, and the estimation method for the standard error of the area under the curve.

- |                        |   |
|------------------------|---|
| <b>CUTOFF(INCLUDE)</b> | <i>Positive classification of test result values includes the cutoff values. Note that the positive test result direction is controlled by the TESTPOS keyword. This is the active default.</i>   |
| <b>CUTOFF(EXCLUDE)</b> | <i>Positive classification of test result values excludes the cutoff values. This distinction leads to the different sets of cutoff values, but none of the output (chart or table) is affected at this moment because the user cannot choose the cutoff values for the classification.</i> |
| <b>TESTPOS(LARGE)</b>  | <i>The user can specify which direction of the test result variable indicates increasing strength of conviction that the subject is test positive. The larger the test result value is, the more positive the test result is. LARGE is the active default.</i>                              |

<b>TESTPOS(SMALL)</b>	<i>The smaller the test result value is, the more positive the test result is.</i>
<b>CI(n)</b>	<i>Confidence level in the (two-sided) asymptotic confidence interval of the area. The user can specify any confidence level in (0, 100) for the asymptotic confidence interval created by /PRINT = SE. The active default parameter value is 95.</i>
<b>DISTRIBUTION(FREE)</b>	<i>The method of calculating the standard error of the area under the curve. When FREE, the standard error of the area under the curve is estimated nonparametrically—that is, without any distributional assumption.</i>
<b>DISTRIBUTION(NEGEXPO)</b>	<i>The standard error is estimated assuming the bi-negative exponential distribution. This latter option is valid only when the number of positive actual state observations equals the number of negative actual state observations.</i>

## ***PRINT Subcommand***

The PRINT subcommand controls the display of optional table and chart output. Note that the area under the curve is always displayed.

<b>SE (MODELQUALITY)</b>	<i>Standard error of the area estimate. In addition to the standard error of the estimated area under the curve, the asymptotic 95% (or other user-specified confidence level) confidence interval as well as the asymptotic <math>p</math> value under the null hypothesis that the true area = 0.5 are calculated. The optional (MODELQUALITY) specification produces a simple bar chart that displays the value of the lower bound of the confidence interval of the area under the curve, expressed as a proportion, on a scale of 0 to 1. Note that the area under the curve statistic is asymptotically normally distributed.</i>
<b>COORDINATES</b>	<i>Coordinate points of the ROC curve along with the cutoff values. Pairs of sensitivity and 1 – specificity values are given with the cutoff values for each curve.</i>

## ***PLOT Subcommand***

The PLOT subcommand controls the display of chart output.

<b>CURVE(REFERENCE)</b>	<i>The ROC curve chart is displayed. The keyword CURVE is the active default. In addition, the user has an option to draw the diagonal reference line (sensitivity = 1 – specificity) by the bracketed parameter REFERENCE.</i>
<b>NONE</b>	<i>No charts are displayed.</i>

# SAMPLE

```
SAMPLE {decimal value}  
      {n FROM m      }
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
SAMPLE .25.
```

## **Overview**

SAMPLE permanently draws a random sample of cases for processing in all subsequent procedures. For a temporary sample, use a TEMPORARY command before SAMPLE.

### **Basic Specification**

The basic specification is either a decimal value between 0 and 1 or the sample size followed by keyword FROM and the size of the active dataset.

- To select an approximate percentage of cases, specify a decimal value between 0 and 1.
- To select an exact-size random sample, specify a positive integer that is less than the file size, and follow it with keyword FROM and the file size.

### **Operations**

- SAMPLE is a permanent transformation.
- Sampling is based on a pseudo-random-number generator that depends on a seed value that is established by the program. On some implementations of the program, this number defaults to a fixed integer, and a SAMPLE command that specifies  $n$  FROM  $m$  will generate the identical sample whenever a session is rerun. To generate a different sample each time, use the SET command to reset SEED to a different value for each session. See the SET command for more information.
- If sampling is done by using the  $n$  FROM  $m$  method, and the TEMPORARY command is specified, successive samples will not be the same because the seed value changes each time that a random-number series is needed within a session.
- A proportional sample (a sample that is based on a decimal value) usually does not produce the exact proportion that is specified.
- If the number that is specified for  $m$  following FROM is less than the actual file size, the sample is drawn only from the first  $m$  cases.
- If the number following FROM is greater than the actual file size, the program samples an equivalent proportion of cases from the active dataset.
- If SAMPLE follows SELECT IF, SAMPLE samples only cases that are selected by SELECT IF.

- If `SAMPLE` precedes `SELECT IF`, cases are selected from the sample.
- If more than one `SAMPLE` command is specified in a session, each command acts on the sample that was selected by the preceding `SAMPLE` command.
- If `N OF CASES` is used with `SAMPLE`, the program reads as many records as required to build the specified  $n$  cases. It makes no difference whether `N OF CASES` precedes or follows `SAMPLE`.

### **Limitations**

`SAMPLE` cannot be placed in a `FILE TYPE-END FILE TYPE` or `INPUT PROGRAM-END INPUT PROGRAM` structure. `SAMPLE` can be placed nearly anywhere following these commands in a transformation program. For more information, see the topic [Commands and Program States](#) on p. 2196..

## **Examples**

### ***Sampling an Approximate Percentage of Cases***

```
SAMPLE .25.
```

- This command samples approximately 25% of the cases in the active dataset.

### ***Sampling a Specific Number of Cases***

```
SAMPLE 500 FROM 3420.
```

- The active dataset must have 3420 or more cases to obtain a random sample of exactly 500 cases.
- If the file contains fewer than 3420 cases, proportionally fewer cases are sampled.
- If the file contains more than 3420 cases, a random sample of 500 cases is drawn from the first 3420 cases.

### ***Sampling Subgroups***

```
DO IF SEX EQ 'M'.
SAMPLE 1846 FROM 8000.
END IF.
```

- `SAMPLE` is placed inside a `DO IF-END IF` structure to sample subgroups differently. Assume that this survey is for 10,000 people in which 80% of the sample is male, while the known universe is 48% male. To obtain a sample that corresponds to the known universe and that maximizes the size of the sample, 1846 ( $48/52 * 2000$ ) males and all females must be sampled. The `DO IF` structure restricts the sampling process to the males.

# SAVE

```
SAVE OUTFILE='filespec'  
  
  [/VERSION={3**}]  
    {2 }  
  
  [/UNSELECTED=[{RETAIN}]  
    {DELETE}]  
  
  [/KEEP={ALL** }] [/DROP=varlist]  
    {varlist}  
  
  [/RENAME=(old varlist=new varlist)...]  
  
  [/MAP]  [{COMPRESSED }]  
    {UNCOMPRESSED}  
  
  [/NAMES]  
  
  [/PERMISSIONS={READONLY }  
    {WRITEABLE}]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
SAVE OUTFILE='/data/empl.sav'.
```

## Overview

SAVE produces a data file in IBM® SPSS® Statistics format, which contains data plus a dictionary. The dictionary contains a name for each variable in the data file plus any assigned variable and value labels, missing-value flags, and variable print and write formats. The dictionary also contains document text that was created with the DOCUMENTS command.

XSAVE also creates SPSS Statistics data files. The difference is that SAVE causes data to be read, while XSAVE is not executed until data are read for the next procedure.

See [SAVE TRANSLATE](#) on p. 1822 for information about saving data files that can be used by other programs.

## Options

**Compatibility with Earlier Releases.** You can save a data file that can be read by releases prior to 7.5.

**Variable Subsets and Order.** You can use the DROP and KEEP subcommands to save a subset of variables and reorder the variables that are saved.

**Filtered Cases.** If filtering is in effect, you use the UNSELECTED subcommand to specify inclusion or exclusion of filtered cases. By default, all cases are included.

**Variable Names.** You can use the `RENAME` subcommand to rename variables as they are copied into the SPSS Statistics data file.

**Variable Map.** To confirm the names and order of the variables that are saved in the data file, use the `MAP` subcommand. `MAP` displays the variables that are saved in the data file next to their corresponding names in the active dataset.

**Data Compression.** You can use the `COMPRESSED` or `UNCOMPRESSED` subcommand to write the data file in compressed or uncompressed form.

### ***Basic Specification***

The basic specification is the `OUTFILE` subcommand, which specifies a name for the data file to be saved.

### ***Subcommand Order***

- Subcommands can be specified in any order.

### ***Syntax Rules***

- `OUTFILE` is required and can be specified only once. If `OUTFILE` is specified more than once, only the last specified `OUTFILE` is in effect.
- `KEEP`, `DROP`, `RENAME`, and `MAP` can each be used as many times as needed.
- Only one of the subcommands `COMPRESSED` or `UNCOMPRESSED` can be specified per `SAVE` command.

### ***Operations***

- `SAVE` is executed immediately and causes the data to be read.
- The new data file dictionary is arranged in the same order as the active dataset dictionary, unless variables are reordered with the `KEEP` subcommand. Documentary text from the active dataset dictionary is always saved unless it is dropped with the `DROP DOCUMENTS` command before `SAVE`.
- New variables that were created by transformations and procedures that occurred prior to the `SAVE` command are included in the new data file, and variables that were altered by transformations are saved in their modified form. Results of any temporary transformations that immediately precede the `SAVE` command are included in the file; scratch variables are not included.
- The active dataset is still available for transformations and procedures after `SAVE` is executed.
- `SAVE` processes the dictionary first and displays a message that indicates how many variables will be saved. After the data are written, `SAVE` indicates how many cases were saved. If the second message does not appear, the file was probably not completely written.

## Examples

### Renaming Variables for a Saved Version

```
GET FILE='/hubempl.sav'.
SAVE OUTFILE='/data/empl88.sav' /RENAME=(AGE=AGE88) (JOB CAT=JOB CAT88).
```

- The GET command retrieves the IBM® SPSS® Statistics data file *hubempl.sav*.
- The RENAME subcommand renames variable *AGE* to *AGE88* and variable *JOB CAT* to *JOB CAT88*.
- SAVE causes the data to be read and saves a new data file with the filename *empl88.sav*. The original data file *hubempl.sav* is not changed.

### Performing Temporary Transformations Prior to Saving

```
GET FILE='/data/hubempl.sav'.
TEMPORARY.
RECODE DEPT85 TO DEPT88 (1,2=1) (3,4=2) (ELSE=9).
VALUE LABELS DEPT85 TO DEPT88 1 'MANAGEMENT' 2 'OPERATIONS' 9 'UNKNOWN'.
SAVE OUTFILE='/data/hubtemp.sav'.
CROSSTABS DEPT85 TO DEPT88 BY JOB CAT.
```

- The GET command retrieves the SPSS Statistics data file *hubempl.sav*.
- The TEMPORARY command indicates that RECODE and VALUE LABELS are in effect only for the next command that reads the data (SAVE).
- The RECODE command recodes values for all variables between and including *DEPT85* and *DEPT88* on the active dataset.
- The VALUE LABELS command specifies new labels for the recoded values.
- The OUTFILE subcommand on SAVE specifies *hubtemp.sav* as the new data file. The *hubtemp.sav* file will include the recoded values for *DEPT85* to *DEPT88* and the new value labels.
- The CROSSTABS command crosstabulates *DEPT85* to *DEPT88* with *JOB CAT*. Because the RECODE and VALUE LABELS commands were temporary, the CROSSTABS output does not reflect the recoding and new labels.
- If XSAVE were specified instead of SAVE, the data would be read only once. Both the saved data file and the CROSSTABS output would reflect the temporary recoding and labeling of the department variables.

## OUTFILE Subcommand

OUTFILE specifies the data file to be saved. OUTFILE is required and can be specified only once. If OUTFILE is specified more than once, only the last OUTFILE is in effect.

## **VERSION Subcommand**

VERSION allows you to save a data file that can be opened in releases prior to 7.5. If VERSION is not specified or is specified with no value, the default of 3 is used, which saves the file in a format that cannot be read by releases prior to 7.5. Specify 2 to save a file that is compatible with earlier releases.

## **Variable Names**

For data files with variable names that are longer than eight bytes in version 10.x or 11.x, unique, eight-byte versions of variable names are used (but the original variable names are preserved for use in release 12.0 or later). In releases prior to version 10.0, the original long variable names are lost if you save the data file.

## **UNSELECTED Subcommand**

UNSELECTED determines whether cases that were excluded on a previous FILTER or USE command are to be retained or deleted in the saved data file. The default is RETAIN. The UNSELECTED subcommand has no effect when the active dataset does not contain unselected cases.

- |               |  |
|---------------|--|
| <b>RETAIN</b> | <i>Retain the unselected cases.</i> All cases in the active dataset are saved. This setting is the default when UNSELECTED is specified by itself. |
| <b>DELETE</b> | <i>Delete the unselected cases.</i> Only cases that meet the FILTER or USE criteria are saved in the data file.                                    |

## **DROP and KEEP Subcommands**

DROP and KEEP are used to save a subset of variables. DROP specifies the variables that are not to be saved in the new data file; KEEP specifies the variables that are to be saved in the new data file; variables that are not named on KEEP are dropped.

- Variables can be specified in any order. The order of variables on KEEP determines the order of variables in the data file. The order on DROP does not affect the order of variables in the data file.
- Keyword ALL on KEEP refers to all remaining variables that were not previously specified on KEEP or DROP. ALL must be the last specification on KEEP.
- If a variable is specified twice on the same subcommand, only the first mention is recognized.
- Multiple DROP and KEEP subcommands are allowed. If a variable is specified that is not in the active dataset, or that has been dropped because of a previous DROP or KEEP subcommand, an error results, and the SAVE command is not executed.
- Keyword TO can be used to specify a group of consecutive variables in the active file.

### **Example**

```
GET FILE='/data/personnel.sav'.  
COMPUTE TENURE=(12-CMONTH +(12*(88-CYEAR)))/12.
```



```

COMPUTE JTENURE=(12-JMONTH +(12*(88-JYEAR)))/12.
VARIABLE LABELS      TENURE 'Tenure in Company'
                    JTENURE 'Tenure in Grade'.
SAVE OUTFILE='/data/personnel88.sav' /DROP=GRADE STORE
  /KEEP=LNAME NAME TENURE JTENURE ALL.

```

- The variables *TENURE* and *JTENURE* are created by `COMPUTE` commands and are assigned variable labels by the `VARIABLE LABELS` command. *TENURE* and *JTENURE* are added to the end of the active dataset.
- `DROP` excludes variables *GRADE* and *STORE* from file *personnel88.sav*. `KEEP` specifies that *LNAME*, *NAME*, *TENURE*, and *JTENURE* are the first four variables in file *personnel88.sav*, followed by all remaining variables not specified on `DROP`. These remaining variables are saved in the same sequence as they appear in the original file.

## **RENAME Subcommand**

`RENAME` changes the names of variables as they are copied into the new data file.

- The specification on `RENAME` is a list of old variable names followed by an equals sign and a list of new variable names. The same number of variables must be specified on both lists. Keyword `TO` can be used in the first list to refer to consecutive variables in the active dataset and can be used in the second list to generate new variable names. The entire specification must be enclosed in parentheses.
- Alternatively, you can specify each old variable name individually, followed by an equals sign and the new variable name. Multiple sets of variable specifications are allowed. The parentheses around each set of specifications are optional.
- `RENAME` does not affect the active dataset. However, if `RENAME` precedes `DROP` or `KEEP`, variables must be referred to by their new names on `DROP` or `KEEP`.
- Old variable names do not need to be specified according to their order in the active dataset.
- Name changes take place in one operation. Therefore, variable names can be exchanged between two variables.
- Multiple `RENAME` subcommands are allowed.

### **Examples**

```

SAVE OUTFILE='/data/empl88.sav'
  /RENAME AGE=AGE88 JOBCAT=JOBCAT88.

```

- `RENAME` specifies two name changes for the file *empl88.sav*: The variable *AGE* is renamed to *AGE88*, and the variable *JOBCAT* is renamed to *JOBCAT88*.

```

SAVE OUTFILE='/data/empl88.sav'
  /RENAME (AGE JOBCAT=AGE88 JOBCAT88).

```

- The name changes are identical to the changes in the previous example: *AGE* is renamed to *AGE88*, and *JOBCAT* is renamed to *JOBCAT88*. The parentheses are required with this method.

## **MAP Subcommand**

MAP displays a list of the variables in the saved data file and their corresponding names in the active dataset.

- The only specification is keyword MAP.
- Multiple MAP subcommands are allowed. Each MAP subcommand maps the results of subcommands that precede it (but not results of subcommands that follow it).

### **Example**

```
GET FILE='/data/hubempl.sav'.  
SAVE OUTFILE='/data/empl88.sav' /RENAME=(AGE=AGE88) (JOB CAT=JOB CAT88)  
/KEEP=LNAME NAME JOB CAT88 ALL /MAP.
```

- MAP confirms the new names for *AGE* and *JOB CAT* and the order of variables in the *empl88.sav* file (*LNAME*, *NAME*, and *JOB CAT88*, followed by all remaining variables from the active dataset).

## **COMPRESSED and UNCOMPRESSED Subcommands**

COMPRESSED saves the file in compressed form. UNCOMPRESSED saves the file in uncompressed form. In a compressed file, small integers (from -99 to 155) are stored in one byte instead of the eight bytes that are used in an uncompressed file.

- The only specification is the keyword COMPRESSED or UNCOMPRESSED.
- Compressed data files occupy less disk space than uncompressed data files.
- Compressed data files take longer to read than uncompressed data files.
- The GET command, which reads IBM® SPSS® Statistics data files, does not need to specify whether the files that it reads are compressed or uncompressed.
- Only one of the subcommands COMPRESSED or UNCOMPRESSED can be specified per SAVE command. COMPRESSED is usually the default, though UNCOMPRESSED may be the default on some systems.

## **NAMES Subcommand**

For all variable names that are longer than eight bytes, NAMES displays the eight-byte equivalents that will be used if you read the data file into a version prior to release 12.0.

---

## ***PERMISSIONS Subcommand***

The `PERMISSIONS` subcommand sets the operating system read/write permissions for the file.

- |                  |  |
|------------------|--|
| <b>READONLY</b>  | <i>File permissions are set to read-only for all users. The file cannot be saved by using the same file name with subsequent changes unless the read/write permissions are changed in the operating system or the subsequent <code>SAVE</code> command specifies <code>PERMISSIONS=WRITEABLE</code>.</i> |
| <b>WRITEABLE</b> | <i>File permissions are set to allow writing for the file owner. If file permissions were set to read-only for other users, the file remains read-only for them.</i>   |

Your ability to change the read/write permissions may be restricted by the operating system.

## ***Data File Compatibility with Previous Releases***

A data file saved in Unicode mode (see [SET command, Unicode subcommand](#)) cannot be read by versions of IBM® SPSS® Statistics prior to 16.0. To save a Unicode data file in a format that can be read by earlier releases, open the file in code page mode and re-save it. The file will be saved in the encoding based on the current locale setting (see [SET command, LOCALE subcommand](#)). Some data loss may occur if the file contains characters not recognized by the current locale.

# SAVE DATA COLLECTION

This command is available only on Windows operating systems.

```
SAVE DATA COLLECTION
  /OUTFILE='filespec'
  /METADATA='filespec'
  [/UNSELECTED={RETAIN**}]
    {DELETE}
  [/KEEP={ALL** }] [/DROP=varlist]
    {varlist}
  [/MAP]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release history**

Release 15.0

- Command introduced as SAVE DIMENSIONS.

Release 18.0

- SAVE DIMENSIONS deprecated. Command name changed to SAVE DATA COLLECTION.

## **Example**

```
SAVE DATA COLLECTION
  /OUTFILE=' /data/survey.sav'
  /METADATA=' /data/survey.mdd' .
```

## **Overview**

SAVE DATA COLLECTION produces IBM® SPSS® Statistics data files and IBM® SPSS® Data Collection metadata files that you can use to read the SPSS Statistics data file into Data Collection applications such as mrTables and mrInterview. This is particularly useful when “roundtripping” data between SPSS Statistics and Data Collection applications. For example, you can read an IBM® SPSS® Data Collection Interviewer Server data source into SPSS Statistics, calculate some new variables, and then save the data in a form that can be read by mrTables, without loss of any of the original metadata attributes.

### **Options**

**Variable subsets and order.** You can use the `DROP` and `KEEP` subcommands to save a subset of variables and reorder the variables that are saved.

**Filtered cases.** If filtering is in effect, you use the `UNSELECTED` subcommand to specify inclusion or exclusion of filtered cases. By default, all cases are included.

**Variable map.** To confirm the names and order of the variables that are saved in the SPSS Statistics data file, use the `MAP` subcommand. `MAP` displays the variables that are saved in the SPSS Statistics data file next to their corresponding names in the active dataset.

### **Basic specification**

The basic specification is the `OUTFILE` subcommand that specifies a name for the SPSS Statistics data file to be saved and the `METADATA` subcommand that specifies the name for the Data Collection metadata file.

### **Subcommand order**

- Subcommands can be specified in any order.

### **Syntax rules**

- `OUTFILE` and `METADATA` are required and can be specified only once.
- `KEEP`, `DROP`, `MAP`, and `UNSELECTED` are optional.

### **Operations**

For new variables and datasets not created from Data Collection data sources, SPSS Statistics variable attributes are mapped to Data Collection metadata attributes in the metadata file according to the methods described in the SAV DSC documentation in the IBM® SPSS® Data Collection Developer Library.

If the active dataset was created from a Data Collection data source:

- The new metadata file is created by merging the original metadata attributes with metadata attributes for any new variables, plus any changes to original variables that might affect their metadata attributes (for example, addition of, or changes to, value labels).
- For original variables read from the Data Collection data source, any metadata attributes not recognized by SPSS Statistics are preserved in their original state. For example, SPSS Statistics converts grid variables to regular SPSS Statistics variables, but the metadata that defines these grid variables is preserved when you save the new metadata file.
- If any Data Collection variables were automatically renamed to conform to SPSS Statistics variable naming rules, the metadata file maps the converted names back to the original Data Collection variable names.

The presence or absence of value labels can affect the metadata attributes of variables and consequently the way those variables are read by Data Collection applications. If value labels have been defined for any nonmissing values of a variable, they should be defined for all

nonmissing values of that variable; otherwise, the unlabeled values will be dropped when the data file is read by Data Collection.

### **Limitations**

This feature is only available with SPSS Statistics installed on Microsoft Windows operating systems, and is only available in local analysis mode. This feature is not available in distributed analysis mode using SPSS Statistics Server.

To write Data Collection metadata files, you must have the following items installed:

- .NET framework. To obtain the most recent version of the .NET framework, go to <http://www.microsoft.com/net>.
- IBM® SPSS® Data Collection Survey Reporter Developer Kit. For information on obtaining a compatible version of SPSS Survey Reporter Developer Kit, go to [support.spss.com](http://support.spss.com) (<http://support.spss.com>).

## **OUTFILE subcommand**

OUTFILE specifies the IBM® SPSS® Statistics data file to be saved.

- OUTFILE is required and can be specified only once.
- File specifications should be enclosed in quotes.
- The standard extension for SPSS Statistics data files is *.sav*. You should explicitly specify the extension; it will not be automatically appended to the file name.

## **METADATA subcommand**

METADATA specifies the IBM® SPSS® Data Collection metadata file to be saved. See the Operations section of the [Overview](#) for more information.

- METADATA is required and can be specified only once.
- File specifications should be enclosed in quotes.
- The standard extension for a Data Collection metadata file is *.mdd*. You should explicitly specify the extension; it will not be automatically appended to the file name.

## **UNSELECTED subcommand**

UNSELECTED determines whether cases that were excluded on a previous FILTER or USE command are to be retained or deleted in the IBM® SPSS® Statistics data file. The default is RETAIN. The UNSELECTED subcommand has no effect when the active dataset does not contain unselected cases. Since the metadata file does not contain case data, this subcommand has no effect on the contents of the metadata file.

- |               |   |
|---------------|---|
| <b>RETAIN</b> | <i>Retain the unselected cases.</i> All cases in the active dataset are saved. This is the default.                             |
| <b>DELETE</b> | <i>Delete the unselected cases.</i> Only cases that meet the FILTER or USE criteria are saved in the SPSS Statistics data file. |

## ***DROP and KEEP subcommands***

DROP and KEEP are used to save a subset of variables. DROP specifies the variables that are not to be saved in the new data file; KEEP specifies the variables that are to be saved in the new data file; variables that are not named on KEEP are dropped.

- Variables can be specified in any order. The order of variables on KEEP determines the order of variables in the data file. The order on DROP does not affect the order of variables in the data file.
- Keyword ALL on KEEP refers to all remaining variables that were not previously specified on KEEP or DROP. ALL must be the last specification on KEEP.
- If a variable is specified twice on the same subcommand, only the first mention is recognized.
- Multiple DROP and KEEP subcommands are allowed. If a variable is specified that is not in the active dataset or that has been dropped because of a previous DROP or KEEP subcommand, an error results, and the SAVE DATA COLLECTION command is not executed.
- Keyword TO can be used to specify a group of consecutive variables in the active file.
- If the active dataset was created from a IBM® SPSS® Data Collection data source, any original variables defined as grid or array elements in the Data Collection data source are retained in the metadata file, even if those variables are not included in the IBM® SPSS® Statistics data file. Thus, the original grid or array structure is preserved, but there will be no case data for any variables not included in the SPSS Statistics data file.

### ***Example***

```
SAVE DATA COLLECTION
  /OUTFILE=' /data/survey.sav'
  /METADATA=' /data/survey.mdd'
  /DROP gridVar7
  /KEEP gridVar1 to gridVar8.
```

Assuming that the order of the variables in the active dataset is *gridVar1*, *gridVar2*, *gridVar3*,...*gridVar8* and that all eight variables are grid variables in the original Data Collection data source, *gridVar7* will be dropped from *survey.sav*, but the original metadata for *gridVar7* will be preserved in *survey.mdd*.

## ***MAP subcommand***

MAP displays a list of the variables in the IBM® SPSS® Statistics data file and their corresponding names in the active dataset.

- The only specification is keyword MAP.
- Multiple MAP subcommands are allowed. Each MAP subcommand maps the results of the DROP and KEEP subcommands that precede it (but not results of subcommands that follow it).

### ***Example***

```
SAVE DATA COLLECTION
  /OUTFILE=' /data/survey.sav'
  /METADATA=' /data/survey.mdd'
  /DROP var7 /MAP
```

---

*SAVE DATA COLLECTION*

*/KEEP var1 to var8 /MAP.*



# SAVE MODEL

```
SAVE MODEL OUTFILE='filename'  
  
  [/KEEP={ALL**          }]  
        {model names}  
        {procedures }  
  
  [/DROP={model names}]  
        {procedures }  
  
  [/TYPE={MODEL**}]  
        {COMMAND}
```

\*\*Default if the subcommand is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
SAVE MODEL OUTFILE='ACFMODE.DAT' .
```

## Overview

SAVE MODEL saves the models that are created by certain procedures in a model file. The saved model file can be read later in the session or in another session with the READ MODEL command.

## Options

You can save a subset of models into the file using the DROP and KEEP subcommands. You can use the TYPE subcommand to control whether models are specified by model name or by the name of the procedure that generated them.

## Basic Specification

The basic specification is the OUTFILE subcommand followed by a filename.

- By default, SAVE MODEL saves all currently active models in the specified file. Each model that is saved in the file includes information such as the procedure that created it, the model name, the specified variable names, subcommands and specifications used, and parameter estimates. The names of the models are either the default *MOD\_n* names or the names that are assigned on the MODEL NAME command. In addition to the model specifications, the TSET settings that are currently in effect are saved.

## Subcommand Order

- Subcommands can be specified in any order.

**Syntax Rules**

- If a subcommand is specified more than once, only the last subcommand is executed.

**Operations**

- `SAVE MODEL` is executed immediately.
- Model files are designed to be written and read by specific procedures and should not be edited.
- The active models are not affected by the `SAVE MODEL` command.
- `DATE` specifications are not saved in the model file.
- Models are not saved in IBM® SPSS® Statistics data files.
- The following procedures can generate models that can be saved with the `SAVE MODEL` command: `AREG`, `ARIMA`, `EXSMOOTH`, `SEASON`, and `SPECTRA` in the Forecasting module; `ACF`, `CASEPLOT`, `CCF`, `CURVEFIT`, `PACF`, `PLOT`, and `TSPLOT` in the Core system; and `WLS` and `2SLS` in the Regression module.

**Limitations**

- A maximum of one filename can be specified.
- The `READ MODEL` command can only read models created by a `SAVE MODEL` command on the same operating system/platform.

**OUTFILE Subcommand**

`OUTFILE` names the file where models will be stored and is the only required subcommand.

- The only specification on `OUTFILE` is the name of the model file.
- The filename should be enclosed in quotes.
- Only one filename can be specified.
- You can store models in other directories by specifying a fully qualified filename.

**KEEP and DROP Subcommands**

`KEEP` and `DROP` allow you to save a subset of models. By default, all currently active models are saved.

- `KEEP` specifies models to be saved in the model file.
- `DROP` specifies models that are not saved in the model file.
- Models can be specified by using individual model names or the names of the procedures that created them. To use procedure names, you must specify `COMMAND` on the `TYPE` subcommand.
- Model names are either the default `MOD_n` names or the names that are assigned with `MODEL NAME`.
- If you specify a procedure name on `KEEP`, all models that are created by that procedure are saved; on `DROP`, any models created by that procedure are not included in the model file.
- Model names and procedure names cannot be mixed on a single `SAVE MODEL` command.

- If more than one KEEP or DROP subcommand is specified, only the last subcommand is executed.
- You can specify the keyword ALL on KEEP to save all models that are currently active. This setting is the default.

**Example**

```
SAVE MODEL OUTFILE='ACFCCF.DAT'  
/KEEP=ACF1 ACF2.
```

- In this example, only models *ACF1* and *ACF2* are saved in model file *ACFCCF.DAT*.

## **TYPE Subcommand**

TYPE indicates whether models are specified by model name or procedure name on DROP and KEEP.

- One keyword, MODEL or COMMAND, can be specified after TYPE.
- MODEL is the default and indicates that models are specified as model names.
- COMMAND indicates that the models are specified by procedure name.
- TYPE has no effect if KEEP or DROP is not specified.
- The TYPE specification applies only to the current SAVE MODEL command.

**Example**

```
SAVE MODEL OUTFILE='CURVE1.DAT'  
/KEEP=CURVEFIT  
/TYPE=COMMAND.
```

- This command saves all models that were created by the CURVEFIT procedure into the model file *CURVE1.DAT*.

# SAVE TRANSLATE

```
SAVE TRANSLATE

[{/OUTFILE=file          }]
{/CONNECT=ODBC connect string}

[{/UNENCRYPTED}]
{/ENCRYPTED  }

[/TYPE={CSV  }]
      {DB2  }
      {DB3  }
      {DB4  }
      {ODBC }
      {PC   }
      {SAS  }
      {STATA}
      {SYM  }
      {SLK  }
      {TAB  }
      {WKS  }
      {WK1  }
      {WK3  }
      {XLS  }

[/VERSION={1}]
      {2}
      {3}
      {4}
      {5}
      {6}
      {7}
      {9}
      {8}
      {12}
      {X}

[/ENCODING={'Locale'}]
      {'System'}
      {'value'  }

[/FIELDNAMES]

[/CELLS={VALUES}]
      {LABELS}

[/TEXTOPTIONS [DELIMITER='char'] [QUALIFIER='char']
      [DECIMAL={DOT  }]
      {COMMA}
      [FORMAT={PLAIN  }]]
      {VARIABLE}

[/EDITION={INTERCOOLED}
      {SE      }

[/PLATFORM={ALPHA}]
      {WINDOWS}
      {UNIX}

[{/VALFILE=filename}]

[/TABLE = 'database table name']

[/SQL = 'SQL statement' [/SQL = 'SQL statement']]

[{/APPEND }]
{/REPLACE}

[/RENAME=(old varlist=new varlist)[(...)]
```

```

[/KEEP={ALL      }]
      {varlist}

[/DROP=varlist]

[/UNSELECTED=[ {RETAIN}]
              {DELETE}]

[/MISSING={IGNORE}]
          {RECODE}

[ {/COMPRESSED  }]
[ {/UNCOMPRESSED}]

[/MAP]

```

- OUTFILE is not valid for TYPE=ODBC but is required for all other types.
- CONNECT is required for TYPE=ODBC but is invalid for all other types.
- APPEND, SQL, ENCRYPTED, and TABLE are available only for TYPE=ODBC.
- ENCODING is available only for TYPE=SAS and TYPE=STATA. For tab-delimited (TYPE=TAB) and CSV (TYPE=CSV) text files, the encoding is UTF-8 in Unicode mode or the encoding of the current code page in code page mode. For more information on Unicode mode, see [UNICODE Subcommand](#).
- FIELDNAMES is available for spreadsheet formats and text data formats.
- PLATFORM and VALFILE are available only for TYPE=SAS, and PLATFORM is required.
- CELLS is available for TYPE=XLS /VERSION=8, TYPE=TAB, and TYPE=CSV.
- TEXTOPTIONS is available only for TYPE=TAB and TYPE=CSV.
- EDITION is available only for TYPE=STATA.
- COMPRESSED and UNCOMPRESSED are available only for TYPE=PC.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

#### Release 14.0

- Value STATA added to list for TYPE subcommand.
- EDITION subcommand introduced for TYPE=STATA.
- SQL subcommand introduced.
- MISSING subcommand introduced.
- Field/column names specified on the RENAME subcommand can contain characters (for example, spaces, commas, slashes, plus signs) that are not allowed in IBM® SPSS® Statistics variable names.
- Continuation lines for connection strings on the CONNECT subcommand do not need to begin with a plus sign.

#### Release 15.0

- ENCRYPTED subcommand introduced.

- Value `CSV` added to list for `TYPE` subcommand.
- `TEXTOPTIONS` subcommand introduced for `TYPE=CSV` and `TYPE=TAB`.

#### Release 16.0

- `VERSION=12` introduced for writing data in Excel 2007 XLSX format with `TYPE=XLS`.

#### Release 17.0

- `UNENCRYPTED` subcommand introduced.

#### Release 18.0

- `VERSION=9` introduced for writing SAS 9+ files with `TYPE=SAS`.

#### Release 19

- `ENCODING` subcommand introduced.

#### **Example**

```
SAVE TRANSLATE OUTFILE='/data/sales.xls'.
```

## **Overview**

`SAVE TRANSLATE` translates the active dataset into a file that can be used by other software applications. Supported formats are 1-2-3, Symphony, Multiplan, Excel, dBASE II, dBASE III, dBASE IV, tab-delimited text files, comma-delimited text files, SAS, Stata, SPSS/PC+ data files, and ODBC database sources. For ODBC database sources, it can also manipulate database tables.

### **Options**

**Variable Subsets.** You can use the `DROP` and `KEEP` subcommands to specify variables to omit or retain in the resulting file.

**Variable Names.** You can use the `RENAME` subcommand to rename variables as they are copied to the external file format. For spreadsheets and text data files, use `FIELDNAMES` to include the variable names in the file.

**Variable Map.** To confirm the names and order of the variables that are saved in the resulting file, use the `MAP` subcommand. `MAP` displays the variables that are saved in the file next to their corresponding names in the active dataset.

**Value Labels.** For Excel spreadsheets and tab-delimited and comma-delimited text files, you can use the `CELLS` subcommand to export value labels rather than values in spreadsheet files. For SAS files, you can use the `VALFILE` subcommand to create a SAS syntax file containing your data's value labels. For Stata files, value labels for numeric variables are automatically saved as Stata value labels.

### **Limitations**

- A maximum of 2,048 cases can be saved to 1-2-3 Release 1A; a maximum of 8,192 cases can be saved to 1-2-3 Release 2.0 or Symphony files.

- A maximum of 16,384 cases can be saved to Excel 2.1. For Excel 95, Excel 97 and Excel 2007, if the number of cases exceeds the single sheet maximum, multiple sheets are created in the workbook.
- A maximum of 256 variables can be saved to Excel 2.1 Excel 95, and Excel 97; a maximum of 16,000 variables can be saved to Excel 2007.
- A maximum of 65,535 cases and 32 variables can be saved to dBASE II, a maximum of one billion cases (subject to disk space availability) and 128 variables can be saved to dBASE III, and a maximum of one billion cases (subject to disk space availability) and 255 variables can be saved to dBASE IV.
- A maximum of 2,047 variables can be saved to Stata 5–6 and Intercooled Stata 7–8. A maximum of 32,767 variables can be saved to Stata SE 7–8.
- A maximum of 32,767 variables can be saved to SAS 6-8.

## Operations

- The active dataset remains available after `SAVE TRANSLATE` is executed.
- If the active dataset contains more variables than the file can receive, `SAVE TRANSLATE` writes the maximum number of variables that the file can receive.

## Spreadsheets

Variables in the active dataset become columns, and cases become rows in the spreadsheet file.

- If you specify `FIELDNAMES`, variable names become the first row.
- To save value labels instead of data values, use the `CELLS` subcommand. [For more information, see the topic CELLS Subcommand on p. 1833.](#)
- System-missing values are translated to `#NULL!` in Excel files.

IBM® SPSS® Statistics formats are translated as follows:

SPSS Statistics	1-2-3/Symphony	Excel
Number	Fixed	0.00;#,##0.00;...
COMMA	Comma	0.00;#,##0.00;...
DOLLAR	Currency	\$#,##0_);...
DATE	Date	d-mmm-yy
TIME	Time	hh:mm:ss
String	Label	General

## dBASE

Variables in the active dataset become fields, and cases become records in the database file.

- Characters that are allowed in variable names, but not in dBASE field names, are translated to colons in dBASE II and underscores in dBASE III and dBASE IV.

- Numeric variables containing the system-missing value are translated to \*\*\*\* in dBASE III and dBASE IV and 0 in dBASE II.
- The width and precision of translated numeric variables are taken from the print format—the total number of characters for the number is taken from the width of the print format, and the number of digits to the right of the decimal point is taken from the decimals in the print format. To adjust the width and precision, use the `PRINT FORMATS` command before using `SAVE TRANSLATE`. Values that cannot be converted to the given width and precision are converted to missing values.
- A maximum of 65,535 cases and 32 variables can be saved to dBASE II, a maximum of one billion cases (subject to disk space availability) and 128 variables can be saved to dBASE III, and a maximum of one billion cases (subject to disk space availability) and 255 variables can be saved to dBASE IV.

Variable formats are translated to dBASE formats as follows:

IBM® SPSS® Statistics	dBASE
Number	Numeric
String	Character
Dollar	Numeric
Comma	Numeric

### ***Comma-Delimited (CSV) Text Files***

Variables in the active dataset become columns, and cases become rows in the text file.

- If you specify `FIELDNAMES`, variable names are written in the first row.
- To save value labels instead of data values, use the `CELLS` subcommand. [For more information, see the topic `CELLS` Subcommand on p. 1833.](#)
- Use the `TEXTOPTIONS` subcommand to override the default delimiter, qualifier, decimal indicator, or data value formatting. [For more information, see the topic `TEXTOPTIONS` Subcommand on p. 1833.](#)
- System-missing values are translated to a blank space in text files.
- In Unicode mode, CSV files are saved in UTF-8 format. In code page mode, CSV files are saved in the code page of the current locale. [For more information on encoding, see `UNICODE` Subcommand.](#)

*Note:* Tab characters embedded in string values are preserved as tab characters in the tab-delimited file. No distinction is made between tab characters embedded in values and tab characters that separate values.

### ***Tab-Delimited Text Files***

Variables in the active dataset become columns, and cases become rows in the text file.

- If you specify `FIELDNAMES`, variable names are written to the first row.



- To save value labels instead of data values, use the `CELLS` subcommand. For more information, see the topic `CELLS Subcommand` on p. 1833.
- Use the `TEXTOPTIONS` subcommand to override the default decimal indicator or data value formatting. For more information, see the topic `TEXTOPTIONS Subcommand` on p. 1833.
- System-missing values are translated to a blank space in text files.
- In Unicode mode, tab-delimited files are saved in UTF-8 format. In code page mode, tab-delimited files are saved in the code page of the current locale. For more information on encoding, see `UNICODE Subcommand`.

## SAS Files

Data can be saved in numerous SAS data file formats. A SAS transport file is a sequential file that is written in SAS transport format and can be read by SAS with the `XPORT` engine and `PROC COPY` or the `DATA` step.

- Certain characters that are allowed in IBM® SPSS® Statistics variable names are not valid in SAS, such as `@`, `#`, and `$`. These illegal characters are replaced with an underscore when the data are exported.
- SPSS Statistics variable names that contain multibyte characters (e.g., Japanese or Chinese characters) are converted to variable names of the general form `Vnnn`, where `nnn` is an integer value.
- SPSS Statistics variable labels containing more than 40 characters are truncated when exported to a SAS v6 file.
- Where they exist, SPSS Statistics variable labels are mapped to the SAS variable labels. If no variable label exists in the SPSS Statistics data, the variable name is mapped to the SAS variable label.
- SAS allows only one value for missing, whereas SPSS Statistics allows the definition of numerous missing values. As a result, all missing values in SPSS Statistics are mapped to a single missing value in the SAS file.

The following table shows the variable type matching between the original data in SPSS Statistics and the exported data in SAS.

SPSS Statistics Variable Type	SAS Variable Type	SAS Data Format
Numeric	Numeric	12
Comma	Numeric	12
Dot	Numeric	12
Scientific Notation	Numeric	12
Date	Numeric	Date
Date (Time)	Numeric	Time
Date (Date-Time)	Numeric	DateTime
Dollar	Numeric	12
Custom Currency	Numeric	12
String	Character	\$8

## Stata Files

- Data can be written in Stata 5–8 format and in both Intercooled and SE format (versions 7 and 8 only).
- Data files that are saved in Stata 5 format can be read by Stata 4.
- The first 80 bytes of variable labels are saved as Stata variable labels.
- For numeric variables, the first 80 bytes of value labels are saved as Stata value labels. Value labels are dropped for string variables, non-integer numeric values, and numeric values greater than an absolute value of 2,147,483,647.
- For versions 7 and 8, the first 32 bytes of variable names in case-sensitive form are saved as Stata variable names. For earlier versions, the first eight bytes of variable names are saved as Stata variable names. Any characters other than letters, numbers, and underscores are converted to underscores.
- IBM® SPSS® Statistics variable names that contain multibyte characters (for example, Japanese or Chinese characters) are converted to variable names of the general form *Vnnn*, where *nnn* is an integer value.
- For versions 5–6 and Intercooled versions 7–8, the first 80 bytes of string values are saved. For Stata SE 7–8, the first 244 bytes of string values are saved.
- For versions 5–6 and Intercooled versions 7–8, only the first 2,047 variables are saved. For Stata SE 7–8, only the first 32,767 variables are saved.

SPSS Statistics Variable Type	Stata Variable Type	Stata Data Format
Numeric	Numeric	g
Comma	Numeric	g
Dot	Numeric	g
Scientific Notation	Numeric	g
Date*, Datetime	Numeric	D_m_Y
Time, DTime	Numeric	g (number of seconds)
Wkday	Numeric	g (1–7)
Month	Numeric	g (1–12)
Dollar	Numeric	g
Custom Currency	Numeric	g
String	String	s

\*Date, Adate, Edate, SDate, Jdate, Qyr, Moyr, Wkyr

## SPSS/PC+ System Files

Variables are saved as they are defined. The resulting file is given the extension *.sys* if no extension is explicitly specified. The dictionary is saved so that labels, formats, missing value specifications, and other dictionary information are preserved. This format is available only on Windows operating systems.

## ODBC Database Sources

The following rules apply when writing to a database with `TYPE=ODBC`:

- If you specify a table name that does not exist in the database, a new table with that name is created.
- If any case cannot be stored in the database for any reason, an error is returned. Therefore, either all cases are stored or no cases are stored.
- At insert time, a check is performed to see whether the value that is being stored in a column is likely to cause an overflow. If so, the user is warned about the overflow and is informed that a *SYSMIS* is stored instead.
- If any variable names in the active dataset contain characters that are not allowed by the database, they are replaced by an underscore. If this process causes a duplicate variable name, a new variable name is generated.

By default, IBM® SPSS® Statistics variable formats are mapped to database field types based on the following general scheme. Actual database field types may vary, depending on the database.

SPSS Statistics Variable Format	Database Field Type
Numeric	Float or Double
Comma	Float or Double
Dot	Float or Double
Scientific Notation	Float or Double
Date	Date or Datetime or Timestamp
Datetime	Datetime or Timestamp
Time, DTime	Float or Double (number of seconds)
Wkday	Integer (1–7)
Month	Integer (1–12)
Dollar	Float or Double
Custom Currency	Float or Double
String	Char or Varchar

- For database Datetime/Timestamp formats that don't allow fractional seconds, fraction seconds in SPSS Statistics Datetime values are truncated.
- The default width of Char database data types for SPSS Statistics string variables is the defined width of the string variable. If the defined width of the string exceeds the maximum width allowed by the database, the values are truncated to the maximum width, and a warning is issued.
- You can override the default settings by using the SQL subcommand to create or update tables. [For more information, see the topic SQL Subcommand on p. 1836.](#)

## TYPE Subcommand

`TYPE` indicates the format of the resulting file.

- `TYPE` can be omitted for spreadsheet files if the file extension that is named on `OUTFILE` is the default for the type of file that you are saving.

---

*SAVE TRANSLATE*

- `TYPE` with the keyword `DB2`, `DB3`, or `DB4` is required for translating to dBASE files.
- `TYPE` takes precedence over the file extension.

<b>ODBC</b>	<i>Database accessed via ODBC.</i>
<b>XLS</b>	<i>Excel.</i>
<b>CSV</b>	<i>Comma-delimited text data files.</i>
<b>TAB</b>	<i>Tab-delimited text data files.</i>
<b>SAS</b>	<i>SAS data files and SAS transport files.</i>
<b>STATA</b>	<i>Stata data files.</i>
<b>DB2</b>	<i>dBASE II.</i>
<b>DB3</b>	<i>dBASE III or dBASE III PLUS.</i>
<b>DB4</b>	<i>dBASE IV.</i>
<b>WK1</b>	<i>1-2-3 Release 2.0.</i>
<b>WKS</b>	<i>1-2-3 Release 1.4.</i>
<b>SYM</b>	<i>Symphony releases.</i>
<b>SLK</b>	<i>Multiplan (symbolic format).</i>
<b>PC</b>	<i>SPSS/PC+ system files. This format is available only on Windows operating systems.</i>

**Example**

```
SAVE TRANSLATE OUTFILE='PROJECT.XLS' /TYPE=XLS (/VERSION=8) .
```

- `SAVE TRANSLATE` translates the active dataset into the Excel workbook file named *PROJECT.XLS*.

**VERSION Subcommand**

`VERSION` specifies the file version for multiversion applications. For example, this subcommand is necessary to differentiate between different versions of Excel.

- `VERSION` is required for all applications listed below, except Excel.
- If `VERSION` is not specified with `TYPE=XLS`, the Excel file is saved in version 2.1 format.

<b>Version</b>	<b>Application</b>
1	Symphony Release 1.0
2	Excel 2.1, Symphony Release 2.0, dBASE II
3	dBASE III or dBASE III PLUS
4	dBASE IV
5	Excel 95 Workbook, Stata 4–5
6	SAS v6, Stata 6
7	SAS v7–8, Stata 7
8	Excel 97–2003 Workbook, Stata 8
9	SAS 9
12	Excel 2007 Workbook
X	SAS transport file

**Example**

```
SAVE TRANSLATE OUTFILE='STAFF.XLS' /TYPE=XLS /VERSION=8.
```

- SAVE TRANSLATE creates an Excel workbook file in the version 97 format.

**ENCODING Subcommand**

The ENCODING subcommand specifies the character encoding for SAS and Stata data files.

- The ENCODING subcommand is only valid with TYPE=SAS and TYPE=STATA.
- The subcommand name is preceded by a slash and followed by an optional equals sign and a quoted value.
- The quoted value can be “Locale” or “System” or one of the values in the Encodings column in See [Table 220-1](#) on p. 1831.. For SAS 9 files, the value can also be “UTF-8”.
- The default is “Locale”, which is the current IBM® SPSS® Statistics locale, except for SAS 9 files saved in Unicode mode. [For more information, see the topic LOCALE Subcommand on p. 1876.](#)
- In Unicode mode with TYPE=SAS and VERSION=9, the default is “UTF-8”. [For more information, see the topic UNICODE Subcommand on p. 1878.](#)
- The ENCODING setting also applies to the value labels file specified on the optional VALFILE subcommand for TYPE=SAS.

**Example**

```
SAVE TRANSLATE
  /OUTFILE=' /data/sasdata.sas7bdat '
  /VALFILE=' /data/saslabels.sas '
  /TYPE=SAS /VERSION=7 /PLATFORM=WINDOWS
  /ENCODING='Windows-1252'.
```

Table 220-1  
Character Encoding

Character Set	Encoding
SPSS Statistics Locale	Locale
Operating System Locale	System
Western European	
Western	ISO-8859-1
Western	ISO-8859-15
Western	IBM-850
Western	Windows-1252
Celtic	ISO-8859-14
Greek	ISO-8859-7
Greek	Windows-1253
Nordic	ISO-8859-10
Eastern European	
Baltic	Windows-1257
Central European	IBM-852

Central European	ISO-8859-2
Cyrillic	IBM-855
Cyrillic	ISO-8859-5
Cyrillic	Windows-1251
Cyrillic/Russian	CP-866
Eastern Asian	
Chinese Simplified	GBK
Chinese Simplified	ISO-2022-CN
Chinese Traditional	Big5
Chinese Traditional	EUC-TW
Japanese	EUC-JP
Japanese	ISO-2022-JP
Japanese	Shift-JIS
Korean	EUC-KR
Southeastern and Southwestern Asian	
Thai	Windows-874
Turkish	IBM-857
Turkish	ISO-8859-9
Middle Eastern	
Arabic	Windows-1256
Arabic	IBM-864
Hebrew	ISO-8859-8
Hebrew	Windows-1255
Hebrew	IBM-862

## ***OUTFILE Subcommand***

OUTFILE assigns a name to the file to be saved. The only specification is the name of the file. On some operating systems, file specifications should be enclosed in quotes.

### ***Example***

```
SAVE TRANSLATE OUTFILE='STAFF.DBF' /TYPE=DB3.
```

- SAVE TRANSLATE creates a dBASE III file called *STAFF.DBF*. The TYPE subcommand is required to specify the type of dBASE file to save.

## ***FIELDNAMES Subcommand***

FIELDNAMES writes variable names to the first row in spreadsheets and tab-delimited and comma-delimited text files.

### ***Example***

```
SAVE TRANSLATE OUTFILE='/data/csvfile.csv' /TYPE=CSV /FIELDNAMES.
```

## **CELLS Subcommand**

For TYPE=XLS (/VERSION=8), TYPE=TAB, and TYPE=CSV, the CELLS subcommand specifies whether data values or value labels are saved.

**CELLS=VALUES.** *Saves data values.* This is the default.

**CELLS=LABELS.** *Saves value labels instead of values.* For values with no corresponding value label, the value is saved.

### **Example**

```
SAVE TRANSLATE OUTFILE='/data/sales.xls'  
  /TYPE=XLS /VERSION=8  
  /CELLS=LABELS.
```

## **TEXTOPTIONS Subcommand**

For TYPE=CSV, the TEXTOPTIONS subcommand controls the delimiter and qualifier. For TYPE=CSV and TYPE=TAB, it controls the decimal indicator and the data value formatting. For all other types, this subcommand is ignored.

**DELIMITER="char".** *Specifies the delimiter (separator) between values.* The value must be a single character (single byte), enclosed in quotes (single or double quotes). The default is either a comma or semicolon, based on the decimal indicator. If the decimal indicator is a period, then the default separator is a comma. If the decimal indicator is a comma, then the default separator is a semicolon. The value cannot be the same as the QUALIFIER or DECIMAL value. This setting only applies for TYPE=CSV. For other types, it is ignored.

**QUALIFIER="char".** *Specifies the qualifier to use to enclose values that contain the delimiter character.* The default is a double quote. The value must be a single character (single byte), enclosed in quotes (single or double quotes). Values that contain the qualifier character will also be enclosed by the qualifier and qualifiers within the value will be doubled. The value cannot be the same as the DELIMITER or DECIMAL value. This setting only applies for TYPE=CSV. For other types, it is ignored.

**DECIMAL=DOT|COMMA.** *Specifies the value to use as the decimal indicator for numeric values.* The default is the current IBM® SPSS® Statistics decimal indicator. This setting has no effect on the decimal indicator used with FORMAT=VARIABLE for comma, dollar, custom currency, and dot formats. (Comma, dollar, and custom currency always use the period as the decimal indicator; dot always uses the comma as the decimal indicator.) The value cannot be the same as the DELIMITER or QUALIFIER value. (If DECIMAL=DOT, the default delimiter is a comma; if DECIMAL=COMMA, the default delimiter is a semicolon.)

- **DOT.** *Use a period (.) as the decimal indicator.*
- **COMMA.** *Use a comma (,) as the decimal indicator.*

The default decimal indicator is the current SPSS Statistics decimal indicator. The current SPSS Statistics decimal indicator can be set with `SET LOCALE` (which uses the OS locale decimal indicator) or `SET DECIMAL` and can be shown with `SHOW DECIMAL`. The `TEXTOPTIONS DECIMAL` setting overrides the default decimal indicator in the current `SAVE TRANSLATE` command but does not change the default/current SPSS Statistics decimal indicator, and it has no effect on the determination of the default delimiter for `TYPE=CSV`.

**FORMAT=PLAIN|VARIABLE.** *Specifies the data formats to use when writing out data values.*

- **PLAIN.** *Remove all “extraneous” formatting from numeric values.* This is the default. For example, a dollar format value of \$12,345.67 is written simply as 12345.67 (or 12345,67, depending on the decimal indicator). The number of decimal positions for each value is the number of decimal positions necessary to preserve the entire value (and different values of the same variable may have a different number of decimal positions). Dates are written in the general format mm/dd/yyyy. Times are written in the general format hh:mm:ss.
- **VARIABLE.** *Use the print format for each variable to write data values.* For example, a value of 12345.67 with a print format of `DOLLAR10.2` is written as \$12,345.67 (and is qualified with double quotes or the user-specified qualifier if a comma is the value delimiter).

Print formats can be set with the `PRINT FORMATS` or `FORMATS` commands and displayed with `DISPLAY DICTIONARY`.

## ***EDITION Subcommand***

The `EDITION` subcommand specifies the edition for Stata files. `EDITION` only applies to Stata version 7 or later.

- |                    |   |
|--------------------|---|
| <b>INTERCOOLED</b> | Saves the data file in Stata Intercooled format. This setting is the default. Only the first 2,047 variables are saved, and only the first 80 bytes of string values are saved. |
| <b>SE</b>          | Saves the data file in Stata SE format. The first 32,767 variables are saved, and the first 244 bytes of string values are saved.   |

### ***Example***

```
SAVE TRANSLATE OUTFILE='//data/newdata.dta' /TYPE=STATA
/VERSION=8 /EDITION=SE.
```

This subcommand is ignored for any file types other than Stata version 7 or later.

## ***PLATFORM Subcommand***

The `PLATFORM` subcommand is required for all SAS file types, with the exception of SAS transport files. Enter the keyword corresponding to the platform for which the target SAS file is intended. There is no default value for this command. Choose from the following keywords: `WINDOWS`, `ALPHA`, `UNIX`.

### ***Example***

```
SAVE TRANSLATE OUTFILE='STAFF.SD7' /TYPE=SAS /VERSION=7
```



```
/PLATFORM=WINDOWS /VALFILE='LABELS.SAS'.
```

## VALFILE Subcommand

The VALFILE subcommand, which is available only for SAS file formats, excluding the SAS transport file format, saves value labels to a SAS syntax file. This syntax file is used in conjunction with the saved SAS data file to re-create value labels in SAS.

- The syntax file has an *.sas* file extension.

### Example

```
SAVE TRANSLATE OUTFILE='STAFF.SD7' /TYPE=SAS /VERSION=7
/PLATFORM=WINDOWS /VALFILE='LABELS.SAS'.
```

- SAVE TRANSLATE saves the current data as a SAS v7 file and creates a SAS syntax file named *LABELS.SAS*, which contains the value labels for the data.

## ODBC Database Subcommands

The CONNECT, ENCRYPTED, TABLE, SQL, and APPEND subcommands are used with TYPE=ODBC.

### CONNECT Subcommand

CONNECT identifies the database name and other parameters for TYPE=ODBC.

- The database must already exist. You cannot create a new database with SAVE TRANSLATE (although you can create new tables in an existing database).
- If you are unsure what the connect string should be, you can use the Export to Database Wizard to generate SAVE TRANSLATE command syntax. If the connection string is specified on multiple lines, each line must be enclosed in quotes.

### Example

```
SAVE TRANSLATE
  /TYPE=ODBC
  /CONNECT
  'DSN=MS Access Database;DBQ=C:\examples\data\dm_demo.mdb;'
  'DriverId=25;FIL=MS Access;MaxBufferSize=2048;PageTimeout=5;'
  /TABLE='mytable'.
```

### ENCRYPTED and UNENCRYPTED Subcommands

The ENCRYPTED subcommand indicates that the password in the CONNECT string is encrypted. The default is UNENCRYPTED. (Both the Database Wizard and Export to Database Wizard generate encrypted passwords.)

The ENCRYPTED and UNENCRYPTED subcommands have no additional specifications.

## TABLE Subcommand

TABLE identifies the table name for TYPE=ODBC.

- Table names should be enclosed in quotes.
- You can replace an existing table or append new records to an existing table.
- You can create a new table in the database by specifying a table name that does not currently exist in the database.

### Example

```
SAVE TRANSLATE
  /TYPE=ODBC
  /CONNECT
  'DSN=MS Access Database;DBQ=C:\examples\data\dm_demo.mdb;'
  'DriverId=25;FIL=MS Access;MaxBufferSize=2048;PageTimeout=5;'
  /TABLE='mytable'.
```

## SQL Subcommand

The SQL subcommand provides the ability to issue any SQL directives that are needed in the target database. It can be used, for example, to define joins or alter table properties in the database to include new columns or modify column properties.

- Each SQL statement must be enclosed in quotes.
- You can use multiple lines for a single SQL statement by using multiple quoted strings connected with plus signs (the standard string concatenation symbol).
- Each quoted line cannot exceed 256 characters.
- Multiple SQL statements can be included by using multiple SQL subcommands.
- Table and field specifications in SQL statements refer to tables and fields available in the database, not datasets and variables available in the IBM® SPSS® Statistics session (although in many cases the names may be the same).
- Regardless of the position of the SQL subcommand, the SQL statements are executed last, after all other actions executed by the SAVE TRANSLATE command.

### Example: Adding New Columns to an Existing Table

```
SAVE TRANSLATE /TYPE=ODBC
  /CONNECT=
  'DSN=MS Access Database;DBQ=C:\examples\data\dm_demo.mdb;'
  'DriverId=25;FIL=MS Access;MaxBufferSize=2048;PageTimeout=5;'
  /TABLE = 'NewColumn'
  /KEEP ID income_score
  /REPLACE
  /SQL='ALTER TABLE CombinedTable ADD COLUMN income_score REAL'
  /SQL='UPDATE CombinedTable INNER JOIN NewColumn ON ' +
  'CombinedTable.ID=NewColumn.ID SET ' +
  'CombinedTable.income_score=NewColumn.income_score'.
```

- The TABLE, KEEP, and REPLACE subcommands create or replace a table named *NewColumn* that contains two variables: a key variable (*ID*) and a computed score (*income\_score*).

- The first SQL subcommand, specified on a single line, adds a column to an existing table that will contain values of the computed variable *income\_score*. At this point, all we have done is create an empty column in the existing database table, and the fact that both database tables and the active dataset use the same name for that column is merely a convenience for simplicity and clarity.
- The second SQL subcommand, specified on multiple lines with the quoted strings concatenated with plus signs, adds the *income\_score* values from the new table to the existing table, matching rows (cases) based on the value of the key variable *ID*.

The end result is that an existing table is modified to include a new column containing the values of the computed variable.

### **Example: Specifying Data Types for a New Table**

```
SAVE TRANSLATE /TYPE=ODBC
/CONNECT='DSN=MS Access Database;DBQ=c:\temp\temp.mdb;DriverId=25;FIL=MS'+
' Access;MaxBufferSize=2048;PageTimeout=5;'
/TABLE=tempTable /REPLACE
/SQL='CREATE TABLE NewTable(numVar double, intVar integer, dollarVar currency)'
/SQL='INSERT INTO NewTable(numVar, intVar, dollarVar) SELECT * FROM tempTable'
/SQL='DROP TABLE tempTable'.
```

- The TABLE subcommand creates a new table that contains variables in the active dataset with the default database data types. In this example, the original variables have SPSS Statistics variable formats of F8.2, F4.0, and Dollar10.2 respectively, but the default database type for all three is double.
- The first SQL subcommand creates another new table with data types explicitly defined for each field. At this point, this new table contains no data.
- The second SQL subcommand inserts the data from the *tempTable* into *NewTable*. This does not affect the data types previously defined for *NewTable*, so *intVar* will have a data type of integer and *dollarVar* will have a data type of currency.
- The last SQL subcommand deletes *tempTable*, since it is no longer needed.

## **APPEND Subcommand**

APPEND appends rows (cases) to an existing database table after type and variable name validations. There must be a matching column in the table for each IBM® SPSS® Statistics variable. If a column that can correctly store a variable is not found, a failure is returned. If the table contains more columns than the number of SPSS Statistics variables, the command still stores the data in the table. The variable names and column names must match exactly. A variable can be stored in a column as long as the column type is a type that can store values of the variable type. A column of any numeric type (short integer, integer, float, double, etc.) is valid for a numeric variable, and a column of any character type is valid for a string variable.

- APPEND is valid only for TYPE=ODBC.
- APPEND and REPLACE are mutually exclusive.

*Note:* APPEND can only add rows to a table, not columns. If you want to add columns (fields) to a database table, see the [SQL subcommand](#).

## **REPLACE Subcommand**

REPLACE gives permission to overwrite an existing file (or database table) of the same name. REPLACE takes no further specifications.

- SAVE TRANSLATE will not overwrite an existing file without an explicit REPLACE subcommand. The default behavior is to not overwrite.
- For database tables (TYPE=ODBC), APPEND and REPLACE are mutually exclusive.

*Note:* For database tables, REPLACE destroys an existing table and replaces it with the data written by the SAVE TRANSLATE command, which may lead to loss of database-specific information, such as the designation of key variables (fields) and data formats. If you want to update an existing table with new values in some existing rows or by adding additional fields to the table, see the [SQL subcommand](#).

## **UNSELECTED Subcommand**

UNSELECTED determines whether cases that were excluded on a previous FILTER or USE command are to be retained or deleted. The default is RETAIN. The UNSELECTED subcommand has no effect when the active dataset does not contain unselected cases.

- RETAIN**     *Retain the unselected cases.* All cases in the active dataset are saved. This setting is the default when UNSELECTED is specified by itself.
- DELETE**     *Delete the unselected cases.* Only cases that meet the FILTER or USE criteria are saved.

## **DROP and KEEP Subcommands**

Use DROP or KEEP to include only a subset of variables in the resulting file. DROP specifies a set of variables to exclude. KEEP specifies a set of variables to retain. Variables that are not specified on KEEP are dropped.

- Specify a list of variable, column, or field names separated by commas or spaces.
- KEEP does not affect the order of variables in the resulting file. Variables are kept in their original order.
- Specifying a variable that is not in the active dataset or that has been dropped because of a previous DROP or KEEP subcommand results in an error, and the SAVE command is not executed.

### **Example**

```
SAVE TRANSLATE OUTFILE='ADDRESS.DBF' /TYPE=DB4 /DROP=PHONENO, ENTRY.
```

- SAVE TRANSLATE creates a dBASE IV file named ADDRESS.DBF, dropping the variables PHONENO and ENTRY.

## **RENAME Subcommand**

RENAME changes the names of variables as they are copied into the resulting file.

- The specification on `RENAME` is a list of old variable names followed by an equals sign and a list of new variable names. The same number of variables must be specified on both lists. The keyword `TO` can be used in the first list to refer to consecutive variables in the active dataset and can be in the second list to generate new variable names. The entire specification must be enclosed in parentheses.
- Alternatively, you can specify each old variable name individually, followed by an equals sign and the new variable name. Multiple sets of variable specifications are allowed. The parentheses around each set of specifications are optional.
- New names cannot exceed 64 bytes. Characters not allowed in IBM® SPSS® Statistics variable names can be used in new names for the target file, but if the name contains special characters (e.g., spaces, commas, slashes, plus signs) the name must be enclosed in quotes.
- `RENAME` does not affect the active dataset. However, if `RENAME` precedes `DROP` or `KEEP`, variables must be referred to by their new names on `DROP` or `KEEP`.
- Old variable names do not need to be specified according to their order in the active dataset.
- Name changes take place in one operation. Therefore, variable names can be exchanged between two variables.
- Multiple `RENAME` subcommands are allowed.

### Examples

```
SAVE TRANSLATE OUTFILE='STAFF.SYM' VERSION=2 /FIELDNAMES
  /RENAME AGE=AGE88 JOBCAT=JOBCAT88.
```

- `RENAME` renames the variable `AGE` to `AGE88` and renames `JOBCAT` to `JOBCAT88` before they are copied to the first row of the spreadsheet.

```
SAVE TRANSLATE OUTFILE='STAFF.SYM' VERSION=2 /FIELDNAMES
  /RENAME (AGE JOBCAT=AGE88 JOBCAT88).
```

- The name changes are identical to the changes in the previous example: `AGE` is renamed to `AGE88`, and `JOBCAT` is renamed to `JOBCAT88`. The parentheses are required with this method.

## MISSING Subcommand

The `MISSING` subcommand controls the treatment of user-missing values. By default, user-missing values are treated as regular valid values in the target file.

- IGNORE**     *Treat user-missing values as regular valid values.* This setting is the default.
- RECODE**    *Recode numeric user-missing values to system-missing, and recode string user-missing values to blank spaces.*

## MAP Subcommand

`MAP` displays a list of the variables in the resulting file and their corresponding names in the active dataset.

- The only specification is the keyword `MAP`.

---

*SAVE TRANSLATE*

- Multiple `MAP` subcommands are allowed. Each `MAP` subcommand maps the results of subcommands that precede it but not the results of subcommands that follow it.

***Example***

```
GET FILE=HUBEMPL.  
SAVE TRANSLATE OUTFILE='STAFF.SYM' /VERSION=2 /FIELDNAMES  
  /RENAME=(AGE=AGE88) (JOB CAT=JOB CAT88) .
```

- `MAP` is specified to confirm that the variable `AGE` is renamed to `AGE88` and `JOB CAT` is renamed to `JOB CAT88`.

# SCRIPT

```
SCRIPT 'filename' [(quoted string)]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- Scripts run from the `SCRIPT` command now run synchronously with the command syntax stream.

Release 17.0

- Ability to run Python scripts introduced.

## **Overview**

`SCRIPT` runs a script to customize the program or automate regularly performed tasks. You can run a Basic script or a Python script.

### **Basic Specification**

The basic specification is keyword `SCRIPT` with a filename. The filename is required. The optional quoted string, enclosed in parentheses, can be passed to the script.

### **Operations**

- `SCRIPT` runs the specified script, and works in both local and distributed mode. The effect is the same as opening the script file in the script editor (Basic or Python) and running it from there.

### **Limitations**

- Python scripts run from the `SCRIPT` command cannot execute command syntax.
- The `SCRIPT` command is not supported in the IBM® SPSS® Statistics Batch Facility (available with SPSS Statistics Server).

## **Running Basic Scripts That Contain Syntax Commands**

If a Basic script that is run via the `SCRIPT` command contains syntax commands, those commands must be run asynchronously. To run commands asynchronously, set the `bSync` parameter of the `ExecuteCommands` method to `False`, as in:

```
Dim objSpssApp As spsswinLib.Application16  
Dim strCommands As String
```

```
Set objSpssApp=CreateObject("SPSS.Application16")  
  
' Construct and execute syntax commands:  
  
strCommands = "GET FILE = '/data/bank.sav'. " & vbCr  
strCommands = strCommands & "Display Dictionary."  
objSpssApp.ExecuteCommands strCommands, False
```



# SEASON

SEASON is available in the Forecasting option.

```
SEASON VARIABLES= series names
```

```
  [/MODEL={MULTIPLICATIVE**}]  
           {ADDITIVE      }
```

```
  [/MA={EQUAL    }]  
       {CENTERED}
```

```
  [/PERIOD=n]
```

```
  [/APPLY [= 'model name']]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
SEASON VARIABLES = VARX  
  /MODEL=ADDITIVE.
```

## Overview

SEASON estimates multiplicative or additive seasonal factors for time series by using any specified periodicity. SEASON is an implementation of the Census Method I, otherwise known as the ratio-to-moving-average method. See (Makridakis, Wheelwright, and McGee, 1983) and (McLaughlin, 1984).

### Options

**Model Specification.** You can specify either a multiplicative or additive model on the MODEL subcommand. You can specify the periodicity of the series on the PERIOD subcommand.

**Computation Method.** Two methods of computing moving averages are available on the MA subcommand for handling series with even periodicities.

**Statistical Output.** Specify TSET PRINT=BRIEF to display only the initial seasonal factor estimates. TSET PRINT=DETAILED produces the same output as the default.

**New Variables.** To evaluate the displayed averages, ratios, factors, adjusted series, trend-cycle, and error components without creating new variables, specify TSET NEWVAR=NONE prior to SEASON. This can result in faster processing time. To add new variables without erasing the values of previous Forecasting-generated variables, specify TSET NEWVAR=ALL. This saves all new variables that are generated during the current session to the active dataset and may require extra processing time. When the default (TSET NEWVAR=CURRENT) is in effect, only variables from the current procedure are saved to the active dataset, and the suffix #n is used to distinguish

variables that are generated by different series. `TSET MXNEWVAR` specifies the maximum number of new variables that can be generated by a procedure. The default is 60. The order in which new variables are added to the active dataset's dictionary is *ERR*, *SAS*, *SAF*, and *STC*.

### **Basic Specification**

The basic specification is one or more series names.

- By default, `SEASON` uses a multiplicative model to compute and display moving averages, ratios, seasonal factors, the seasonally adjusted series, the smoothed trend-cycle components, and the irregular (error) component for each specified series (variable). The default periodicity is the periodicity that is established with `TSET PERIOD` or `DATE`.

- Unless the default on `TSET NEWVAR` is changed prior to the procedure, `SEASON` creates four new variables, with the following prefixes, for each specified series (these variables are automatically named, labeled, and added to the active dataset):

**SAF.** *Seasonal adjustment factors.* These values indicate the effect of each period on the level of the series.

**SAS.** *Seasonally adjusted series.* These values are the values that are obtained after removing the seasonal variation of a series.

**STC.** *Smoothed trend-cycle components.* These values show the trend and cyclical behavior that are present in the series.

**ERR.** *Residual or "error" values.* These values are the values that remain after the seasonal, trend, and cycle components have been removed from the series.

### **Subcommand Order**

- Subcommands can be specified in any order.

### **Syntax Rules**

- `VARIABLES` can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each one is executed.

### **Operations**

- The endpoints of the moving averages and ratios are displayed as system-missing in the output.
- Missing values are not allowed anywhere in the series. (You can use the `RMV` command to replace missing values, use `TSET MISSING=INCLUDE` to include user-missing values, and use `USE` to ignore missing observations at the beginning or end of a series. See `RMV` and `USE` for more information.)

If `SPLIT FILE` processing is on, `SEASON` ignores `USE` settings.

### **Limitations**

- A maximum of one `VARIABLES` subcommand is allowed. There is no limit on the number of series that are named on the list.

## **VARIABLES Subcommand**

VARIABLES specifies the series names and is the only required subcommand.

- Each specified series must contain at least four full seasons of data.

## **MODEL Subcommand**

MODEL specifies whether the seasonal decomposition model is multiplicative or additive.

- The specification on MODEL is the keyword MULTIPLICATIVE or ADDITIVE.
- If more than one keyword is specified, only the first keyword is used.
- MULTIPLICATIVE is the default if the MODEL subcommand is not specified or if MODEL is specified without any keywords.

### **Example**

```
SEASON VARIABLES = VARX  
/MODEL=ADDITIVE.
```

- This example uses an additive model for the seasonal decomposition of *VARX*.

## **MA Subcommand**

MA specifies how to treat an even-periodicity series when computing moving averages.

- MA should be specified only when the periodicity is even. When periodicity is odd, the EQUAL method is always used.
- For even-periodicity series, the keyword EQUAL or CENTERED can be specified. CENTERED is the default.
- EQUAL calculates moving averages with a span (number of terms) equal to the periodicity and all points weighted equally.
- CENTERED calculates moving averages with a span (number of terms) equal to the periodicity plus 1 and endpoints weighted by 0.5.
- The periodicity is specified on the PERIOD subcommand.

### **Example**

```
SEASON VARIABLES = VARY  
/MA=CENTERED  
/PERIOD=12.
```

- In this example, moving averages are computed with spans of 13 terms and endpoints weighted by 0.5.

## **PERIOD Subcommand**

PERIOD indicates the size of the period.

- The specification on `PERIOD` indicates how many observations are in one period or season and can be any positive integer.
- If `PERIOD` is not specified, the periodicity that is established on `TSET PERIOD` is in effect. If `TSET PERIOD` is not specified, the periodicity that is established on the `DATE` command is used. If periodicity was not established anywhere, the `SEASON` command will not be executed.

**Example**

```
SEASON VARIABLES = SALES
  /PERIOD=12.
```

- In this example, a periodicity of 12 is specified for *SALES*.

**APPLY Subcommand**

`APPLY` allows you to use a previously defined `SEASON` model without having to repeat the specifications.

- The only specification on `APPLY` is the name of a previous model in quotation marks. If a model name is not specified, the model that was specified on the previous `SEASON` command is used. Model names are either the default `MOD_n` names or the names that are assigned on the `MODEL NAME` command.
- To change one or more model specifications, specify the subcommands of only those portions that you want to change after the `APPLY` subcommand.
- If no series are specified on the command, the series that were originally specified with the model that is being reapplied are used.
- To change the series used with the model, enter new series names before or after the `APPLY` subcommand. If a series name is specified before `APPLY`, the slash before the subcommand is required.

**Example**

```
SEASON VARIABLES = X1
  /MODEL=ADDITIVE.
SEASON VARIABLES = Z1
  /APPLY.
```

- The first command specifies an additive model for the seasonal decomposition of *X1*.
- The second command applies the same type of model to series *Z1*.

**Example**

```
SEASON X1 Y1 Z1
  /MODEL=MULTIPLICATIVE.
SEASON APPLY
  /MODEL=ADDITIVE.
```

- The first command specifies a multiplicative model for the seasonal decomposition of *X1*, *Y1*, and *Z1*.
- The second command applies an additive model to the same three variables.

## **References**

Makridakis, S., S. C. Wheelwright, and V. E. McGee. 1983. *Forecasting: Methods and applications*. New York: John Wiley and Sons.

McLaughlin, R. L. 1984. *Forecasting techniques for decision making*. Rockville, Md.: Control Data Management Institute.

# SELECT IF

```
SELECT IF [(logical expression)]
```

The following relational operators can be used in logical expressions:

Symbol	Definition
EQ or =	Equal to
NE or ~= or != or <>	Not equal to
LT or <	Less than
LE or <=	Less than or equal to
GT or >	Greater than
GE or >=	Greater than or equal to

The following logical operators can be used in logical expressions:

Symbol	Definition
AND or &	Both relations must be true
OR or	Either relation can be true
NOT	Reverses the outcome of an expression

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
SELECT IF (SEX EQ 'MALE').
```

## Overview

`SELECT IF` permanently selects cases for analysis based on logical conditions that are found in the data. These conditions are specified in a *logical expression*. The logical expression can contain relational operators, logical operators, arithmetic operations, and any functions that are allowed in `COMPUTE` transformations. For temporary case selection, specify a `TEMPORARY` command before `SELECT IF`.

### Basic Specification

The basic specification is simply a logical expression.

### Syntax Rules

- Logical expressions can be simple logical variables or relations, or these expressions can be complex logical tests involving variables, constants, functions, relational operators, and logical operators. The logical expression can use any of the numeric or string functions that

are allowed in `COMPUTE` transformations (see `COMPUTE` and see Transformation Expressions on p. 68).

- Parentheses can be used to enclose the logical expression. Parentheses can also be used within the logical expression to specify the order of operations. Extra blanks or parentheses can be used to make the expression easier to read.
- A relation can compare variables, constants, or more complicated arithmetic expressions. Relations cannot be abbreviated. For example, `(A EQ 2 OR A EQ 5)` is valid while `(A EQ 2 OR 5)` is not valid. Blanks (not commas) must be used to separate relational operators from the expressions that are being compared.
- A relation cannot compare a string variable to a numeric value or variable, or vice versa. A relation cannot compare the result of the logical functions `SYSMIS`, `MISSING`, `ANY`, or `RANGE` to a number.
- String values that are used in expressions must be specified in quotation marks and must include any leading or trailing blanks. Lowercase letters are considered distinct from uppercase letters.

### **Operations**

- `SELECT IF` permanently selects cases. Cases that are not selected are dropped from the active dataset.
- The logical expression is evaluated as true, false, or missing. If a logical expression is true, the case is selected; if it is false or missing, the case is not selected.
- Multiple `SELECT IF` commands that are issued prior to a procedure command must all be true for a case to be selected.
- `SELECT IF` should be placed before other transformations for efficiency.
- Logical expressions are evaluated in the following order: numeric functions, exponentiation, arithmetic operators, relational operators, and logical operators. Use parentheses to change the order of evaluation.
- If `N OF CASES` is used with `SELECT IF`, the program reads as many records as required to build the specified *n* cases. It makes no difference whether `N OF CASES` precedes or follows `SELECT IF`.
- System variable `$CASENUM` is the sequence number of a case in the active dataset. Although it is syntactically correct to use `$CASENUM` on `SELECT IF`, it does not produce the expected results. To select a set of cases based on their sequence in a file, create your own sequence variable with the transformation language prior to making the selection (see the Examples on p. 1850).

### **Missing Values**

- If the logical expression is indeterminate because of missing values, the case is not selected. In a simple relational expression, a logical expression is indeterminate if the expression on either side of the relational operator is missing.

- If a compound expression is used in which relations are joined by the logical operator `OR`, the case is selected if either relation is true, even if the other relation is missing.
- To select cases with missing values for the variables within the expression, use the missing-value functions. To include cases with values that have been declared user-missing, along with other cases, use the `VALUE` function.

### **Limitations**

`SELECT IF` cannot be placed within a `FILE TYPE-END FILE TYPE` or `INPUT PROGRAM-END INPUT PROGRAM` structure. `SELECT IF` can be placed nearly anywhere following these commands in a transformation program. For more information, see the topic [Commands and Program States](#) on p. 2196.

## **Examples**

### **Working With Simple Logical Expressions**

```
SELECT IF (SEX EQ 'MALE').
```

- All subsequent procedures will use only cases in which the value of `SEX` is `MALE`.
- Because uppercase and lowercase are treated differently in comparisons of string variables, cases for which the value of `SEX` is `male` are not selected.

```
SELECT IF (INCOME GT 75000 OR INCOME LE 10000).
```

- The logical expression tests whether a case has a value that is either greater than 75,000 or less than or equal to 10,000. If either relation is true, the case is used in subsequent analyses.

```
SELECT IF (V1 GE V2).
```

- This example selects cases where variable `V1` is greater than or equal to `V2`. If either `V1` or `V2` is missing, the logical expression is indeterminate, and the case is not selected.

```
SELECT IF (SEX = 'F' & INCOME <= 10000).
```

- The logical expression tests whether string variable `SEX` is equal to `F` and whether numeric variable `INCOME` is less than or equal to 10,000. Cases that meet both conditions are included in subsequent analyses. If either `SEX` or `INCOME` is missing for a case, the case is not selected.

```
SELECT IF (SYSMIS(V1)).
```

- The logical expression tests whether `V1` is system-missing. If it is system-missing, the case is selected for subsequent analyses.

```
SELECT IF (VALUE(V1) GT 0).
```

- Cases are selected if `V1` is greater than 0, even if the value of `V1` has been declared user-missing.

```
SELECT IF (V1 GT 0).
```

- Cases are not selected if `V1` is user-missing, even if the user-missing value is greater than 0.



```
SELECT IF ((V1-15) LE (V2*(-0.001))).
```

- The logical expression compares whether *V1* minus 15 is less than or equal to *V2* multiplied by  $-0.001$ . If it is true, the case is selected.

```
SELECT IF ((YRMODA(88,13,0) - YRMODA(YVAR,MVAR,DVAR)) LE 30).
```

- The logical expression subtracts the number of days representing the date (*YVAR*, *MVAR*, and *DVAR*) from the number of days representing the last day in 1988. If the difference is less than or equal to 30, the case is selected.

### **Understanding and Changing the Order of Evaluation**

```
SELECT IF (RECEIV GT DUE AND (REVNUS GE EXPNS OR BALNCE GT 0)).
```

- By default, AND is executed before OR. This expression uses parentheses to change the order of evaluation.
- The program first tests whether variable *REVNUS* is greater than or equal to variable *EXPNS*, or variable *BALNCE* is greater than 0. Second, the program tests whether *RECEIV* is greater than *DUE*. If one of the expressions in parentheses is true and *RECEIV* is greater than *DUE*, the case is selected.
- Without the parentheses, the program would first test whether *RECEIV* is greater than *DUE* and *REVNUS* is greater than or equal to *EXPNS*. Second, the program would test whether *BALNCE* is greater than 0. If the first two expressions are true *or* if the third expression is true, the case is selected.

### **Selecting Cases Based on Their Sequence in a File**

```
COMPUTE #CASESEQ=#CASESEQ+1.
SELECT IF (MOD(#CASESEQ,2)=0).
```

- This example computes a scratch variable, *#CASESEQ*, containing the sequence numbers for each case. Every other case is selected, beginning with the second case.
- *#CASESEQ* must be a scratch variable so that it is not reinitialized for every case. An alternative is to use the LEAVE command.

### **Using SELECT IF within a DO IF Structure**

```
DO IF SEX EQ 'M'.
+ SELECT IF PRESTIGE GT 50.
ELSE IF SEX EQ 'F'.
+ SELECT IF PRESTIGE GT 45.
END IF.
```

- The SELECT IF commands within the DO IF structure select males with prestige scores above 50 and females with prestige scores above 45.

# SELECTPRED

SELECTPRED is available in IBM® SPSS® Statistics Server.

```
SELECTPRED dependent variable [(MLEVEL={S})]
                                {O}
                                {N}
    [BY factor list] [WITH covariate list]

[/EXCEPT VARIABLES=varlist]

[/SCREENING [STATUS={ON**}] [PCTMISSING={70** }]]
            {OFF }          {number}
    [PCTEQUAL={95** }]] [PCTUNEQUAL={90** }]]
            {number}          {number}
    [STDDEV={0** }]]
            {number}
    [CV={0.001** }]]
            {number}

[/CRITERIA [SIZE={AUTO** }]] [PCUTOFF={0.05**}]
           {integer}          {number}
    [RANKING={PEARSONCHISQ**}] [TIMER={5** }]]
           {LRCHISQ }          {number}
           {CRAMERSV }
           {LAMBDA }
    [SHOWUNSELECTED={0** }]]
           {integer}

[/MISSING USERMISSING={EXCLUDE**}]
           {INCLUDE }

[/PRINT [CPS**] [EXCLUDED**] [SELECTED**] [NONE]]
        [SUMMARY**[(PEARSONCHISQ) [LRCHISQ) [CRAMERSV) [LAMBDA)]]]

[/PLOT SUMMARY]
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release history

Release 14.0

- Command introduced.

## Example

```
SELECTPRED response_01.
```

## Overview

The SELECTPRED procedure submits a large number of predictor variables and selects a smaller subset for use in predictive modeling procedures. This procedure uses a univariable method that considers each predictor in isolation, as opposed to the multivariable method of selecting procedures that is used by the NAIVEBAYES procedure. The SELECTPRED procedure

supports both categorical and scale dependent variables and accepts very large sets of predictors. SELECTPRED is useful when rapid computing is required.

### **Options**

**Methods.** The SELECTPRED procedure includes a screening step as well as the univariable predictor selection method.

**Missing values.** Cases with missing values for the dependent variable or for all predictors are excluded. The SELECTPRED procedure has an option for treating user-missing values of categorical variables as valid. User-missing values of scale variables are always treated as invalid.

**Output.** SELECTPRED displays pivot table output by default but offers an option for suppressing most such output. The procedure optionally displays lists of categorical and scale predictors—not a model—by way of global macro variables. These global macro variables can be used to represent the subset of selected predictors in subsequent procedures.

### **Basic specification**

The basic specification is the SELECTPRED command followed by a dependent variable.

By default, SELECTPRED determines the measurement level of the dependent variable based on its dictionary setting. All other variables in the dataset — except the weight variable if it is defined — are treated as predictors, with the dictionary setting of each predictor determining its measurement level. SELECTPRED performs the screening step and then selects a subset of predictors by using a univariable method. User-missing values are excluded, and default pivot table output is displayed.

### **Syntax rules**

- All subcommands are optional.
- Only a single instance of each subcommand is allowed.
- An error occurs if a keyword is specified more than once within a subcommand.
- Parentheses, equals signs, and slashes that are shown in the syntax chart are required.
- The command name, subcommand names, and keywords must be spelled in full.
- Empty subcommands are not honored.

### **Operations**

The SELECTPRED procedure begins by excluding the following types of cases and predictors.

- Cases with a missing value for the dependent variable.
- Cases with missing values for all predictors.
- Predictors with missing values for all cases.
- Predictors with the same value for all cases.

A further screening step can be used to exclude the following types of predictors:

- Predictors with a large percentage of missing values.

- Categorical predictors with a large percentage of cases representing a single category.
- Categorical predictors with a large percentage of categories containing one case.
- Scale predictors with a small coefficient of variation (standard deviation divided by mean).

The univariable predictor selection method also ranks each predictor based on its association with the dependent variable and selects the top subset of predictors to use in subsequent modeling procedures. The ranking criterion that is used for any given predictor depends on the measurement level of the dependent variable as well as the measurement level of the predictor.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

#### **Categorical dependent variable.**

- For categorical predictors, the ranking criterion can be the Pearson chi-square p-value, likelihood ratio chi-square p-value, Cramér's  $V$ , or Lambda.
- For scale predictors, the F p-value from a one-way ANOVA is always used as the criterion.
- For mixed predictors, the categorical predictors use the chi-square p-value or the likelihood ratio chi-square p-value, and the scale predictors use the F p-value from a one-way ANOVA.

#### **Scale dependent variable.**

- For categorical predictors, the F p-value from a one-way ANOVA is always used as the criterion.
- For scale predictors, the p-value for a Pearson correlation coefficient is always used.
- For mixed predictors, both types of p-values are used, depending on the predictor type.

#### **Frequency weight**

If a `WEIGHT` variable is specified, its values are used as frequency weights by the `SELECTPRED` procedure.

- Cases with missing weights or weights that are less than 0.5 are not used in the analyses.
- The weight values are rounded to the nearest whole numbers before use. For example, 0.5 is rounded to 1, and 2.4 is rounded to 2.

#### **Limitations**

`SPLIT FILE` settings are ignored by the `SELECTPRED` procedure.

## **Examples**

```
SELECTPRED response_01
  /EXCEPT VARIABLES=custid response_02 response_03
  /MISSING USERMISSING=INCLUDE.
```

- This analysis specifies `response_01` as the dependent variable.

- All other variables are to be considered as possible predictors, with the exception of *custid*, *response\_02*, and *response\_03*.
- User-missing values of categorical variables are treated as valid for the purpose of selecting predictors.
- All other settings fall back to their default values.

## Variable lists

The variable lists specify the dependent variable, any categorical predictors (also known as factors), and any scale predictors (also known as covariates).

- The dependent variable must be the first specification on `SELECTPRED`.
- The dependent variable may not be the weight variable.
- The dependent variable may be followed by the measurement-level specification, which contains, in parentheses, the `MLEVEL` keyword followed by an equals sign and then `S` for scale, `O` for ordinal, or `N` for nominal. `SELECTPRED` treats ordinal and nominal dependent variables equivalently as categorical.
- If a measurement level is specified, it temporarily overrides the dependent variable's setting in the data dictionary.
- If no measurement level is specified, `SELECTPRED` defaults to the dictionary setting.
- If a measurement level is not specified, and no setting is recorded in the data dictionary, a numeric variable is treated as scale and a string variable is treated as categorical.
- A string variable may be defined as ordinal or nominal only.
- The names of the factors, if any, must be preceded by the keyword `BY`. If `BY` is specified with no factors, a warning is issued, and the keyword is ignored.
- The names of the covariates, if any, must be preceded by the keyword `WITH`. If `WITH` is specified without covariates, a warning is issued, and the keyword is ignored.
- If the dependent variable or the weight variable is specified within a factor list or a covariate list, the variable is ignored in the list.
- All variables that are specified within a factor or covariate list must be unique. If duplicate variables are specified within a list, the duplicates are ignored.
- If duplicate variables are specified across the factor and covariate lists, an error is issued.
- The universal keywords `TO` and `ALL` may be specified in the factor and covariate lists.
- If neither `BY` nor `WITH` is specified, all variables in the active dataset except the dependent variable, the weight variable, and any variables that are specified on the `EXCEPT` subcommand, are treated as predictors. If the dictionary setting of a predictor is nominal or ordinal, the predictor is treated as a factor. If the dictionary setting is scale, the predictor is treated as a covariate.
- The dependent variable and factor variables can be numeric or string.
- The covariates must be numeric.

## **EXCEPT subcommand**

The EXCEPT subcommand lists any variables that the SELECTPRED procedure should exclude from the factor or covariate lists on the command line. This subcommand is useful if the factor or covariate lists contain a large number of variables—specified by using the TO or ALL keyword, for example—but there are a few variables (for example, Case ID) that should be excluded.

- The EXCEPT subcommand ignores duplicate variables, the dependent variable, and variables that are not specified on the command line's factor or covariate lists.
- There is no default variable list on the EXCEPT subcommand.

## **SCREENING subcommand**

The SCREENING subcommand specifies settings for the screening step, which excludes unsuitable predictors.

<b>STATUS=option</b>	<i>Perform the screening step.</i> Valid options are ON or OFF. If OFF is specified, any other SCREENING specifications are ignored. The default value is ON.
<b>PCTMISSING=number</b>	<i>Percentage of missing values in a predictor.</i> If the percentage of cases representing a single category in a predictor is greater than the specified value, the predictor is excluded. The specified value must be a positive number that is less than or equal to 100. The default value is 70.
<b>PCTEQUAL=number</b>	<i>Percentage of cases representing a single category in a categorical predictor.</i> If the percentage of cases representing a single category in a predictor is greater than the specified value, the predictor is excluded. This setting applies only to categorical predictors. The specified value must be a positive number that is less than or equal to 100. The default value is 95.
<b>PCTUNEQUAL=number</b>	<i>Percentage of categories containing one case in a categorical predictor.</i> If the percentage of a predictor's categories containing only one case is greater than the specified value, the predictor is excluded. This setting applies only to categorical predictors. The specified value must be a positive number that is less than or equal to 100. The default value is 90.
<b>STDDEV=number</b>	<i>Minimum standard deviation for a scale predictor.</i> If the standard deviation is less than the specified value, the predictor is excluded. This setting applies only to scale predictors. The specified value must be a non-negative number. The default value is 0, which turns off the standard deviation check.
<b>CV=number</b>	<i>Coefficient of variation for a scale predictor.</i> A predictor's coefficient of variation is defined as its standard deviation divided by its mean. If the absolute value of the coefficient of variation is less than the specified value, the predictor is excluded. This setting applies only to scale predictors and applies only if the mean is not equal to 0. The specified value must be a positive number. The default value is 0.001.

## **CRITERIA subcommand**

The CRITERIA subcommand specifies computational settings for the SELECTPRED procedure.

**SIZE keyword**

The **SIZE** keyword specifies the number of predictors to select. Value **AUTO** indicates that the number should be computed automatically. Alternatively, a positive integer less than the number of unique predictors on the **SELECTPRED** command may be specified. **AUTO** is the default.

**RANKING keyword**

The **RANKING** keyword specifies the ranking criterion that is used for categorical predictors if the target is categorical.

- PEARSONCHISQ** *Rank predictors based on Pearson chi-square p-value.* Predictors are sorted in ascending order of Pearson chi-square p-values. This criterion is the default.
- LRCHISQ** *Rank predictors based on likelihood ratio chi-square p-value.* Predictors are sorted in ascending order of likelihood ratio chi-square p-values.
- CRAMERSV** *Rank predictors based on Cramer's V.* Predictors are sorted in descending order of Cramer's *V*. If this criterion is used when the target is categorical but predictors are mixed (that is, some categorical, some scale), the **PEARSONCHISQ** default is used instead.
- LAMBDA** *Rank predictors based on Lambda (asymmetric).* Predictors are sorted in descending order of Lambda. If this criterion is used when the target is categorical but predictors are mixed; that is, some categorical, some scale, then the **PEARSONCHISQ** default is used instead.

**PCUTOFF keyword**

The **PCUTOFF** keyword specifies the cutoff whenever a p-value is used as a ranking criterion. More particularly, if p-values are used to rank predictors, only those predictors for which the p-value is less than the **PCUTOFF** value may be selected. A positive number that is less than or equal to 1 may be specified. Specifying 1 turns off the limit on p-values. The default is 0.05.

**SHOWUNSELECTED keyword**

The **SHOWUNSELECTED** keyword specifies a limit on the number of unselected predictors to display in tables and charts. If the specified number *n* is greater than 0, the top *n* unselected predictors are displayed. A non-negative integer may be specified. The default is 0.

**TIMER keyword**

The **TIMER** keyword specifies the maximum number of minutes during which the **SELECTPRED** procedure can be run. If the time limit is exceeded, the procedure is terminated, and no results are given. Any number that is greater than or equal to 0 may be specified. Specifying 0 turns the timer off completely. The default is 5.

**MISSING Subcommand**

The **MISSING** subcommand controls whether user-missing values for categorical variables are treated as valid values. By default, user-missing values for categorical variables are treated as invalid.

- User-missing values for scale variables are always treated as invalid.
- System-missing values for any variables are always treated as invalid.

**USERMISSING=EXCLUDE**     *User-missing values for categorical variables are treated as invalid. This setting is the default.*

**USERMISSING=INCLUDE**     *User-missing values for categorical variables are treated as valid values.*

## ***PRINT subcommand***

The `PRINT` subcommand indicates the tabular output to display. If `PRINT` is not specified, the default tables are displayed. If `PRINT` is specified, only the requested tables are displayed.

### ***CPS Keyword***

The `CPS` keyword displays the case processing summary table, which summarizes the number of cases that are included and excluded in the analysis.

### ***EXCLUDED keyword***

The `EXCLUDED` keyword displays a table of screened predictors. The table lists screened predictors by type (categorical or scale) and reason for being screened. The table includes predictors that are screened due to missing or constant values for all cases, as well as any additional screening that is requested on the `SCREENING` subcommand.

### ***SUMMARY keyword***

The `SUMMARY` keyword displays a statistical summary of selected and unselected predictors. Selected predictors are always displayed. The maximum number of unselected predictors that can be displayed is determined by the `SHOWUNSELECTED` option on the `CRITERIA` subcommand.

The `SUMMARY` keyword may be followed, optionally, by one or more ranking criterion keywords in parentheses. When the target is categorical, and predictors are categorical or of mixed type, the `RANKING` option on the `CRITERIA` subcommand specifies the ranking criterion actually used by the algorithm to select predictors. The parenthesized list that follows the `SUMMARY` keyword displays additional ranking criteria for comparative purposes. The criterion that is actually used is always displayed, irrespective of whether it is specified following `SUMMARY`.

If a parenthesized list follows `SUMMARY`, then keywords in the list may be given in any order. If a specified keyword is inapplicable, it is ignored, and a warning is issued.

**PEARSON-CHISQ**     *Display the Pearson chi-square statistics.*

**LRCHISQ**     *Display the likelihood ratio chi-square statistics.*

**CRAMERSV**     *Display the Cramer's V statistics.*

**LAMBDA**     *Display the Lambda statistics.*



***SELECTED keyword***

The `SELECTED` keyword displays a table of selected predictors by type (categorical or scale).

***NONE keyword***

The `NONE` keyword suppresses all tabular output except the Notes table and any warnings. This keyword may not be specified with any other keywords.

***PLOT subcommand***

The `PLOT` subcommand and `SUMMARY` keyword display a chart that summarizes the selected and unselected predictors. Selected predictors are always displayed. The maximum number of unselected predictors that can be displayed is determined by the `SHOWUNSELECTED` option on the `CRITERIA` subcommand.

# SET

*Note:* Square brackets used in the SET syntax chart are required parts of the syntax and are not used to indicate optional elements. All subcommands are optional.

```
SET  WORKSPACE={6144**}  MXCELLS={AUTOMATIC**}
      {n }                {n }

FORMAT={F8.2**}  CTEMPLATE {NONE**}
      {Fw.d }          {'full path and name'}

TLOOK={NONE**}
      {'full path and name'}

ONUMBERS={LABELS**}  OVARs={LABELS**}
      {VALUES }      {NAMES }
      {BOTH }        {BOTH }

TFIT={BOTHLARGE**}  TNUMBERS={LABELS**}  TVARS={LABELS**}
      {LABELS }      {VALUES }      {NAMES }
      {BOTH }        {BOTH }        {BOTH }

RNG={MC**}
      {MT }

SEED={2000000**}
      {RANDOM }
      {n }

MTINDEX={2000000**}
      {RANDOM }
      {n }

EPOCH={AUTOMATIC}
      {begin year}

{ERRORS } = {LISTING**}
{RESULTS } = {NONE}

{PRINTBACK} = {NONE**}
{MESSAGES } = {LISTING}

MEXPAND={ON**}  MPRINT={OFF**}  MNEST={50**}  MITERATE={1000**}
      {OFF }      {ON }          {n }          {n }

BLANKS={SYSMIS**}  UNDEFINED={WARN**}
      {value }      {NOWARN}

MXWARNS={10**}  MXLOOPS={40**}  MXERRS={100**}
      {n }        {n }          {n }

BLOCK={X'2A'**}
      {X'hexchar' }
      {'character'}

BOX={X'2D7C2B2B2B2B2B2B2B2B2B2B'**}
      {X'hexstring' }
      {'character' }

CCA={'-,,, ' }  CCB={'-,,, ' }  CCC={'-,,, ' }
      {'format' }  {'format' }  {'format' }

CCD={'-,,, ' }  CCE={'-,,, ' }
      {'format' }  {'format' }

DECIMAL={COMMA}
      {DOT }

CACHE={20**}
```

```

        {n }
HEADER={NO** }
        {YES }
        {BLANK}

LENGTH={59** } WIDTH={80** }
        {n }           {255 }
        {NONE}         {n }

SMALL=n

OLANG=output language

DEFOLANG=default output language

SCALEMIN=n

SORT={EXTERNAL** }
        {INTERNAL }
        {SS }

LOCALE='localeid'

THREADS=n

MCACHE=n

UNICODE={OFF** }
        {ON }

FUZZBITS={6** }
        {n }

MIOOUTPUT = [OBSERVED** IMPUTED** POOLED** DIAGNOSTICS**]

ROWSBREAK={100** }
        {n }
        {NONE }

CELLSBREAK={10000** }
        {n }

TOLERANCE={1** }
        {n }

ZCOMPRESSION={YES}
        {NO }

CMPTRANS={YES}
        {NO }

XVERSION={DEFAULT** }
        {'version'}

OATTRS={OLANG** }
        {ENG }
        {BOTH }

DIGITGROUPING={NO** }
        {YES }

TABLERENDER={FULL** }
        {LIGHT }

REPDEFER={NO }
        {YES }

```

\*\* Default setting at installation.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

**Release History**

## Release 13.0

- RNG and MTINDEX subcommands introduced.
- Default for MXERRS subcommand increased to 100.
- SORT subcommand introduced.
- LOCALE subcommand introduced.

## Release 14.0

- Default for WORKSPACE subcommand increased to 6144.

## Release 15.0

- LABELS replaces VALUES as the default for the TNUMBERS subcommand.
- JOURNAL subcommand is obsolete and no longer supported.
- Value EXTERNAL added to list for SORT subcommand, replacing the value INTERNAL as the default. Value SS is deprecated.

## Release 16.0

- MCACHE subcommand introduced.
- THREADS subcommand introduced.
- UNICODE subcommand introduced.

## Release 16.0.1

- BOTHLARGE keyword introduced for the TFIT subcommand.

## Release 17.0

- FUZZBITS subcommand introduced.
- MIOOUTPUT subcommand introduced.

## Release 18.0

- ROWSBREAK, CELLSBREAK, and TOLERANCE subcommands introduced for controlling display of large pivot tables.
- ZCOMPRESSION subcommand introduced.
- COMPRESSION subcommand is obsolete and ignored.
- REPDEFER subcommand introduced.

## Release 19.0

- XVERSION subcommand introduced.
- OATTRS subcommand introduced.
- DIGITGROUPING subcommand introduced.
- TABLERENDER subcommand introduced.
- CMPTRANS subcommand introduced.

## Overview

Many of the running options in the program can be tailored to your own preferences with the `SET` command. The default settings for these options vary from system to system. To display the current settings, use the `SHOW` command. A setting that is changed by `SET` remains in effect for the entire working session unless changed again by another `SET` command. The `PRESERVE` command saves the current settings so that you can return to them later in the session by using the `RESTORE` command. `PRESERVE` and `RESTORE` are especially useful with the macro facility.

### Options

**Memory Management.** Use the `WORKSPACE` subcommand to dynamically allocate memory when some procedures indicate memory shortage. Use the `MXCELLS` subcommand to increase maximum cell numbers for a pivot table.

**Output Format.** Use the `FORMAT` subcommand to change the default (F8.2) print and write formats used for numeric variables. Use the `TLOOK` and `CTEMPLATE` subcommands to specify a TableLook file and/or a chart template file. (*Note:* TableLooks created in earlier versions of IBM® SPSS® Statistics cannot be used in version 16.0 or later.) Use the `ONUMBERS`, `OVARs`, `TNUMBERS`, and `TVARs` subcommands to define default display of variables in the outline or pivot tables. Use the `TFIT` subcommand to specify default column widths. Use the `DIGITGROUPING` subcommand to specify whether the current SPSS Statistics locale's digit grouping format is applied to numeric values in pivot tables, charts, and Model Viewer items, as well as in the Data Editor.

**Samples and Random Numbers.** You can use the `RNG`, `SEED`, and `MTINDEX` subcommands to change the random number generator and initialization value.

**Output Destination.** You can use the `ERRORS`, `MESSAGES`, `PRINTBACK`, and `RESULTS` subcommands to send error messages, resource-utilization messages, command printback, and the output from your commands to your screen and/or to a file. You can also suppress each of these items by using the keyword `NONE`.

**Journal Files.** The `JOURNAL` subcommand is obsolete. You can only set the journal location via the Options dialog (Edit menu>Options>File Locations tab).

**Macro Displays.** You can use the `MEXPAND`, `MITERATE`, and `MNEST` subcommands to control macro expansion, the maximum number of loop iterations, and nesting levels within a macro. You can also use the `MPRINT` subcommands to control the display of the variables, commands, and parameters that a macro uses.

**Blanks and Undefined Input Data.** You can use the `BLANKS` subcommand to specify the value that the program should use when it encounters a completely blank field for a numeric variable. You can also use `UNDEFINED` to turn off the warning message that the program issues when it encounters an invalid value for a numeric variable.

**Maximum Warnings.** You can use `MXWARNs` to limit the warning messages for each set of commands that read the data, after which further warnings are suppressed.

**Maximum Loops.** You can use `MXLOOPs` to raise or lower the maximum number of iterations that are allowed for the `LOOP-END LOOP` structure.

**Scratch File Compression.** You can specify whether scratch files are kept in compressed or uncompressed form using the `ZCOMPRESSION` subcommand.

**Custom Currency Formats.** You can customize currency formats for your own applications using the `CCA`, `CCB`, `CCC`, `CCD`, and `CCE` subcommands. For example, you can display currency as French francs rather than American dollars.

**Cache File.** The `CACHE` subcommand creates a complete copy of the active data file in temporary disk space after a specified number of changes in the active data file. Caching the active data file can improve performance.

**Large Pivot Table Display.** The `ROWSBREAK`, `CELLSBREAK`, and `TOLERANCE` subcommands control the display of large pivot tables in the Viewer. By default, tables with many rows are displayed in sections of 100 rows. To turn off this feature and always display the complete table regardless of size, specify `ROWSBREAK=NONE`. [For more information, see the topic ROWSBREAK, CELLSBREAK, and TOLERANCE Subcommands on p. 1880.](#)

### ***Basic Specification***

The basic specification is at least one subcommand.

### ***Subcommand Order***

Subcommands can be specified in any order.

### ***Syntax Rules***

- You can specify as many subcommands as needed. Subcommands must be separated by at least one space or slash.
- Only one keyword or argument can be specified for each subcommand.
- `SET` can be used more than once in the command sequence.
- `YES` and `ON` are aliases for each other. `NO` and `OFF` are aliases for each other.

### ***Operations***

- Settings that are specified on `SET` remain in effect until they are changed by another `SET` command or until the current session is ended.
- Each time that `SET` is used, only the specified settings are changed. All other settings remain at their previous settings or the default.
- Where filenames are specified, they must include the full path. Relative file specifications, file handles, and filenames without a path are not allowed.

## ***Example***

```
SET BLANKS=0 / UNDEFINED=NOWARN.
```

- `BLANKS` specifies 0 as the value that the program should use when it encounters a completely blank field for a numeric variable.
- `UNDEFINED=NOWARN` suppresses the message that is displayed whenever anything other than a number or a blank is encountered as the value for a numeric variable.

## ***WORKSPACE and MXCELLS Subcommands***

`WORKSPACE` allocates more memory for some procedures when you receive a message indicating that the available memory has been used up or indicating that only a given number of variables can be processed. `MXCELLS` increases the maximum number of cells you can create for a new pivot table when you receive a warning that a pivot table cannot be created because it exceeds the maximum number of cells that are allowed.

- `WORKSPACE` allocates workspace memory in kilobytes for some procedures that allocate only one block of memory. The default is 6144.
- Do not increase the workspace memory allocation unless the program issues a message that there is not enough memory to complete a procedure.
- Use `MXCELLS` with caution. Set `MXCELLS` at a number higher than the limit indicated in the warning message that you receive. After the table is created, restore the number to the default.
- The memory-allocation or cell maximum number increase takes effect as soon as you run the `SET` command.

*Note:* The `MXMEMORY` subcommand is no longer supported.

## ***FORMAT Subcommand***

`FORMAT` specifies the default print and write formats for numeric variables. This default format applies to numeric variables that are defined on `DATA LIST` in freefield format and to all numeric variables that are created by transformation commands (unless a format is explicitly specified).

- The specification must be a simple `F` format. The default is `F8.2`.
- You can use the `PRINT FORMATS`, `WRITE FORMATS`, and `FORMATS` commands to change print and write formats.
- Format specifications on `FORMAT` are output formats. When specifying the width, enough positions must be allowed so that any punctuation characters, such as decimal points, commas, and dollar signs, can be included.
- If a numeric data value exceeds its width specification, the program still attempts to display some value. The program rounds decimal values, removes punctuation characters, tries scientific notation, and finally, if there is still not enough space, produces asterisks indicating that a value is present but cannot be displayed in the assigned width.

## ***TLOOK and CTEMPLATE Subcommands***

`TLOOK` and `CTEMPLATE` specify a file that is used to define the table and chart appearance in the output. The default for either command is `NONE`, which produces tables and charts that use the system defaults.

- **TLOOK** determines the properties of output tables that are produced. The properties include the borders, placement of titles, column and row labels, text font, and column and cell formats. (*Note:* TableLooks created in earlier versions of IBM® SPSS® Statistics cannot be used in version 16.0 or later.)
- **CTEMPLATE** determines the properties of output charts and plots. The properties include line style, color, fill pattern, and text font of relevant chart elements (such as frames, titles, labels, and legends).
- The specification on **TLOOK** or **CTEMPLATE** remains in effect until a new **TLOOK** or **CTEMPLATE** is specified.

**NONE**                    *Use the system defaults.* The tables and charts in the output do not use customized properties.

**filename**                *Use the specified file as a template for tables/charts in the output.* For **TLOOK** you should specify a full path, enclosed in quotes (Directory settings on the **CD** and **INSERT** commands do not apply to the template file location for **TLOOK**).

## ***ONUMBERS, OVARS, TNUMBERS, and TVARS Subcommands***

**ONUMBERS**, **OVARS**, **TNUMBERS**, and **TVARS** control how variables are displayed in the outline for pivot table output and in the pivot tables.

- **ONUMBERS** controls the display of variable values in the outline for pivot tables. The default is **LABELS**.
- **OVARS** controls the display of variables in the outline for pivot tables. The default is **LABELS**.
- **TNUMBERS** controls the display of variable values and/or value labels in the pivot tables. The default is **LABELS**.
- **TVARS** controls the display of variable names and/or variable labels in the pivot tables. The default is **LABELS**.

**NAMES**                    *Display variable names.*

**VALUES**                   *Display variable values.*

**LABELS**                   *Display variable labels.*

**BOTH**                     *Display both labels and values for variables or both names and labels for variables.*

## ***TFIT Subcommand***

**TFIT** controls the default column widths of the pivot tables. The default at installation is **BOTHLARGE**.

**BOTH-LARGE**              *Adjust column widths to accommodate both labels and data, except for very large tables.* For very large tables (more than 10,000 cells), column widths are adjusted to accommodate labels only, and data values that are wider than the column labels may be truncated.

**BOTH**                      *Adjust column widths to accommodate both labels and data.*

**LABELS**                   *Adjust column widths to accommodate labels only.* This setting produces compact tables, but data values that are wider than labels may be truncated.



## ***RNG, SEED, and MTINDEX Subcommands***

Two random number generators are available. The generator that is currently in effect is set by the `RNG` subcommand:

<b>RNG=MC</b>	<i>The random number generator that is used in version 12 and previous releases. If you need to reproduce randomized results that were generated in previous releases based on a specified seed value, use this random number generator. This setting is the default.</i>
<b>RNG=MT</b>	<i>Mersenne Twister random number generator. This generator is a newer random number generator that is more reliable for simulation purposes. If reproducing randomized results from version 12 or earlier is not an issue, use this random number generator.</i>

If you need to reproduce the same randomized results in the future, you can set the initialization value for the random number generator. The initialization value applies to the current session only. It is automatically reset to the default at the start of each new session.

<b>SEED={integer   RANDOM}</b>	<i>Initialization value for MC random number generator. The value must be a positive integer that is less than 2,000,000,000 or the keyword <code>RANDOM</code>, which randomly sets the initialization value. The default is 2,000,000.</i>
<b>MTINDEX={value   RANDOM}</b>	<i>Initialization value for the MT random number generator. The value can be any positive or negative value, including fractional values (expressed as decimals), or the keyword <code>RANDOM</code>. The default is 2,000,000.</i>

### ***Example***

```
SET RNG=MT MTINDEX=-12345.678.
```

Some procedures have internal random number generators, including:

[LINEAR](#)

[GENLINMIXED](#)

## ***EPOCH Subcommand***

`EPOCH` defines the 100-year-span dates that are entered with two-digit years and date functions with a two-digit year specification.

<b>AUTOMATIC</b>	<i>100-year span beginning 69 years prior to the current date and ending 30 years after the current date.</i>
<b>begin year</b>	<i>First year of the 100-year span.</i>

### ***Examples***

```
SET EPOCH=1900.
```

- All dates that are entered with two-digit year values are read as years between 1900 and 1999. For example, a date that is entered as 10/29/87 is read as 10/29/1987.

```
SET EPOCH=1980.
```

- Dates that are entered with two-digit year values between 80 and 99 are read as years between 1980 and 1999.
- Dates that are entered with two-digit year values between 00 and 79 are read as years between 2000 and 2079.

## ***ERRORS, MESSAGES, RESULTS, and PRINTBACK Subcommands***

ERRORS, MESSAGES, RESULTS, and PRINTBACK are used with keywords LISTING and NONE to route program output. ERRORS, MESSAGES, and RESULTS apply only to text output. PRINTBACK applies to all commands that are entered in a syntax window or generated from a dialog box during a session.

- ERRORS refers to both error messages and warning messages for text output.
- MESSAGES refers to resource-utilization messages that are displayed with text output, including the heading and the summaries (such as the amount of memory that is used by a command).
- RESULTS refers to the text output that is generated by program commands.
- PRINTBACK refers to command printback in the log in the Viewer window and in the journal file. Syntax is always displayed as part of the Notes in the syntax window.

**LISTING**      *Display output in the designated output window.* This alias is ON or YES. For PRINTBACK, the alias is BOTH. The executed commands are printed back in the journal and displayed in the log in the output window. You can either display an icon only or list all commands.

**NONE**          *Suppress the output.* The alias is NO or OFF.

The default routes vary from operating system to operating system and vary according to the way commands are executed. In windowed environments, the typical defaults are:

<b>Subcommand</b>	<b>Windowed Environments</b>
ERRORS	LISTING
MESSAGES	NONE
PRINTBACK	BOTH
RESULTS	LISTING

## ***JOURNAL Subcommand***

This subcommand is obsolete and no longer supported. To set the location of the journal file, which contains a log of submitted commands and error and warning messages that are generated during a session, and turn the journal on and off:

- ▶ From the menus in any IBM® SPSS® Statistics window, choose:  
Edit > Options
- ▶ On the File Locations tab, specify the journal location and select journal options.

## ***MEXPAND and MPRINT Subcommands***

`MEXPAND` and `MPRINT` control whether macros are expanded and whether the expanded macros are displayed. For more information about macros, see the `DEFINE` command and Using the Macro Facility on p. 2215.

The specifications for `MEXPAND` are:

- ON**     *Expand macros.* This setting is the default.
- OFF**    *Do not expand macros.* The command line that calls the macro is treated like any other command line. If the macro call is a command, it will be executed; otherwise, it will trigger an error message.

The specifications for `MPRINT` are:

- ON**     *Include expanded macro commands in the output.*
- OFF**    *Exclude expanded macro commands from the output.* This is the default.

- `MPRINT` is effective only when `MEXPAND` is `ON` and is independent of the `PRINTBACK` subcommand.

## ***MITERATE and MNEST Subcommands***

`MITERATE` and `MNEST` control the maximum loop traversals and the maximum nesting levels permitted in macro expansions, respectively.

- The specification on `MITERATE` or `MNEST` is a positive integer. The default for `MITERATE` is 1000. The default for `MNEST` is 50.

## ***BLANKS Subcommand***

`BLANKS` specifies the value that the program should use when it encounters a completely blank field for a numeric variable. By default, the program uses the system-missing value.

- `BLANKS` controls only the translation of numeric fields. If a blank field is read with a string format, the resulting value is a blank.
- The value that is specified on `BLANKS` is not automatically defined as a missing value.
- The `BLANKS` specification applies to all numeric variables. You cannot use different specifications for different variables.
- `BLANKS` must be specified before data are read. Otherwise, blanks in numeric fields are converted to the system-missing value (the default) as they are read.

## **UNDEFINED Subcommand**

UNDEFINED controls whether the program displays a warning message when it encounters anything other than a number or a blank as the value for a numeric variable. The default is `WARN`.

- WARN**            *Display a warning message when an invalid value is encountered for a numeric variable. This setting is the default.*
- NOWARN**        *Suppress warning messages for invalid values.*

## **MXERRS Subcommand**

MXERRS controls the maximum number of errors that are allowed in a session. The default is 100.

- MXERRS applies only to command files that are submitted for execution through IBM® SPSS® Statistics Batch Facility (available with SPSS Statistics Server).
- Eligible errors are errors that stop execution of a command but continue the session.
- When the MXERRS limit is exceeded, SPSS Statistics stops processing commands but continues to scan for additional errors.
- In interactive mode or in SPSS Statistics and other windowed environments, MXERRS does not apply.

## **MXWARNS Subcommand**

MXWARNS controls the number of warnings that are issued. The default is 10. The behavior of this setting depends on mode of operation: interactive or IBM® SPSS® Statistics Batch Facility. In this context, “interactive” includes any mode of operation other than SPSS Statistics Batch Facility.

**Interactive.** In interactive mode, MXWARNS limits the number of warnings that are issued for each set of commands that read the data (for example, a group of transformation commands followed by a statistical procedure).

- Exceeding the limit does not halt execution of commands; it simply suppresses further warnings.
- `MXWARNS=0` suppresses all warnings except a warning that further warnings have been suppressed.
- MXWARNS does not affect the display of error messages, and errors do not count toward the warning limit.

**Batch Facility.** In the Batch Facility (available with SPSS Statistics Server), MXWARNS limits the total number of warnings that are allowed in a job.

- When the MXWARNS limit is exceeded, SPSS Statistics stops processing commands but continues to scan for errors.
- When the combined total number of warnings and errors exceeds the MXWARNS limit, SPSS Statistics stops processing commands but continues to scan for errors.

## ***MXLOOPS Subcommand***

MXLOOPS specifies the maximum number of times that a loop that is defined by the LOOP-END LOOP structure is executed for a single case or input record. The default is 40.

- MXLOOPS prevents infinite loops, which may occur if no cutoff is specified for the loop structure (see LOOP-END LOOP).
- MXLOOPS will limit the number of loops for any loop structure that doesn't have an indexing clause, including loops with conditional IF clauses. If a loop has an indexing clause (e.g., LOOP #i=1 to 1000), the indexing clause overrides the MXLOOPS setting.
- When a loop is terminated, control passes to the command immediately following the END LOOP command, even if the END LOOP condition is not yet met.

## ***EXTENSIONS Subcommand***

This subcommand is no longer supported.

## ***BLOCK Subcommand***

BLOCK specifies the character that is used for drawing icicle plots.

- You can specify any single character either as a quoted string or as a quoted hexadecimal pair preceded by the character X.
- The default is X'2A'.

### ***Example***

```
SET BLOCK=' #' .
```

- This command specifies a pound sign (#) as the character to be used for drawing bar charts. The character is specified as a quoted string.

## ***BOX Subcommand***

BOX specifies the characters that are used to draw table borders in text output. The specification is either a 3-character or 11-character quoted string, in which the characters have the following representations:

1	horizontal line	7	upper right corner
2	vertical line	8	left T
3	middle (cross)	9	right T
4	lower left corner	10	top T
5	upper left corner	11	bottom T
6	lower right corner		

- The characters can be specified either as a quoted string or hexadecimal pairs. Specify an *X* before the quoted hexadecimal pairs.
- The defaults vary from system to system. To display the current settings, use the `SHOW` command.

## ***LENGTH and WIDTH Subcommands***

`LENGTH` and `WIDTH` specify the maximum page length and width for the text output, respectively. These two subcommands apply only to text output.

- Length represents number of lines, and width represents number of characters.
- You can specify any length from 40 through 999,999 lines. If a long page length is specified, the program continues to provide page ejects and titles at the start of each procedure and at logical points in the display, such as between crosstabulations.
- To suppress page ejects, use keyword `NONE` on `LENGTH`. The program will insert titles at logical points in the display but will not supply page ejects.
- You can specify any number of characters from 80 through 255 for `WIDTH`.

## ***HEADER Subcommand***

`HEADER` controls whether the output includes headings. The `HEADER` subcommand applies to both default headings and headings that are specified on the `TITLE` and `SUBTITLE` commands. This command applies only to text output from this program. The default is `NO`.

- NO**            *Suppress headings in text output. All general headings, including pagination, are replaced by a single blank line.*
- YES**            *Display headings in text output.*
- BLANK**        *Suppress headings but start a new page.*

## ***CCA, CCB, CCC, CCD, and CCE Subcommands***

You can use the subcommands `CCA`, `CCB`, `CCC`, `CCD`, and `CCE` to specify up to five custom currency formats.

- Each custom currency subcommand defines one custom format and can include four specifications in the following order: a negative prefix, a prefix, a suffix, and a negative suffix.
- The specifications are delimited by either periods or commas, whichever you do not want to use as a decimal point in the format.
- If your custom currency format includes periods or commas that you need to distinguish from delimiters, use a single quotation mark as an escape character before the period or comma that is part of the custom currency format. For example, if the format includes a period but the decimal indicator is a comma, the period must also be used as the delimiter.
- Each currency specification must always contain three commas or three periods. All other specifications are optional.
- Use blanks in the specification only where you want blanks in the format.

- The entire specification must be enclosed in single or double quotation marks. If the format includes a single quotation mark as an escape character, the entire specification must be enclosed in double quotation marks.
- A specification cannot exceed 16 characters (excluding the apostrophes).
- Custom currency formats cannot be specified as input formats on `DATA LIST`. Use them only as output formats in the `FORMATS`, `WRITE FORMATS`, `PRINT FORMATS`, `WRITE`, and `PRINT` commands.

### ***Specifying a Custom Currency Format***

```
SET CCA='- , $ , ' .
```

- A minus sign (-) preceding the first command is used as the negative prefix.
- A dollar sign is specified for the prefix.
- No suffixes are specified (there are two consecutive commas before the closing apostrophe).
- Because commas are used as separators in the specification, the decimal point is represented by a period.

### ***Specifying Multiple Custom Currency Formats***

```
SET CCA='( , , -)' CCB=' , , % , ' CCC='( , $ , , )' CCD='- / - . Df1 . . - ' .
FORMATS VARA(CCA9.0) / VARB(CCB6.1) / VARC(CCC8.0) / VARD(CCD14.2) .
```

- `SET` defines four custom currency formats.
- `FORMATS` assigns these formats to specific variables.

**Table 225-1**  
*Custom currency examples*

	<b>CCA</b>	<b>CCB</b>	<b>CCC</b>	<b>CCD</b>
negative prefix	(	none	(	-/-
prefix	none	none	\$	Df1
suffix	none	%	none	none
negative suffix	-)	none	)	-
separator	,	,	,	.
sample positive number	23,456	13.7%	\$352	Df1 37.419,00
sample negative number	(19,423-)	13.7%	(\$189)	-/-Df1 135,19-

## ***DECIMAL Subcommand***

`DECIMAL` can be used to override the default decimal indicator for reading and writing text data. The default decimal indicator is the OS locale decimal indicator or the decimal indicator for the locale specified on the `LOCALE` subcommand.

**DOT.** The decimal indicator is a period.

**COMMA.** The decimal indicator is a comma.

- A subsequent `LOCALE` subcommand—either on the same or separate `SET` command—will override the `DECIMAL` setting. For more information, see the topic [LOCALE Subcommand](#) on p. 1876.
- This setting only affects reading and writing text data. It has no effect on how data values are displayed in the Data Editor or in results displayed in the Viewer.

## ***CACHE Subcommand***

The `CACHE` subcommand creates a complete copy of the active data file in temporary disk space after a specified number of changes in the active data file. If you have the available disk space, this feature can improve performance. The default number of changes that can occur before the active file is cached is 20.

### ***Example***

```
SET CACHE 10.
```

## ***SMALL Subcommand***

The `SMALL` subcommand controls the display of numbers in scientific notation in output for small decimal values.

### ***Example***

```
SET SMALL = 0.  
SET SMALL = .001.
```

- The first `SET SMALL` command suppresses the display of scientific notation in all output.
- The second `SET SMALL` command will only display scientific notation for values that are less than 0.001.

## ***OLANG Subcommand***

The `OLANG` subcommand controls the language that is used in output. `OLANG` does not apply to simple text output. Available languages may vary. (The General tab in the Options dialog box displays a list of available output languages.) Valid language keywords include `English`, `French`, `German`, `Spanish`, `Italian`, `Japanese`, `Korean`, `TChinese` (traditional Chinese), `SChinese` (simplified Chinese), `Polish`, and `Russian`.

Output that is produced after the command is executed will be in the specified language (if that language is available). Additional language materials may be available for downloading from the SPSS Inc. web site.

Depending on the language, you may also need to switch to [UNICODE](#) mode or change the [LOCALE](#) setting to match the output language for characters to render properly.



**Example**

```
SET OLANG = German.
```

- The language keyword is not case-sensitive.
- Do *not* enclose the language keyword in quotation marks or other string delimiters.

**DEFOLANG Subcommand**

The DEFOLANG subcommand specifies the default output language (if the language that is specified on the OLANG subcommand is not available). The initial default setting is the language of the installed software version. For example, if you install the English version of the software on a Japanese operating system, the default output language is English.

**Example**

```
SET DEFOLANG = JAPANESE.
```

- The language keyword is not case-sensitive.
- Do *not* enclose the language keyword in quotation marks or other string delimiters.

**SCALEMIN Subcommand**

For IBM® SPSS® Statistics data files that were created prior to release 8.0 and data read from external sources, you can specify the minimum number of data values for a numeric variable that is used to classify the variable as continuous (scale) or nominal. Variables with fewer than the specified number of unique values are classified as nominal. All string variables are classified as nominal, regardless of the number of unique values.

There are numerous other conditions that are evaluated prior to applying the minimum number of data values rule when determining to apply the continuous (scale) or nominal measurement level. Conditions are evaluated in the order listed in the table below. The measurement level for the first condition that matches the data is applied.

Condition	Measurement Level
Format is dollar or custom-currency	Continuous
Format is date or time (excluding Month and Wkday)	Continuous
All values of a variable are missing	Nominal
Variable contains at least one non-integer value	Continuous
Variable contains at least one negative value	Continuous
Variable contains no valid values less than 10,000	Continuous
Variable has <i>N</i> or more valid, unique values*	Continuous
Variable has no valid values less than 10	Continuous
Variable has less than <i>N</i> valid, unique values*	Nominal

\* *N* is the user-specified cut-off value. The default is 24.

## ***SORT Subcommand***

By default, IBM® SPSS® Statistics tries to use an external, third-party sorting mechanism, which may reduce processing time with large data sources. The third-party sorting option is available only if you have SPSS Statistics Server. The specific sorting engine is defined by your server administrator. If you are not connected to the SPSS Statistics Server or the SPSS Statistics Server cannot find the third-party sort engine, the built-in sorting mechanism is used.

<b>EXTERNAL</b>	<i>Use the external, third-party sort engine if available.</i> This setting is the default. If the third-party sort engine is not available, this setting is ignored, and the built-in sorting mechanism is used. COXREG and CURVEFIT use built-in sorting regardless of the SORT setting.
<b>INTERNAL</b>	<i>Use the built-in sorting mechanism.</i>
<b>SS</b>	This setting is deprecated. It has the same effect as EXTERNAL.

## ***LOCALE Subcommand***

The LOCALE subcommand allows you to change the locale used for data analysis. By default, the locale is set to the computer's system locale if possible. With the LOCALE subcommand, you can override the default behavior and process data files in other locales without changing your computer's system or user locale.

- The default locale is the current operating system locale.
- The LOCALE subcommand persists. The next time that IBM® SPSS® Statistics is started on your computer, SPSS Statistics will run in that locale.
- If the locale ID does not match the system locale, not all output will be rendered correctly.
- In code page mode, locale changes that also change the encoding can only be made when there are no open datasets. For information on code page mode and Unicode mode, see [UNICODE Subcommand](#).
- You can use SHOW LOCALE to view the current SPSS Statistics locale.

### ***Example***

```
SET LOCALE='Japanese' .
```

- When you are connecting to a server, the relevant locale ID is defined in the *loclmap.xml* file, which is located on the server computer. Check with your server administrator for available locale IDs.
- Locale ID values of the general form *Language\_Country.codepage#*, which were valid in previous releases, are not valid in release 16.0 or later. In most instances, the “Language” portion of the old locale specification should be a valid locale ID in release 16.0 or later. For example, instead of ‘Japanese\_Japan.932’, you could specify simply ‘Japanese’. For a complete list of names and aliases that can be used to specify locale, go to <http://www.iana.org/assignments/character-sets>

Two-letter language abbreviations adhere to the ISO 639-1 standard:

[http://www.loc.gov/standards/iso639-2/php/code\\_list.php](http://www.loc.gov/standards/iso639-2/php/code_list.php)

Two-letter country codes adhere to the ISO 3166 standard: [http://en.wikipedia.org/wiki/ISO\\_3166-1](http://en.wikipedia.org/wiki/ISO_3166-1)

### **Example**

```
SET LOCALE='ja_JP.Shift_JIS'
```

- “ja” is the two-letter language code for Japanese.
- “JP” is the two-letter country code for Japan.
- “Shift\_JIS” is the name of the character set.

## **THREADS Subcommand**

The `THREADS` subcommand specifies the number of threads that multithreaded procedures use when calculating results. By default, IBM® SPSS® Statistics uses an internal algorithm to determine the number of threads for a particular computer. You can change this setting, but the default will often provide the best performance. The default also appears as the recommended setting when you run the `SHOW` command.

Using more than one thread is relevant only if your computer has multiple processors or each processor has multiple cores. Also, this setting affects only selected procedures that take advantage of it. These are:

- CORRELATIONS
- CSCOXREG
- CSLOGISTIC
- CSORDINAL
- CSSELECT
- DISCRIMINANT
- FACTOR
- NOMREG
- PARTIAL CORR
- REGRESSION
- SORT CASES

*Note:* `PRESERVE` and `RESTORE` do not operate on this subcommand.

## **MCACHE Subcommand**

The `MCACHE` subcommand specifies the memory cache (in kilobytes) that multithreaded procedures use when calculating results. By default, IBM® SPSS® Statistics uses an internal algorithm to determine the memory cache for a particular computer. You can change this setting, but the default will often provide the best performance. The default also appears as the recommended setting when you run the `SHOW` command.

Using the memory cache is relevant only if your computer has multiple processors or each processor has multiple cores. Also, this setting affects only selected procedures that take advantage of it. These are:

- CORRELATIONS
- DISCRIMINANT
- FACTOR
- NOMREG
- PARTIAL CORR
- REGRESSION

*Note:* PRESERVE and RESTORE do not operate on this subcommand.

## **UNICODE Subcommand**

SET UNICODE NO|YES controls the default behavior for determining the encoding for reading and writing data files and syntax files.

**NO.** Use the current locale setting to determine the encoding for reading and writing data and command syntax files. This is referred to as **code page mode**. This is the default. The alias is OFF. For information on the current locale setting, see [LOCALE Subcommand](#).

**YES.** Use Unicode encoding (UTF-8) for reading and writing data and command syntax files. This is referred to as **Unicode mode**. The alias is ON.

- You can change the UNICODE setting only when there are no open data sources.
- The UNICODE setting persists across sessions and remains in effect until it is explicitly changed.

There are a number of important implications regarding Unicode mode and Unicode files:

- Data and syntax files saved in Unicode encoding should not be used in releases prior to 16.0. For syntax files, you can specify local encoding when you save the file. For data files, you should open the data file in code page mode and then resave it if you want to read the file with earlier versions.
- When code page data files are read in Unicode mode, the defined width of all string variables is tripled. You can use [ALTER TYPE](#) to automatically adjust the width of all string variables.
- The GET command determines the file encoding for IBM® SPSS® Statistics data files from the file itself, regardless of the current mode setting (and defined string variable widths in code page files are tripled in Unicode mode). [For more information, see the topic GET on p. 832.](#)
- For text data files read with DATA LIST and related commands (for example, REPEATING DATA and FILE TYPE) or written with PRINT or WRITE, you can override the default encoding with the ENCODING subcommand.
- GET DATA uses the default encoding for reading text data files (TYPE=TXT), which is UTF-8 in Unicode mode or the code page determined by the current locale in code page mode.
- OMS uses default encoding for writing text files (FORMAT=TEXT and FORMAT=TABTEXT) and for writing SPSS Statistics data files (FORMAT=SAV).

- By default, `GET STATA`, `GET TRANSLATE`, and `SAVE TRANSLATE` (with the exception of saving SAS 9 format files) read and write data in the current locale code page, regardless of mode. You can use the `ENCODING` subcommand to specify a different encoding.
- By default, `SAVE TRANSLATE` saves SAS 9 files in UTF-8 format in Unicode mode and in the current locale code page in code page mode. You can use the `ENCODING` subcommand to specify a different encoding.
- For syntax files run via `INCLUDE` or `INSERT`, you can override the default encoding with the `ENCODING` subcommand.
- For syntax files, the encoding is set to Unicode *after* execution of the block of commands that includes `SET UNICODE=YES`. You must run `SET UNICODE=YES` separately from subsequent commands that contain Unicode characters not recognized by the local encoding in effect prior to switching to Unicode.

## ***FUZZBITS Subcommand***

`FUZZBITS` specifies the default setting of the *fuzzbits* argument for the `RND` and `TRUNC` functions. The value controls the number of least-significant bits by which the internal representation of the value to be rounded or truncated (expressed as a 64-bit floating point binary) may fall short of the nearest rounding boundary and be rounded up. [For more information, see the topic Arithmetic functions on p. 73.](#)

- The value must be an integer between 0 and 20. The default is 6, which should be sufficient for most applications.
- A value of 0 produces the same results as in release 10. A value of 10 produces the same results as in releases 11 and 12.
- The `FUZZBITS` setting persists across sessions.

## ***MIOUTPUT Subcommand***

`MIOUTPUT` specifies the display of pooled output when a multiply imputed dataset is analyzed.

<b>OBSERVED</b>	<i>Show results for observed data.</i>
<b>IMPUTED</b>	<i>Show results for each imputed dataset.</i>
<b>POOLED</b>	<i>Show final pooled statistics.</i>
<b>DIAGNOSTICS</b>	<i>Display pooling diagnostics for parameters pooled by the univariate method.</i>

- At least one keyword must be specified in square brackets. If a keyword is not specified the corresponding output is suppressed.
- `DIAGNOSTICS` is ignored with a warning if `POOLED` is not specified.
- Default settings depend on licensing. If the Missing Values option is licensed, all options are on by default. If the Missing Values option is not licensed, the defaults are `OBSERVED` and `IMPUTED`, and specifying `POOLED` and `DIAGNOSTICS` keywords are ignored with a warning. That is, without the Missing Values option you can suppress results for original data and imputations, but cannot obtain pooled results.

- The `MIOOUTPUT` setting persists across sessions. It is ignored by tables that do not support splits. It is also ignored when analysis is performed without variable `Imputation_` defined as a split variable.
- `MIOOUTPUT` has no effect on charts or PMML.

## ***ROWSBREAK, CELLSBREAK, and TOLERANCE Subcommands***

These subcommands control the display of large pivot tables in the Viewer. These settings have no effect on printing large pivot tables or exporting output to external formats with command such as `OUTPUT EXPORT` or `OMS`.

**ROWSBREAK=*n*|NONE.** *Sets the number of rows to display.* The default is 100. The value must be an integer between 10 and 1000 or the keyword `NONE`, which turns off this feature and displays the entire table, regardless of size. When the feature is on and large tables are displayed in sections, navigation controls allow you to view different sections of the table.

**CELLSBREAK=*n*.** *Sets the maximum number of cells to display.* The default is 10000. The value must be an integer between 1000 and 100000. If `ROWSBREAK=NONE`, this setting is ignored.

**TOLERANCE=*n*.** *Controls the maximum number of rows of the inner most row dimension of the table to split across displayed views of the table.* For example, if there are six categories in each group of the inner most row dimension, specifying a value of six would prevent any group from splitting across displayed views. The value must be an integer and cannot be greater than the `ROWSBREAK` value. If `ROWSBREAK=NONE`, this setting is ignored.

Several factors can affect the actual number of rows displayed, including:

- If the `CELLSBREAK` value is reached before the `ROWSBREAK` value, then the table is split at that point. For example, if `ROWSBREAK=1000` and `CELLSBREAK=10000`, a table with 20 columns will be displayed in blocks of 500 rows.
- The `TOLERANCE` setting can cause fewer or more rows than the `ROWSBREAK` value to be displayed.

## ***ZCOMPRESSION Subcommand***

*Note:* This command works only if you are connected to the IBM® SPSS® Statistics Server.

`ZCOMPRESSION` is an override mechanism that works with a setting on the SPSS Statistics Server to determine whether scratch files that are created during a session are in compressed or uncompressed form. Scratch files are created by `CACHE`, `SORT`, `AGGREGATE`, transformations, and any other procedure that adds columns to the data file.

- The administrator of the SPSS Statistics Server must give you permission to override the SPSS Statistics Server setting. If you don't have permission, a warning message is displayed and the command is ignored.
- A compressed scratch file reduces the amount of disk I/O compared to an uncompressed scratch file but requires more processing on the SPSS Statistics Server.

- The extra overhead associated with compression is considerably greater when writing than reading a file. Consider using a higher value for `SET CACHE` to reduce the frequency of writing new scratch files containing the entire data set. A higher `SET CACHE` value will also increase the number of times a scratch file is read.
- If you are working with a very large data file (larger than half of the Server's RAM) and compression is on, you may consider issuing an explicit `CACHE` command before running other procedures. Doing this will ensure the procedures take full advantage of the disk I/O benefits.
- The override specification takes effect the next time that a scratch file is written and stays in effect until `SET ZCOMPRESSION` is specified again or until the end of the session.
- The default setting for compressing scratch files is set on the SPSS Statistics Server. There is no syntax default.

**YES**     *Compress scratch files.*

**NO**      *Do not compress scratch files.*

## ***CMPTRANS Subcommand***

*Note:* This command works only if you are connected to the IBM® SPSS® Statistics Server running on Windows Server..

`CMPTRANS` interacts with a setting on the SPSS Statistics Server to determine whether transformation commands (such as `COMPUTE` and `RECODE`) are compiled to machine code when a syntax job is run.

- If the administrator of the SPSS Statistics Server turns on compiled transformations with the SPSS Statistics Server setting *and* `CMPTRANS` is set to `YES`, transformation commands are compiled to machine code when the syntax job is run. The compiled code is used for the subsequent execution of the transformations.
- If the administrator does not turn on compiled transformations, a warning message is displayed and the command is ignored.
- Compiling transformations may improve performance when there are a large number of cases and multiple transformation commands.
- Because there is an overhead involved in compiling the transformations, you should set `CMPTRANS` to `YES` only when there are a large number of cases and multiple transformation commands.

**YES**     *Compile transformations to machine code.*

**NO**      *Do not compile transformations to machine code.*

## ***DIGITGROUPING Subcommand***

`DIGITGROUPING` specifies whether the current IBM® SPSS® Statistics locale's digit grouping format will be applied to numeric values in pivot tables and charts as well as in the Data Editor. For example, in a French locale, with `DIGITGROUPING` set to `YES` the value 34419,57 is displayed as 34 419,57.

The grouping format does not apply to trees, Model Viewer items, numeric values with the DOT or COMMA format, or numeric values with a DOLLAR or custom currency format. It does, however, apply to the display of the days value for numeric values with a DTIME format—for example, to the value of ddd in the format ddd hh:mm.

- YES**     *Apply the current locale's digit grouping format.*  
**NO**       *Do not apply the current locale's digit grouping format.* This is the default.

## ***XVERSION and OATTRS Subcommands***

XVERSION and OATTRS control aspects of OXML and SAV output from the Output Management System (OMS).

- XVERSION specifies the version of the Output XML schema used to generate OXML output from OMS.
- OATTRS specifies whether language invariant versions of text attributes are included in OXML and whether variable names in datasets generated from SAV output are in English. When specified, the English value of a text attribute is stored in the text\_eng attribute which either replaces the text attribute or is in addition to it. The text\_eng attribute allows you to write XPath expressions in a language invariant manner.

*Note:* The settings specified by XVERSION and OATTRS apply to subsequent OMS requests.

The specifications for XVERSION are:

- DEFAULT**     *Use the schema associated with the version of IBM® SPSS® Statistics from which the SET command is executed.* This is the default.
- 'version'**     *A quoted string specifying the output schema version—for example, '1.6'. The version string is the value of the version attribute of the xs:schema element from the associated output schema file. You can specify the version for any of the installed output schema files. Output schema files are located in the SPSS Statistics installation folder.*

The specifications for OATTRS are:

- OLANG**     *The text\_eng attribute is not included in OXML output and variable names in datasets generated from SAV output are translated into the current output language (when available).* This is the default.
- ENG**         *The text attribute is replaced (other than in cell elements) with a text\_eng attribute whose value is the English value of the text attribute. Variable names in datasets generated from SAV output will be in English.* The value ENG requires that the output schema version in use is 1.6 or higher, which is satisfied by the default setting of the XVERSION subcommand.
- BOTH**       *A text\_eng attribute is added to each element (other than a cell element) that has a text attribute. The value of the text\_eng attribute is the English value of the text attribute. Variable names in datasets generated from SAV output will be in English.* The value BOTH requires that the output schema version in use is 1.6 or higher, which is satisfied by the default setting of the XVERSION subcommand.



## ***TABLERENDER Subcommand***

Tables can be rendered as full-featured pivot tables or as lightweight tables.

- FULL**      *Tables are rendered as full-featured pivot tables. This is the default.*
- LIGHT**     *Tables are rendered as lightweight tables. Lightweight tables render faster but cannot be pivoted or edited.*

Lightweight tables render faster than full-featured pivot tables, but they do not support many pivoting and editing features, including:

- Lightweight tables cannot be pivoted. You can display different layers of a multi-layer table, but you cannot move table elements between dimensions (for example, swap rows and columns) or change the display position of rows or columns.
- Lightweight tables cannot be edited. You cannot hide rows or columns, change column widths, modify data values, change cell or table properties, or apply different TableLooks.
- Lightweight tables do not support custom dimension border properties defined in TableLooks. All other TableLook properties are honored at table creation time.
- Since lightweight tables cannot be edited, the Replace feature of the Find and Replace dialog will not replace the specified value in lightweight tables.

To convert a lightweight table to a full-featured pivot table:

- ▶ Right-click the table in the Viewer window or the table editing window.
- ▶ From the context menu, choose Enable Editing

*Note:* Full-featured pivot tables cannot be converted to lightweight tables. If a table is created as — or converted to — a full-featured pivot table, it remains a full-featured pivot table.

*Note:* For procedures that produce tables with layer dimensions, (CROSSTABS, CTABLES, OLAP CUBES), all layer dimensions are collapsed into a single layer dimension in lightweight tables. To separate the layers, convert the table to a full-featured pivot table.

## ***REPDEFER Subcommand***

REPDEFER specifies whether REPOSITORY CONNECT commands (connection requests to an IBM® SPSS® Collaboration and Deployment Services Repository) are deferred until a connection is required.

- YES**      *Defer connection requests until a connection is required. This is the default when executing command syntax with the IBM® SPSS® Statistics Batch Facility (available with SPSS Statistics Server).*
- NO**        *Attempt to connect when a REPOSITORY CONNECT command is encountered. This is the default unless executing command syntax with the SPSS Statistics Batch Facility.*

# SHIFT VALUES

```
SHIFT VALUES VARIABLE=varname RESULT=varname {LEAD=non-negative integer}
                                                {LAG=non-negative integer }
                                                {SHIFT=integer }
[/VARIABLE=varname...]
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 17.0

- Command introduced.

## **Example**

```
SHIFT VALUES VARIABLE=OldVar RESULT=NewVar LEAD=1.
```

## **Overview**

The `SHIFT VALUES` command creates new variables that contain the values of existing variables from preceding or subsequent cases.

### **Basic Specification**

The basic specification is the command name followed by:

**VARIABLE=varname.** This must be a variable that already exists in the active dataset.

**RESULT=newname.** This must be a variable that does not currently exist in the active dataset.

One of the following alternative shift methods:

- **LEAD=n.** Get the value from a subsequent case in the dataset. The value must be a non-negative integer that indicates the number of cases. For example, `LEAD=1` returns the value of the existing variable from the case immediately following the current case.
- **LAG=n.** Get the value from a preceding case in the dataset. The value must be a non-negative integer that indicates the number of cases. For example, `LAG=1` returns the value of the existing variable from the case immediately before the current case.
- **SHIFT=n.** Get the value from a preceding or subsequent case. The value must be an integer that indicates the number of cases. For example, `SHIFT=1` returns the value of the existing variable from the case immediately following the current case, and `SHIFT=-1` returns the value of the existing variable from the case immediately before the current case.

To specify more than one variable on the same command, use a forward slash (/) to separate the full specification for each subsequent variable.

**Syntax Rules**

- The `RESULT` keyword must specify a new variable name. The name cannot be a variable that already exists in the active dataset. For information on variable naming rules, see [Variable Names](#).
- Each keyword must be followed by an equals sign (=) before the value. For example, `RESULT=NewVar`.

**Operations**

- If split file processing is on, the scope of the shift is limited to each split group. A shift value cannot be obtained from a case in a preceding or subsequent split group.
- Filter status is ignored.
- The value of the result variable is set to system-missing for the first or last  $n$  cases in the dataset or split group, where  $n$  is the number of cases specified on `LEAD`, `LAG`, or `SHIFT`. For example, `LAG=1` would set the result variable to system-missing for the first case in the dataset (or first case in each split group).
- A value of 0 for number of cases will copy the value of the original value to the same case for the new variable.
- User-missing values are preserved.
- Dictionary information from the original variable, including defined value labels and user-missing value assignments, is applied to the new variable. (Note: Custom variable attributes are not included.)
- A variable label is automatically generated for the new variable that describes the shift operation that created the variable.

**Example**

```
DATA LIST FREE /Var1 (F1).
BEGIN DATA
0 1 9 0 1
END DATA.
VARIABLE LABEL Var1 'Original Variable'.
VALUE LABELS Var1 0 'No' 1 'Yes' 9 'Missing'.
MISSING VALUES Var1(9).
SHIFT VALUES VARIABLE=Var1 RESULT=Var1_Lead LEAD=1
/VARIABLE=Var1 RESULT=Var1_Lag LAG=1
/VARIABLE=Var1 RESULT=Var1_Shift SHIFT=-1.
```

**Figure 226-1**  
Results displayed in Data View

Var1	Var1_Lead	Var1_Lag	Var1_Shift
0	1	.	.
1	9	0	0
9	0	1	1
0	1	9	9
1	.	0	0

**Figure 226-2**  
*Dictionary information displayed in Variable View*

Name	Type	Width	Decimals	Label	Values	Missing
Var1	Numeric	1	0	Original Variable	{0, No}...	9
Var1_Lead	Numeric	1	0	Lead(Var1,1)	{0, No}...	9
Var1_Lag	Numeric	1	0	Lag(Var1,1)	{0, No}...	9
Var1_Shift	Numeric	1	0	Shift(Var1,-1)	{0, No}...	9

# SHOW

SHOW **ALL** [BLANKS] [BOX] [BLOCK] [CC] [CCA] [CCB]  
[CCC] [CCD] [CCE] [CACHE] [CTEMPLATE] [DECIMAL] [DEFOLANG] [DIGITGROUPING]  
[DIRECTORY] [ENVIRONMENT] [EPOCH] [ERRORS] [EXTPATHS] [FILTER] [FORMAT] [FUZZBITS]  
[HANDLES] [HEADER] [LENGTH] [LICENSE] [LOCALE] [MESSAGES] [MEXPAND] [MIOUTPUT]  
[MITERATE] [MNEST] [MPRINT] [MXCELLS] [MXERRS] [MXLOOPS] [MXWARNS]  
[N] [OATTRS] [OLANG] [ONUMBERS] [OVARS] [PRINTBACK] [REPDEFER] [RESULTS] [RNG]  
[SCALEMIN] [SCOMPRESSION] [SEED] [SMALL] [SORT] [SPLIT] [SYSMIS] [TABLERENDER]  
[TFIT] [TLOOK] [TMSRECORDING] [TNUMBERS] [TVARS] [UNDEFINED] [UNICODE] [VERSION]  
[WEIGHT] [WIDTH] [WORKSPACE] [XVERSION] [ZCOMPRESSION] [\$VARS]

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

### Release 13.0

- BLKSIZE and BUFNO subcommands are obsolete and no longer supported.
- HANDLES subcommand introduced.
- SORT subcommand introduced.

### Release 15.0

- TMSRECORDING subcommand introduced.

### Release 16.0

- UNICODE subcommand introduced.
- MCACHE subcommand introduced.
- THREADS subcommand introduced.

### Release 17.0

- FUZZBITS subcommand introduced.

### Release 18.0

- EXTPATHS subcommand introduced.
- ZCOMPRESSION subcommand introduced.
- COMPRESSION subcommand removed because it is obsolete.
- REPDEFER subcommand introduced.

### Release 19.0

- TABLERENDER subcommand introduced.
- XVERSION subcommand introduced.
- OATTRS subcommand introduced.
- DIGITGROUPING subcommand introduced.
- CMPTRANS subcommand introduced.

### **Example**

SHOW .

## **Overview**

SHOW displays current settings for running options. Most of these settings can be changed by using the SET command.

### **Basic Specification**

The basic specification is simply the command keyword, which displays important current settings (keyword ALL). Some displayed option settings are applicable only when you have options such as Tables and Categories.

### **Subcommand Order**

Subcommands can be named in any order.

### **Syntax**

- If any subcommands are specified, only the requested settings are displayed.
- SHOW can be specified more than once.

## **Example**

SHOW BLANKS /UNDEFINED /MXWARNS .

- BLANKS shows the value to which a completely blank field for a numeric variable is translated.
- UNDEFINED indicates whether a message is displayed whenever the program encounters anything other than a number or a blank as the value for a numeric variable.
- MXWARNS displays the maximum number of warnings that are allowed before a session is terminated.

## Subcommands

The following alphabetical list shows the available subcommands.

<b>ALL</b>	<i>Display important settings that are applicable to your system.</i> This setting is the default.
<b>BLANKS</b>	<i>Value to which a completely blank field for a numeric variable is translated.</i> The default is the system-missing value.
<b>BOX</b>	<i>Characters used to draw boxes.</i> Both character and hexadecimal representations are displayed. The default is X'2D7C2B2B2B2B2B2B2B'. This setting applies only to text output from the program.
<b>BLOCK</b>	<i>Character used to draw bar charts.</i> Both character and hexadecimal representations are displayed. The default is X'2A'. This setting applies only to the text output from the program.
<b>CC</b>	<i>Custom currency formats.</i> CC shows the current custom currency formats that have been defined for CCA, CCB, CCC, CCD, and CCE on SET. In Windows environments, the formats reflect the Regional Settings Properties. You can also request any of these keywords individually.
<b>CACHE</b>	<i>Cache active data file.</i> This setting shows the number of changes in the active data file before a cache file is created. The default is 20.
<b>CMPTRANS</b>	<i>Override SPSS Statistics Server setting for compiled transformations.</i> The setting is either ON or OFF (alias YES or NO). There is no default because the default behavior is determined by the IBM® SPSS® Statistics Server.
<b>CTEMPLATE</b>	<i>Chart template file.</i> The setting is either NONE or a filename.
<b>DECIMAL</b>	<i>Decimal indicator.</i> This setting indicates the current character used as the decimal indicator. DOT indicates that a period is the decimal indicator; COMMA indicates that a comma is the decimal indicator.
<b>DEFOLANG</b>	<i>Default output language.</i> This setting indicates the default output language to use if the language that is specified on the OLANG subcommand is not available. The initial default setting is the language version of the installed software.
<b>DIGITGROUPING</b>	<i>Whether the current SPSS Statistics locale's digit grouping format is applied to numeric output.</i> The setting is YES (alias ON) or NO (alias OFF). The default is NO.
<b>DIRECTORY</b>	<i>Default directory.</i> This setting indicates the root directory that is used to determine the locations of files that are specified with no paths or relative paths. A wide variety of actions can change the current default directory during a session.
<b>ENVIRONMENT</b>	<i>Operating system and computer information.</i> This setting includes information about environment variables, defined paths, domain, etc.
<b>EPOCH</b>	<i>Range of years for date-format variables and date functions entered with a two-digit year value.</i> AUTOMATIC indicates a 100-year range beginning 69 years before the current year and 30 years after the current year.
<b>ERRORS</b>	<i>Error messages for text output.</i> The setting can be LISTING (alias YES or ON) or NONE (alias NO or OFF).
<b>EXTENSIONS</b>	No longer supported.
<b>EXTPATHS</b>	<i>Displays the locations for installing and searching for custom dialogs, extension commands, and R packages.</i> When more than one location exists, location 1 takes precedence over location 2, etc. The location R_HOME (in R_HOME\library) specifies the directory where R is installed.
<b>FILTER</b>	<i>Filter status.</i> This setting indicates whether filtering is currently in effect (FILTER command) and indicates the filter variable that is in use (if any).
<b>FORMAT</b>	<i>Default print and write formats for numeric variables that are defined on DATA LIST in freefield format and all numeric variables created by transformation commands.</i> The default is F8.2.

<b>FUZZBITS</b>	<i>Default setting of the fuzzbits argument for the RND and TRUNC functions. The default is 6.</i>
<b>HANDLES</b>	<i>Displays a list of all file handles currently in effect. File handles are created with the FILE HANDLE command.</i>
<b>HEADER</b>	<i>Headings for text output. The setting is YES, NO, or BLANK. The default is NO.</i>
<b>JOURNAL</b>	No longer supported.
<b>LENGTH</b>	<i>Maximum page length for output. The default is 59. This setting applies only to the text output from the program.</i>
<b>LICENSE</b>	<i>Licensed components, expiration date, release number, and maximum number of users permitted by the license.</i>
<b>LOCALE</b>	<i>Operating system locale setting and codepage. In Windows operating systems, locale is set in the Regional Options of the Control Panel.</i>
<b>MCACHE</b>	<i>Memory cache size in kilobytes. The default and recommended setting are based on your specific hardware.</i>
<b>MESSAGES</b>	<i>Resource-utilization messages for text output. The setting can be LISTING (alias YES or ON) or NONE (alias NO or OFF).</i>
<b>MEXPAND</b>	<i>Macro expansion. The setting is either ON (alias YES) or OFF (alias NO). The default is ON.</i>
<b>MIOUTPUT</b>	<i>Displays the settings for pooled output when a multiply imputed dataset is analyzed.</i>
<b>MITERATE</b>	<i>Maximum loop iterations permitted in macro expansions. The default is 1000.</i>
<b>MNEST</b>	<i>Maximum nesting level for macros. The default is 50.</i>
<b>MPRINT</b>	<i>Inclusion of expanded macros in the output. The setting is either ON (alias YES) or OFF (alias NO). The default is OFF.</i>
<b>MXCELLS</b>	<i>Maximum number of cells that are allowed for a new pivot table. The default is AUTOMATIC, which allows the number to be determined by the available memory.</i>
<b>MXERRS</b>	<i>Maximum number of errors allowed and number of errors so far in current session. In most implementations of SPSS Statistics, the maximum number of errors defined on SET MXERRS is ignored. However, the information that SHOW MXERRS provides about number of errors in the current session can be useful.</i>
<b>MXLOOPS</b>	<i>Maximum executions of a loop on a single case. The default is 40.</i>
<b>MXMEMORY</b>	No longer supported.
<b>MXWARNS</b>	<i>Maximum number of warnings and errors that are shown for text output. The default is 10.</i>
<b>N</b>	<i>Unweighted number of cases in the active dataset. N displays UNKNOWN if a active dataset has not yet been created. N cannot be changed with SET.</i>
<b>OATTRS</b>	<i>Whether OXML and SAV output from OMS contains English-only attribute values and English-only variable names respectively. The setting is OLANG, ENG, or BOTH. The default is OLANG.</i>
<b>OLANG</b>	<i>Output language for pivot tables.</i>
<b>ONUMBERS</b>	<i>Display of variable values in the outline for pivot tables. The settings can be LABELS, VALUES, and BOTH.</i>
<b>OVARS</b>	<i>Display of variables as headings. The settings can be LABELS, NAMES, and BOTH.</i>
<b>PRINTBACK</b>	<i>Command printback. The setting can be BOTH (alias LISTING, YES, or ON) or NONE (alias NO or OFF). The default is BOTH at system installation.</i>
<b>REPDEFER</b>	<i>Whether connection requests to an IBM® SPSS® Collaboration and Deployment Services Repository are deferred until needed. The setting is YES (alias ON) or NO (alias OFF). The default is NO unless executing command syntax with the SPSS Statistics Batch Facility (available with SPSS Statistics Server), in which case the default is YES.</i>



<b>RESULTS</b>	<i>Output from commands.</i> This setting is not applicable to output that is displayed in pivot tables. The setting can be LISTING (alias YES or ON) or NONE (alias NO or OFF).
<b>RNG</b>	<i>Random number generator.</i> Indicates which random number generator is currently being used.
<b>SCALEMIN</b>	<i>For data files that were created in versions prior to version 8.0, the minimum number of unique values for a numeric variable that is used to classify the variable as scale.</i> This setting affects only pre-8.0 data files that are opened in later versions.
<b>SCOMPRESSION</b>	<i>Compression of SPSS Statistics data files.</i> This setting can be overridden by the COMPRESSED or UNCOMPRESSED subcommands on the SAVE or XSAVE commands. The default setting varies by system. SCOMPRESSION cannot be changed with SET.
<b>SEED</b>	<i>Seed for the random-number generator.</i> The default is generally 2,000,000 but may vary by system.
<b>SMALL</b>	<i>Decimal value to control display of scientific notation in output.</i>
<b>SORT</b>	<i>Sorting mechanism that is currently in effect: SPSS Statistics or external, third-party (if available).</i>
<b>SPLIT</b>	<i>Split file status.</i>
<b>SYSMIS</b>	<i>The system-missing value.</i> SYSMIS cannot be changed with SET.
<b>TABLERENDER</b>	<i>Method for rendering tables.</i> The settings can be FULL or LIGHT. Lightweight tables render faster but cannot be pivoted or edited.
<b>TFIT</b>	<i>Adjust column widths in pivot tables.</i> The settings can be BOTH (label and data) and LABELS.
<b>THREADS</b>	<i>Number of threads.</i> The default and recommended setting are based on your specific hardware.
<b>TLOOK</b>	<i>Pivot table template file.</i> The setting can be either NONE or a filename.
<b>TMSRECORDING</b>	<i>TMS recording status.</i> YES indicates that a TMS block for writing transformations to PMML is currently in effect. NO indicates that TMS is not currently in effect. <a href="#">For more information, see the topic TMS BEGIN on p. 1965.</a>
<b>TNUMBERS</b>	<i>Display of variable values in pivot tables.</i> The settings can be VALUES, LABELS, and BOTH.
<b>TVARS</b>	<i>Display of variables as headings.</i> The settings can be NAMES, LABELS, and BOTH.
<b>UNDEFINED</b>	<i>Warning message for undefined data.</i> YES is the value associated with the default setting of SET UNDEFINED=WARN.NO is the value associated with SET UNDEFINED=NOWARN, which suppresses messages but does not alter the count of warnings toward the MXWARNS total.
<b>UNICODE</b>	<i>Unicode mode status.</i> The setting is either YES (alias ON) or NO (alias OFF).
<b>VERSION</b>	<i>Version number and creation date.</i>
<b>WEIGHT</b>	<i>Variable that is used to weight cases.</i> WEIGHT can be specified for SHOW only and cannot be changed with SET.
<b>WIDTH</b>	<i>Maximum page width for the output.</i> The default is 132 columns for batch mode and 80 for interactive mode. This setting applies only to text output from the program.
<b>WORKSPACE</b>	<i>Special workspace memory limit in kilobytes.</i> The default is 6148.
<b>XVERSION</b>	<i>Version of Output XML schema used to generate OXML output from OMS.</i> The setting is DEFAULT or the version number (if the version differs from the default one). The default is the value DEFAULT.

---

*SHOW*

**ZCOMPRESSION**      *Override Server setting for compression of scratch files. The setting is either ON or OFF (alias YES or NO). There is no default because the default behavior is determined by the SPSS Statistics Server.*

**\$VARS**                *Values of system variables. \$VARS cannot be changed with SET.*

# ***SORT CASES***

```
SORT CASES [BY] varlist[({A})] [varlist...]  
           {D}
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## ***Release History***

Release 17.0

- Added support for SET THREADS.

## ***Example***

```
SORT CASES BY DIVISION (A) STORE (D) .
```

## ***Overview***

SORT CASES reorders the sequence of cases in the active dataset based on the values of one or more variables. You can optionally sort cases in ascending or descending order, or you can use combinations of ascending and descending order for different variables.

## ***Basic Specification***

The basic specification is a variable or list of variables that are used as sort keys. By default, cases are sorted in ascending order of each variable, starting with the first variable that is named. For each subsequent variable, cases are sorted in ascending order within categories of the previously named variables.

## ***Syntax Rules***

- Keyword BY is optional.
- BY variables can be numeric or string but not scratch, system, or temporary variables.
- You can explicitly request the default sort order (ascending) by specifying A or UP in parentheses after the variable name. To sort cases in descending order, specify D or DOWN.
- An order specification (A or D) applies to all variables in the list, up to the previous order specification. If you combine ascending and descending order on the same SORT CASES command, you may need to specify the default A explicitly.

## ***Operations***

- SORT CASES first sorts the file according to the first variable that is named. For subsequent variables, cases are sorted within categories of the previously named variables.

- The sort sequence is based on the locale-defined order (and is not necessarily the same as the numerical order of the character codes). The default locale is the operating system locale. You can change the locale with [SET LOCALE](#). Use `SHOW LOCALE` to display the current locale.
- This procedure uses the multithreaded options specified by `SET THREADS`.

#### **Limitations**

- You cannot specify more than 64 sort variables.

## **Examples**

```
SORT CASES BY DIVISION (A) STORE (D) .
```

- Cases are sorted in ascending order of variable *DIVISION*. Cases are further sorted in descending order of *STORE* within categories of *DIVISION*. `A` must be specified so that `D` applies to *STORE* only.

```
SORT CASES DIVISION STORE (A) AGE (D) .
```

- Cases are sorted in ascending order of *DIVISION*. Keyword `BY` is not used in this example.
- Cases are further sorted in ascending order of *STORE* within values of *DIVISION*. Specification `A` applies to both *DIVISION* and *STORE*.
- Cases are further sorted in descending order of *AGE* within values of *STORE* and *DIVISION*.

## ***SORT CASES with Other Procedures***

- In `AGGREGATE`, cases are sorted in order of the break variable or variables. You do not have to use `SORT CASES` prior to running `AGGREGATE`, because the procedure does its own sorting.
- You can use `SORT CASES` in conjunction with the `BY` keyword in `ADD FILES` to interleave cases with the same variables but from different files.
- With `MATCH FILES`, cases must be sorted in the same order for all files that you combine.
- With `UPDATE`, cases must be sorted in ascending order of the key variable or variables in both the master file and all transaction files.
- You can use the `PRINT` command to check the results of a `SORT CASES` command. To be executed, `PRINT` must be followed by a procedure or `EXECUTE`.

# ***SORT VARIABLES***

```
SORT VARIABLES [BY] {NAME          } [({A})]
                   {TYPE           }      {D}
                   {FORMAT         }
                   {LABEL          }
                   {VALUES         }
                   {MISSING        }
                   {MEASURE        }
                   {ROLE           }
                   {COLUMNS       }
                   {ALIGNMENT      }
                   {ATTRIBUTE name}
```

If there are any pending transformations, this command reads the active dataset and causes execution of any pending commands. Otherwise, it takes effect immediately without reading the active dataset.

## ***Release History***

Release 16.0.

- Command introduced.

Release 18.0.

- ROLE keyword introduced.

## ***Example***

```
SORT VARIABLES BY NAME (A) .
```

## ***Overview***

SORT VARIABLES sorts the variables in the active dataset based on the values of the selected dictionary attribute.

### ***Basic Specification***

The basic specification is the command name SORT VARIABLES followed by the name of a single dictionary attribute.

### ***Syntax Rules***

- Only one dictionary attribute can be specified.
- The BY keyword is optional.
- You can explicitly request the default sort order (ascending) by specifying A or UP in parentheses after the attribute name. To sort variables in descending order, specify D or DOWN.

**Sorting Options**

Variables can be sorted based on the values of any one of the following dictionary attributes:

**NAME.** *Sort variables by variable names.* The primary sort is alphabetical, but trailing digits are sorted numerically within ties on the preceding part. For example, *V2* comes before *V10* in ascending order.

**TYPE.** *Group variables by type (numeric or string) and sort string variables by defined width.* In default ascending order, numeric variables come before string variables, and shorter string variables come before longer string variables.

**FORMAT.** *Group variables by format (for example, Date, Dollar, String).*

**LABEL.** *Sort variables in alphabetical order by variable labels.* In default ascending order, all variables without defined variable labels come before variables with defined variable labels. [For more information, see the topic VARIABLE LABELS on p. 2130.](#)

**VALUES.** *Sort variables by value labels.* In default ascending order, all variables without defined value labels come before variables with defined value labels. [For more information, see the topic VALUE LABELS on p. 2116.](#)

**MISSING.** *Group variables by defined missing values.* In default ascending order, all variables without defined missing values come before variables with defined missing values. [For more information, see the topic MISSING VALUES on p. 1197.](#)

**MEASURE.** *Sort variables by measurement level.* The default ascending order is: nominal, ordinal, scale. [For more information, see the topic VARIABLE LEVEL on p. 2132.](#)

**ROLE.** *Sort variables by role.* Some dialogs support predefined roles that can be used to pre-select variables for analysis. [For more information, see the topic Overview on p. 2133.](#)

**COLUMNS.** *Sort variables by column width.* Column width is the width of the column that displays the variable in Data View of the Data Editor. [For more information, see the topic VARIABLE WIDTH on p. 2135.](#)

**ALIGNMENT.** *Group variables by alignment.* The default ascending order is: left, right, center. [For more information, see the topic VARIABLE ALIGNMENT on p. 2126.](#)

**ATTRIBUTE name.** *Sort variables by values of the specified custom variable attribute name.* In default ascending order, all variables without values for the named attribute come before variables with values for the named attribute. [For more information, see the topic VARIABLE ATTRIBUTE on p. 2127.](#)

# SPCHART

SPCHART is available in the Statistics Base option.

```
SPCHART
  [/TEMPLATE='filename']

  [/TITLE='line 1' ['line 2']]
  [/SUBTITLE='line 1']
  [/FOOTNOTE='line 1' ['line 2']]

  {[/XR={var BY var          } ] }
    {var var [var var...][BY var]} [(XBARONLY)]
  { /XS= {var BY var          } ] }
    {var var [var var...][BY var]} [(XBARONLY)]
  { /IR= var [BY var]          } }
  { /I= var [BY var]          } }
  { /NP= {var BY var          } }
    {COUNT(var) N({var }) [BY var]}
    {value }
  { /P= {var BY var          } }
    {COUNT(var) N({var }) [BY var]}
    {value }
  { /C= {var BY var          } }
    {COUNT(var) N({var }) [BY var]}
    {value }
  { /U= {var BY var          } }
    {COUNT(var) N({var }) [BY var]}
    {value }

  { /CPCHART= {var BY var    } }
    {var var [var...]}
    {var }

  [/STATISTICS = [CP] [CPL] [CPU] [K] [CPK] [CR] [CPM]
    [CZL] [CZU] [CZMIN] [CZMAX] [CZOUT]
    [PP] [PPL] [PPU] [PPK] [PR] [PPM]
    [PZL] [PZU] [PZMIN] [PZMAX] [PZOUT]
    [AZOUT] [N] [MEAN] [STDDEV] [CAPSIGMA]
    [LSL] [USL] [TARGET] [AZLOUT] [AZUOUT]
    [CZLOUT] [CZUOUT] [PZLOUT] [PZUOUT] ]

  [/RULES = [ALL] [UCL] [R2UPPER] [R4UPPER] [R8UPPER]
    [R8LOWER] [R4LOWER] [R2LOWER] [LCL] [TRUP] [TRDOWN] [ALTERNATING] ]

  [/ID=var]

  [/CAPSIGMA = [{RBAR }]]
    {SBAR }
    {MRBAR }
    {WITHIN}

  [/SPAN={2**}]
    {n }

  [{/CONFORM }=value]
  {/NONCONFORM}

  [/SIGMAS={3**}]
    {n }

  [/MINSAMPLE={2**}]
    {n }

  [/LSL=value]    [/USL=value]

  [TARGET = value]

  [/MISSING={NOREPORT**} [{EXCLUDE**} ]
    {REPORT } {INCLUDE } ]
```

```

[/NORMAL = { CAPABILITY** } ]
           { PERFORMANCE }
           { BOTH         }
           { NONE         }

[/REFERENCE = { [SL**] [SIGMAS] } ]
             { NONE         }

```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

#### Release 15.0

- (XBARONLY) keyword introduced on XR and XS subcommands.
- RULES subcommand introduced.
- ID subcommand introduced.

#### Release 19.0

- CPCHART subcommand introduced.
- NORMAL subcommand introduced.
- REFERENCE subcommand introduced.
- Following keywords introduced on STATISTICS subcommand: N, MEAN, STDDEV, CAPSIGMA, LSL, USL, TARGET, AZLOUT, AZUOUT, CZLOUT, CZUOUT, PZLOUT, PZUOUT.

### **Example**

```

SPCHART /TEMPLATE='cntr.sgt'
        /IR=subsize.

```

## **Overview**

SPCHART generates several types of high-resolution control charts. A control chart plots a quality characteristic that is measured or computed from a sample versus the sample number or time. This technique is a widely used process-control technique for testing the hypothesis that the process is in a state of statistical control. All control charts display four series:

- The process line representing the quality characteristic for each sample.
- The center line indicating the average value of the quality characteristic that corresponds to the in-control state.
- Two horizontal lines showing the upper control limit and lower control limit.

Control charts are used for improving productivity, preventing defects and unnecessary process adjustments, and gathering information about process capability.



SPCHART produces X-bar, R, s, individuals, and moving range charts as well as np, p, c, and u charts. It also produces a process capability chart. You may need to transform your data to conform to the required data organization that is described under each chart type subcommand.

Control charts are available only on systems where high-resolution display is available.

### **Options**

**Titles and Footnotes.** You can use the `TITLE`, `SUBTITLE`, and `FOOTNOTE` subcommands to specify a title, subtitle, and footnote for the chart.

**Chart Type.** You can request a specific type of chart by using the `XR`, `XS`, `IR`, `I`, `NP`, `P`, `C`, `U`, or `CPCHART` subcommand.

**Templates.** You can specify a template, using the `TEMPLATE` subcommand, to override the default chart attribute settings on your system.

**Control Limits.** You can specify a sigma value on the `SIGMAS` subcommand to modify the calculated upper and lower control limits. You can also use the `USL` and `LSL` subcommands to specify fixed limits. The upper and lower limits that you specify will be displayed simultaneously with the calculated control limits. *Note:* Calculated control limits are not displayed on process capability charts.

**Control Rules.** You can specify control rules that help you quickly identify out-of-control points.

### **Basic Specification**

The basic specification is a chart type subcommand with a variable specification that describes the process data.

### **Subcommand Order**

Subcommands can be specified in any order.

### **Syntax Rules**

- Only one chart type subcommand can be specified.
- Keyword `SPAN` is used only with `IR` and `I` subcommands.
- Keyword `CONFORM` or `NONCONFORM` is used only with `NP` and `P` subcommands.
- Keywords `NORMAL` and `REFERENCE` are used only with the `CPCHART` subcommand.
- Keywords `LSL` and `USL` are required when the `CPCHART` subcommand is specified.

### **Operations**

- All control charts generated by SPCHART contain four basic series: the process, the center line, the upper control line, and the lower control line.
- The chart title, subtitle, and footnote are assigned as they are specified on `TITLE`, `SUBTITLE`, and `FOOTNOTE` subcommands. For control charts, the default chart title is Control Chart, followed by the label of the process variable, and a one-line default footnote displays the sigma level. For a process capability chart, the default chart title is Process Capability and

there is no default footnote. For all chart types, if split-file processing is in effect then a default subtitle is added that provides split-file information.

- For control charts, the category variable label is used as the title for the *category axis*. If no variable label is defined, the variable name is used. If no category variable is defined, the title is null.
- For control charts, the category variable value labels are used as the *category axis* labels. If no value labels are defined, values are used. If no category variable is defined, integer values from 1 to  $n$  are used, where  $n$  is the number of subgroups or units plotted.
- For control charts, all series are plotted as lines. When a series has a constant value across all samples, the value is reported in the legend entry for the series.
- For all chart types, case weights are not honored when each case is a subgroup. Case weights are honored when each case is a unit and when the weights are integers. When weighted data are used in an individuals chart, replicated cases are plotted on the control chart.
- For control charts, the calculated control limits are always displayed and can be suppressed only by editing the chart in a chart window.
- You can specify preset control limits for an X-bar or I chart, as some industries often do. The specified control limits are displayed simultaneously with the calculated limits.

### **Limitations**

- Control charts cannot have fewer than 2 points or more than 3000 points.
- The subgroup size in X-bar and range charts cannot exceed 100.
- The span for individuals charts is limited to 100.
- For process capability charts, the subgroup size cannot exceed 100 when /CAPSIGMA=RBAR, and the span cannot exceed 100 when /CAPSIGMA=MRBAR. In addition, the subgroup size or span cannot be less than 2.

## **Example**

```
SPCHART /TEMPLATE='cntr.sgt'
/IR=subsize.
```

- This command generates an individuals chart and a moving range chart. The process variable *subsize* is a numeric variable that measures the size variation of the product.
- Both charts use the attributes that are defined for the template that is saved in *cntr.sgt*.
- The default span (2) and sigma value (3) are used.
- Because no BY variable is specified, the  $x$  axis is labeled by sequence numbers.

### **X-bar and R-charts**

```
SPCHART
/XR=ph BY time
/STATISTICS=CP CPU CPL K CPM CZOUT PP PPU PPL PPM PZOUT AZOUT
/RULES=All
/CAPSIGMA=RBAR
/SIGMAS=3
/USL=5.5
```

```
/LSL=4.5
/TARGET=5.0
/MINSAMPLE=2.
```

- The procedure produces an x-bar and R chart for *ph* by *time*.
- The `RULES` subcommand specifies that all rules for identifying out-of-control points should be used.
- The `STATISTICS` subcommand requests various process capability and process performance indices, in addition to measures for assessing normality.
- The `USL` and `LSL` subcommands specify upper and lower control limits of 5.5 and 4.5, respectively.
- The `TARGET` subcommand defines the target value that is used in computing CpM and PpM.
- All other options are set to their default values.

### ***Descriptive Heading for First Example***

```
SPCHART
/P=COUNT( defects ) N( sampled )BY lot
/RULES=ALL
/SIGMAS=3.
```

- The procedure produces a p chart using the variable *lot* to define subgroups, *sampled* to define the number of items sampled per lot, and *defects* to define the number of defective units per sample.
- The `RULES` subcommand specifies that all rules for identifying out-of-control points should be used.
- All other options are set to their default values.

## ***TEMPLATE Subcommand***

`TEMPLATE` uses an existing file as a template and applies it to all charts generated by the current `SPCHART` command. The template overrides the default settings that are used to create any charts, and the specifications on the current `SPCHART` command override the template. Templates are created in the Chart Editor by saving an existing chart as a template.

### ***Example***

```
SPCHART /TEMPLATE='/templates/cntr.sgt'
/IR=subsize.
```

## ***TITLE, SUBTITLE, and FOOTNOTE Subcommands***

`TITLE`, `SUBTITLE`, and `FOOTNOTE` specify lines of text that are placed at the top or bottom of the control chart.

- One or two lines of text can be specified for `TITLE` or `FOOTNOTE`, and one line of text can be specified for `SUBTITLE`.
- Each line of text must be enclosed in quotes. The maximum length of any line is 72 characters.

- The default font sizes and types are used for the title, subtitle, and footnote.
- By default, the title, subtitle, and footnote are left-aligned with the y axis.
- For control charts, if you do not specify `TITLE` the default is Control Chart followed by the label of the process variable. For a process capability chart the default title is Process Capability.
- If you do not specify `SUBTITLE`, the subtitle provides the split-file information if split-file processing is in effect; otherwise, it is null, which leaves more space for the chart.
- For control charts, if you do not specify `FOOTNOTE` the sigma level is identified as the first line of the footnote. There is no default footnote for a process capability chart.

### Example

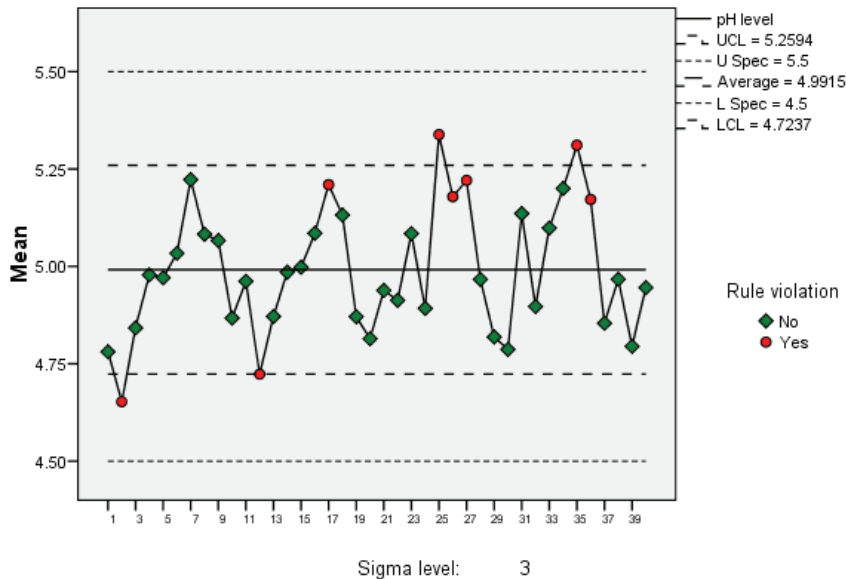
```
SPCHART TITLE = 'Wheel Production'
/SUBTITLE = 'Process Control'
/IR=SUBSIZE.
```

## XR and XS Subcommands

XR produces an X-bar chart and an R chart. XS produces an X-bar chart and an s chart. X-bar, R, and s charts are control charts for continuous variables, such as size, weight, length, and temperature.

An X-bar chart plots the mean of each subgroup. The center line indicates the mean of subgroup means. The control limits are calculated from subgroup means, numbers, standard deviations, and the user-specified `SIGMAS` value. The following figure shows an X-bar chart.

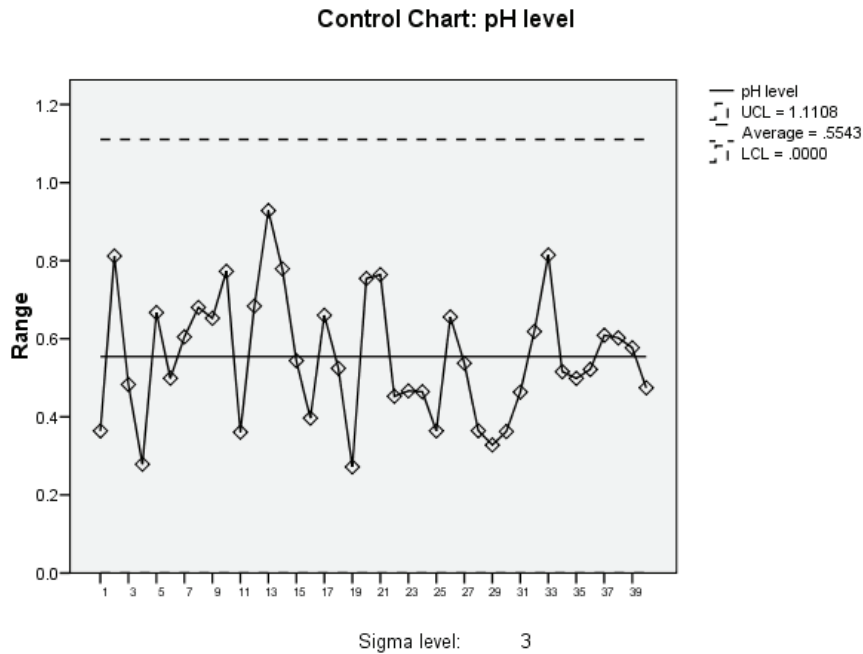
Figure 230-1  
X-bar chart



An R chart plots range values (maximum-minimum) of successive subgroups. The center line indicates the mean of subgroup ranges. The control limits are calculated from subgroup ranges, numbers, and the user-specified `SIGMAS` value. The R chart tests whether the process variability

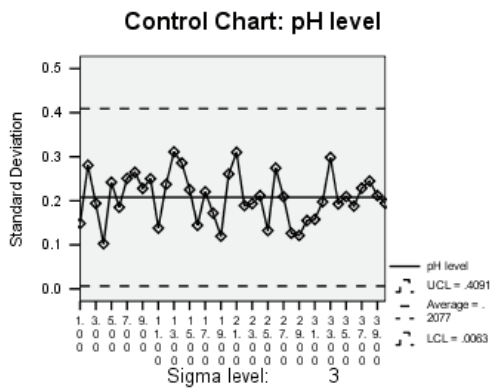
is in control. When the subgroup size is relatively small (4, 5, or 6), the range method yields almost as good an estimator of the variance as does the subgroup variance. The following figure shows an R chart.

Figure 230-2  
R chart



An s chart plots subgroup standard deviations. The center line indicates the mean of subgroup standard deviations. The control limits are calculated from subgroup standard deviations, numbers, and the user-specified SIGMAS value. The s chart tests whether the process variability is in control, especially when the subgroup size is moderate to large. The following figure shows an s chart.

Figure 230-3  
An s chart



## Data Organization

For X-bar, R, or s charts, data can be organized where each case is a unit or where each case is a subgroup.

- If each case is a unit with a subgroup identifier, cases are assigned to a category according to the value of the identifier. Table 230-1 is an example of this type of data organization. The data do not have to be sorted by subgroup. A BY variable (the subgroup identifier) is required to sort and aggregate data and label the process variable.
- If each case is a subgroup, there are as many variables as individuals within one sample. A sample identifier is not required. When there is a sample identifier, it is used for labeling. Table 230-2 shows this type of organization.

Table 230-1

*Each case is a unit for X-bar, R, and s charts*

Subgroup	Length
8:50	6.35
11:30	6.39
8:50	6.40
11:30	6.46
8:50	6.32
11:30	6.37
8:50	6.39
11:30	6.36
...	...

Table 230-2

*Each case is a subgroup for X-bar, R, and s charts*

Subgroup	N1	N2	N3	N4
8:50	6.35	6.40	6.32	6.39
11:30	6.39	6.46	6.37	6.36
...	...	...	...	...

## Variable Specification

If data are organized as shown in Table 230-1, the variable specifications on XR and XS subcommands are

```
VAR BY VAR
```

The variable that is specified before BY is the process variable, which is the variable that contains values for all instances to be plotted (for example, *LENGTH* in Table 230-1). The variable that is specified after BY is the category variable or the BY variable, which is the subgroup identifier (for example, *SUBGROUP* in Table 230-1). The process variable must be numeric, while the category variable can be of any type. The chart is sorted by the category variable.

If data are organized as shown in Table 230-2, the variable specifications on XR and XS subcommands are

```
VAR VAR [VAR...] [BY VAR]
```

Each of the variables that is specified before `BY` represents an instance to be plotted (for example, *N1* to *N3* in Table 230-2). At least two variables are required, and each variable must be numeric. Keyword `BY` and the category variable (for example, *SUBGROUP* in Table 230-2) are optional; if specified, the category variable provides labels for the *category axis* and can be any type of variable. If omitted, the *category axis* is labeled from 1 to the number of variables that are specified before keyword `BY`.

### Example

```
SPCHART /TEMPLATE='ctrl.sgt'
/XR SUBSIZE BY SHIFT.
```

- The data are organized as shown in Table 230-1. *SUBSIZE* is a numeric variable that measures the part size. *SHIFT* contains the subgroup identifier (work shift number).
- The chart template is stored in the chart file *ctrl.sgt*.

## (XBARONLY) Keyword

(`XBARONLY`) suppresses the R or s secondary charts. If this keyword is omitted, the R or s chart will be generated with the X-bar chart.

### Example

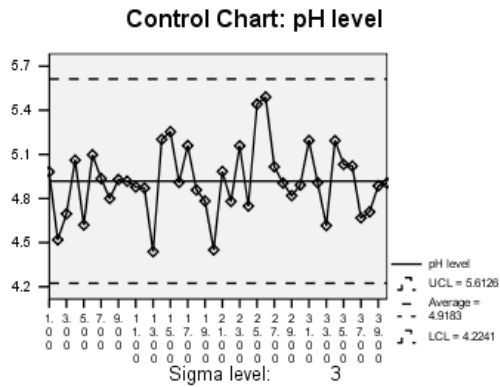
```
SPCHART
/XR=ph BY time (XBARONLY)
/CAPSIGMA=RBAR
/SIGMAS=3
/MINSAMPLE=2.
```

## I and IR Subcommands

`I` produces an individuals chart, and `IR` produces an individuals chart and a moving range chart. Both types are control charts for continuous variables, such as size, weight, length, and temperature.

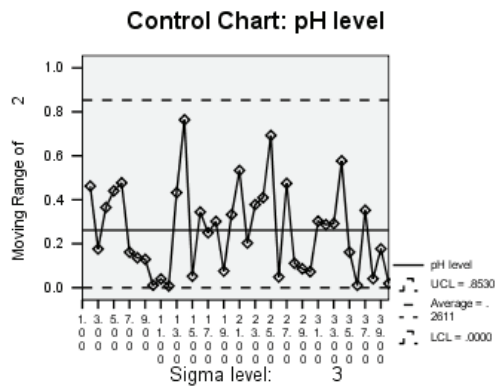
An individuals chart plots each individual observation on a control chart. The center line indicates the mean of all individual values, and the control limits are calculated from the mean of the moving ranges, the span, and the user-specified `SIGMAS` value. Individuals charts are often used with moving range charts to test process variability when the subgroup size is 1. This situation occurs frequently when automated inspection and measurement technology is used and every manufactured unit is analyzed. The situation also occurs when the process is so slow that a larger subgroup size becomes impractical. The following figure shows an individuals chart.

Figure 230-4  
Individuals chart



A moving range chart plots moving ranges of  $n$  successive observations on a control chart, where  $n$  is the specified span (see SPAN Subcommand on p. 1919). The center line is the mean of moving ranges, and the control limits are calculated from the ranges, the span, and the user-specified SIGMAS value (see SIGMAS Subcommand on p. 1919). The following figure shows a moving range chart.

Figure 230-5  
Moving range chart



## Data Organization

For individuals charts and moving range charts, data must be organized so that each case is a unit. Cases are not sorted or aggregated before plotting.

## Variable Specification

The variable specification for I or IR subcommand is

VAR [BY VAR]



You must specify the process variable that contains the value for each individual observation. Each observation is plotted for the individuals chart. The range of  $n$  consecutive observations (where  $n$  is the value that is specified on the `SPAN` subcommand) is calculated and plotted for the moving range chart. The range data for the first  $n-1$  cases are missing, but the mean and the limit series are not missing.

Keyword `BY` and the category variable are optional. When specified, the category variable is used for labeling the *category axis* and can be any type of variable. If omitted, the *category axis* is labeled 1 to the number of individual observations in the process variable.

### **Example**

```
SPCHART /TEMPLATE='ctrl.sgt'  
/IR=SUBSIZE.
```

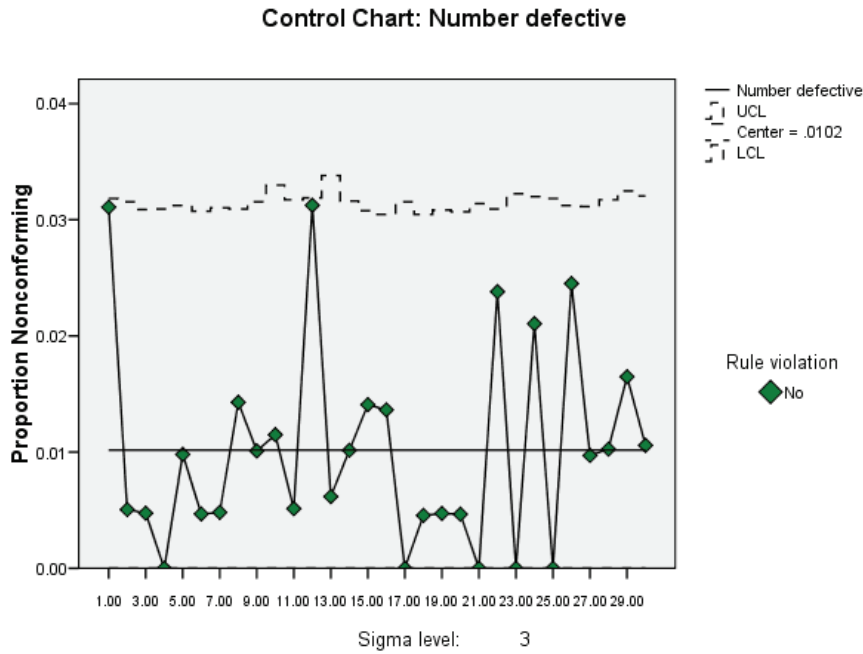
- This command requests an individuals chart and a moving range chart.
- The default span (2) and sigma value (3) are used.

## ***P and NP Subcommands***

`P` produces a p chart and `NP` produces an np chart. Both charts are control charts for attributes. That is, these charts use data that can be counted, such as the number of nonconformities and the percentage of defects.

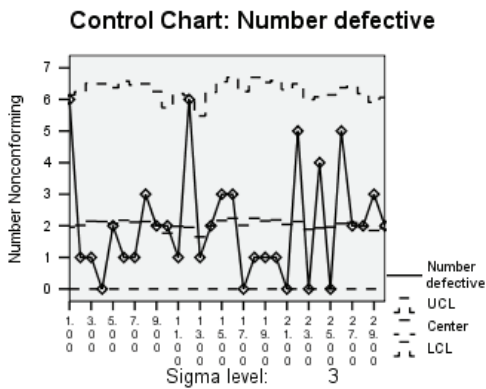
A p chart plots the **fraction nonconforming** on a control chart. Fraction nonconforming is the proportion of nonconforming or defective items in a subgroup to the total number of items in that subgroup. This measurement is expressed as a decimal or, occasionally, as a percentage. The center line of the control chart is the mean of the subgroup fractions, and the control limits are based on a binomial distribution and can be controlled by the user-specified `SIGMAS` value.

Figure 230-6  
A p chart



An np chart plots the number nonconforming rather than the fraction nonconforming. The center line is the mean of the numbers of nonconforming or defective items. The control limits are based on the binomial distribution and can be controlled by the user-specified SIGMAS value. When the subgroup sizes are unequal, np charts are not recommended.

Figure 230-7  
A np chart



**Data Organization**

Data for p and np charts can be organized where each case is a unit or where each case is a subgroup.

- If each case is a unit with a conformity status variable and a subgroup identifier, cases are assigned to a category by the value of the subgroup identifier. Table 230-3 is an example of this type of data organization. The data do not have to be sorted. A `BY` variable (the subgroup identifier) is required to sort and aggregate data and label the *category axis*.
- If each case is a subgroup, one variable contains the total number of items within a subgroup, and one variable contains the total number of nonconforming or defective items in the subgroup. The subgroup identifier is optional. If specified, the subgroup identifier is used for labeling purposes. Table 230-4 is an example of this type of data organization. The data are the same as the data that are used in Table 230-3.

Table 230-3

Each case is a unit for *p* and *np* charts

Subgroup	Outcome
January	Cured
January	Cured
January	Cured
January	Relapse
February	Relapse
February	Cured
February	Relapse
February	Relapse
...	...

Table 230-4

Each case is a subgroup for *p* and *np* charts

Subgroup	Relapse	N
January	1	4
February	3	4
...	...	...

## Variable Specification

If data are organized as illustrated in Table 230-3, the variable specification on `P` or `NP` subcommands is

```
VAR BY VAR
```

The variable that is specified before `BY` is the status variable (for example, *OUTCOME* in Table 230-3). The value of this variable determines whether an item is considered conforming or nonconforming. The status variable can be any type, but if it is a string, the value that is specified on `CONFORM` (or `NONCONFORM`) must be enclosed in quotes (see `CONFORM` and `NONCONFORM` Subcommands on p. 1919). The variable that is specified after `BY` is the category variable and can be any type of variable. The chart is sorted by values of the category variable.

If data are organized as shown in Table 230-4, the variable specification on `P` or `NP` is

```
COUNT (VAR) N ({VAR}) [BY VAR]
                {VAL}
```

The variable that is specified on keyword `COUNT` is the variable that contains the number of nonconforming or defective items (for example, *RELAPSE* in Table 230-4). The specification on keyword `N` is either the variable that contains the sample size or a positive integer for a constant size across samples (for example, *N* in Table 230-4). The `COUNT` variable cannot be larger than the `N` variable for any given subgroup; if it is larger, the subgroup is dropped from calculation and plotting. Keyword `BY` and the category variable are optional. When specified, the category variable is used for *category axis* labels; otherwise, the *category axis* is labeled 1 to the number of subgroups. Cases are unsorted for the control chart.

## C and U Subcommands

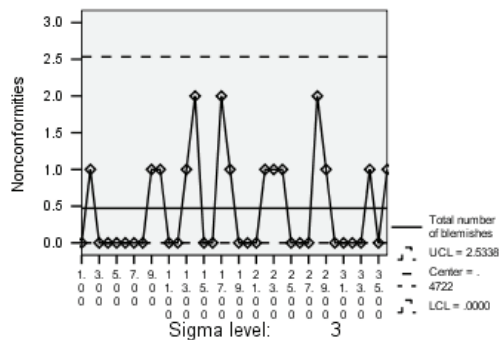
`C` produces a *c* chart and `U` produces a *u* chart. Both charts are control charts for attributes. That is, the charts use data that can be counted.

A *c* chart plots the total number of defects or nonconformities in each subgroup. A defect or nonconformity is one specification that an item fails to satisfy. Each nonconforming item has at least one defect, but any nonconforming item may have more than one defect. The center line of the *c* chart indicates the mean of the defect numbers of all subgroups. The control limits are based on the Poisson distribution and can be controlled by the user-specified `SIGMAS` value. When the sample sizes are not equal, *c* charts are not recommended.

Figure 230-8

A *c* chart

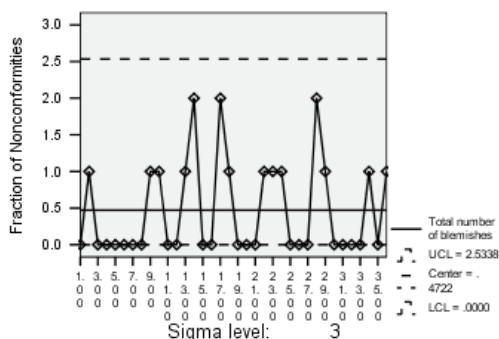
Control Chart: Total number of blemishes



A *u* chart plots the average number of defects or nonconformities per inspection unit within a subgroup. Each subgroup contains more than one inspection unit. The center line of the *u* chart indicates the average number of defects per unit of all subgroups. The control limits are based on Poisson distribution and can be controlled by the user-specified `SIGMAS` value.

Figure 230-9  
A u chart

Control Chart: Total number of blemishes



## Data Organization

Data for c and u charts can be organized where each case is a unit or where each case is a subgroup.

- If each case is a unit with a variable containing the number of defects for that unit and a subgroup identifier, cases are assigned to each subgroup by the value of the identifier. Table 230-5 is an example of this type of data organization. Data do not have to be sorted by subgroup. A BY variable (the subgroup identifier) is required to sort and aggregate data and to label the *category axis*.
- If each case is a subgroup, one variable contains the total number of units within the subgroup, and one variable contains the total number of defects for all units within the subgroup. The subgroup identifier is optional. When specified, the subgroup identifier is used as *category axis* labels; otherwise, the number 1 to the number of subgroups are used to label the *category axis*. Table 230-6 is an example of this method of data organization. The data are the same as the data in Table 230-5.

Table 230-5

Each case is a unit for c and u charts

ID	Subgroup	Count
1	January	0
2	January	2
3	January	0
4	January	0
5	February	5
6	February	1
7	February	0
8	February	0
...	...	...

Table 230-6

Each case is a subgroup for c and u charts

Subgroup	Relapses	N
JANUARY	1	4

Subgroup	Relapses	N
FEBRUARY	3	4
...	...	...

### **Variable Specification**

If data are organized as shown in Table 230-5, the variable specification on C and U subcommands is

```
VAR BY VAR
```

The variable that is specified before keyword BY contains the number of defects in each unit (for example, *COUNT* in Table 230-5). The variable must be numeric. The variable that is specified after keyword BY is the subgroup identifier (for example, *SUBGROUP* in Table 230-5). This variable can be any type of variable. The chart is sorted by values of the subgroup identifier.

If data are organized as shown in Table 230-6, the variable specification on C and U subcommands is

```
COUNT (VAR) N({VAR}) [BY VAR]
                {VAL}
```

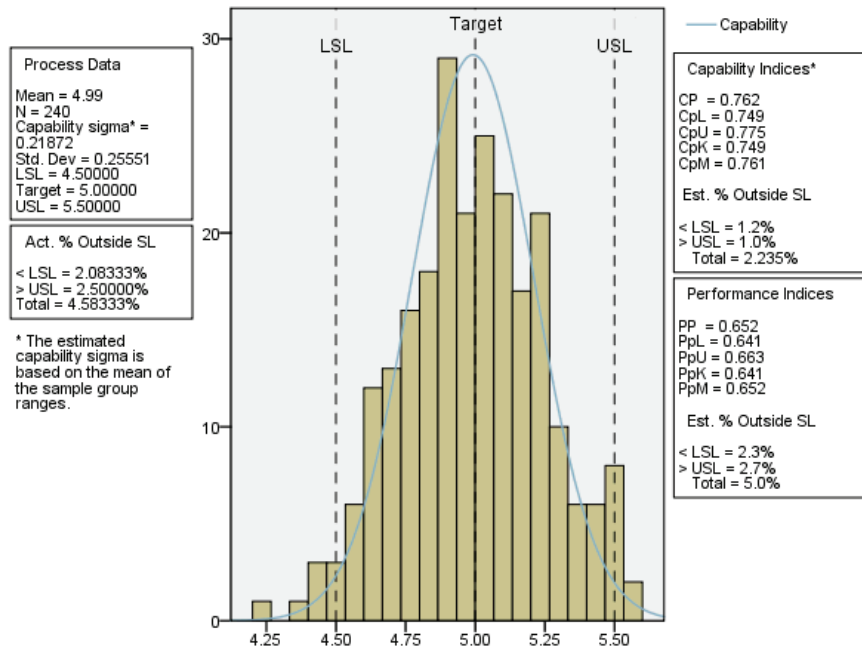
The specification is the same as the specification for p and np charts.

### **CPCHART Subcommand**

CPCHART produces a process capability chart. The default process capability chart consists of a histogram of the process values, a superimposed normal curve based on the capability sigma and the process mean, and vertical reference lines at the fixed upper and lower control limits and target value (if specified). It also displays the process capability indices, the process performance indices, and process data such as the process mean and process standard deviation.

- When CPCHART is specified the LSL and USL subcommands are required. The TARGET subcommand is optional with CPCHART.
- You can control which normal curves are superimposed on the histogram with the NORMAL subcommand.
- You can display vertical reference lines at the plus or minus three sigma level, or a user-specified sigma level, with the REFERENCE subcommand.
- All of the statistics displayed on the process capability chart are available in table form from the STATISTICS subcommand.
- Process capability charts have a default title of Process Capability. You can specify your own title using the TITLE and SUBTITLE subcommands.

Figure 230-10  
Process capability chart



## Data Organization

The data can be organized in the following ways:

- Each case is a unit with a subgroup identifier that is specified with a BY variable. Table 230-7 is an example of this type of data organization. The data do not have to be sorted by subgroup.
- Each case is a subgroup and there are as many variables as individuals within one sample. Table 230-8 shows this type of organization.
- Each case is a unit and there are no subgroups. Table 230-9 shows this type of organization. Cases are not sorted or aggregated before plotting.

Table 230-7

Each case is a unit with a subgroup identifier

Subgroup	Length
8:50	6.35
11:30	6.39
8:50	6.40
11:30	6.46
8:50	6.32
11:30	6.37
8:50	6.39
11:30	6.36
...	...

Table 230-8

*Each case is a subgroup*

N1	N2	N3	N4
6.35	6.40	6.32	6.39
6.39	6.46	6.37	6.36
...	...	...	...

Table 230-9

*Each case is a unit and there are no subgroups***pH**

4.9

4.5

4.7

5.1

4.6

5.1

4.9

4.8

...

## Variable Specification

If data are organized as shown in Table 230-7, the variable specification on the CPCHART subcommand is

```
var BY var
```

The variable that is specified before BY is the process variable, which is used to generate the histogram (for example, *length* in Table 230-7). The variable that is specified after BY is the subgroup identifier (for example, *subgroup* in Table 230-7). The process variable must be numeric, while the subgroup identifier variable can be of any type.

If data are organized as shown in Table 230-8, the variable specification on the CPCHART subcommand is

```
var var [var...]
```

Each variable contains process values for a particular subgroup. Values from all subgroups are used to generate the histogram. At least two variables are required, and each variable must be numeric.

If data are organized as shown in Table 230-9, the variable specification on the CPCHART subcommand is

```
var
```

The specified variable contains the process values used to generate the histogram, and must be numeric.



**Example**

```
SPCHART /TEMPLATE='cpchart.sgt'
/CPCHART length BY shift.
```

- The data are organized as shown in Table 230-7. *length* is a numeric variable that measures the length of a part. *shift* contains the subgroup identifier (work shift number).
- The chart uses the chart template from *cpchart.sgt*.

**STATISTICS Subcommand**

Any keyword may be specified in any place in the subcommand, but for conceptual clarity, the keywords are organized as follows: the Process Capability Indices, the Process Performance Indices, the Process Data, and the Measure(s) for Assessing Normality.

- This subcommand is silently ignored if the chart is not an XR, XS, IR, I, or CPCHART chart.
- A duplicated subcommand name causes a syntax error.
- A duplicated keyword is silently ignored.
- There is no default keyword or parameter value.

**The Process Capability Indices**

<b>CP</b>	<i>Capability of the process.</i>
<b>CPU</b>	<i>The distance between the process mean and the upper specification limit scaled by capability sigma.</i>
<b>CPL</b>	<i>The distance between the process mean and the lower specification limit scaled by capability sigma.</i>
<b>K</b>	<i>The deviation of the process mean from the midpoint of the specification limits. This measurement is computed independently of the estimated capability sigma.</i>
<b>CPK</b>	<i>Capability of the process related to both dispersion and centeredness. It is the minimum of CpU and CpL. If only one specification limit is provided, we compute and report a unilateral CpK instead of taking the minimum.</i>
<b>CR</b>	<i>The reciprocal of CP.</i>
<b>CPM</b>	<i>An index relating capability sigma and the difference between the process mean and the target value. A target value must be specified on the TARGET subcommand by the user.</i>
<b>CZU</b>	<i>The number of capability sigmas between the process mean and the upper specification limit.</i>
<b>CZL</b>	<i>The number of capability sigmas between the process mean and the lower specification limit.</i>
<b>CZMIN</b>	<i>The minimum number of capability sigmas between the process mean and the specification limits.</i>
<b>CZMAX</b>	<i>The maximum number of capability sigmas between the process mean and the specification limits.</i>
<b>CZOUT</b>	<i>The estimated percentage outside the specification limits. The standard normal approximation is based on CZ<sub>U</sub> and CZ<sub>L</sub>.</i>
<b>CZLOUT</b>	<i>The estimated percentage outside the lower specification limit. The standard normal approximation is based on CZ<sub>L</sub>.</i>
<b>CZUOUT</b>	<i>The estimated percentage outside the upper specification limit. The standard normal approximation is based on CZ<sub>U</sub>.</i>

- For each of the keywords (other than CPK), both the LSL subcommand and the USL subcommand must be specified. Otherwise, the keyword(s) are ignored, and a syntax warning is issued. For CPK, at least one of the LSL and USL subcommands must be specified.
- If the TARGET subcommand is not specified, the keyword CPM is ignored, and a syntax warning is issued.

### **The Process Performance Indices**

<b>PP</b>	<i>Performance of the process.</i>
<b>PPU</b>	<i>The distance between the process mean and the upper specification limit scaled by process standard deviation.</i>
<b>PPL</b>	<i>The distance between the process mean and the lower specification limit scaled by process standard deviation.</i>
<b>PPK</b>	<i>Performance of the process related to both dispersion and centeredness. It is the minimum of PpU and PpL. If only one specification limit is provided, we compute and report a unilateral PpK instead of taking the minimum.</i>
<b>PR</b>	<i>The reciprocal of PP.</i>
<b>PPM</b>	<i>An index relating process variance and the difference between the process mean and the target value. A target value must be specified on the TARGET subcommand by the user.</i>
<b>PZU</b>	<i>The number of standard deviations between the process mean and the upper specification limit.</i>
<b>PZL</b>	<i>The number of standard deviations between the process mean and the lower specification limit.</i>
<b>PZMIN</b>	<i>The minimum number of standard deviations between the process mean and the specification limits.</i>
<b>PZMAX</b>	<i>The maximum number of standard deviations between the process mean and the specification limits.</i>
<b>PZOUT</b>	<i>The estimated percentage outside the specification limits. The standard normal approximation is based on PZ<sub>U</sub> and PZ<sub>L</sub>.</i>
<b>PZLOUT</b>	<i>The estimated percentage outside the lower specification limit. The standard normal approximation is based on PZ<sub>L</sub>.</i>
<b>PZUOUT</b>	<i>The estimated percentage outside the upper specification limit. The standard normal approximation is based on PZ<sub>U</sub>.</i>

- For each of the keywords (other than PPK), both the LSL subcommand and the USL subcommand must be specified. Otherwise, we ignore the keyword(s) and issue a syntax warning. For PPK, at least one of the LSL and USL subcommands must be specified.
- If the TARGET subcommand is not specified, the keyword PPM is ignored, and a syntax warning is issued.

### **Process Data**

<b>N</b>	<i>Total sample size.</i>
<b>MEAN</b>	<i>The process mean.</i>
<b>STDDEV</b>	<i>The process standard deviation.</i>
<b>CAPSIGMA</b>	<i>The capability sigma.</i>
<b>LSL</b>	<i>The fixed lower control limit.</i>

<b>USL</b>	<i>The fixed upper control limit.</i>
<b>TARGET</b>	<i>The target value.</i>

### **Measure(s) for Assessing Normality**

<b>AZOUT</b>	<i>The observed percentage outside the specification limits. A point is defined outside the specification limits when its value is greater than or equal to the upper specification limit or is less than or equal to the lower specification limit.</i>
<b>AZLOUT</b>	<i>The observed percentage outside the lower specification limit. A point is defined outside the lower specification limit when its value is less than or equal to the lower specification limit.</i>
<b>AZUOUT</b>	<i>The observed percentage outside the upper specification limit. A point is defined outside the upper specification limit when its value is greater than or equal to the upper specification limit.</i>

- For each of the keywords, both the `LSL` subcommand and the `USL` subcommand must be specified. Otherwise, we ignore the keyword and issue a syntax warning.

### **RULES Subcommand**

`RULES` specifies the rules for identifying out-of-control points. If a point violates any rule, it appears in the primary chart with a different shape and color compared to in-control points. Note that only the last point in a discovered pattern shall be marked for a rule violation, and any and all overlapping patterns are treated as separate patterns. A table of rule violations is also included in the output that displays the points marked in the control chart. If desired, use the `ID` keyword to specify the variable that identifies points in this table.

- Any keyword may be specified in any place in the subcommand.
- A duplicated subcommand name causes a syntax error.
- A duplicated keyword is silently ignored.
- The default keyword is `ALL`.
- If the subcommand is omitted, no control rules are used.
- `RULES` does not apply when generating a process capability chart (`CPCHART` subcommand).

<b>ALL</b>	<i>All rules.</i>
<b>UCL</b>	<i>Greater than +3 sigma.</i>
<b>R2UPPER</b>	<i>2 points out of the last 3 greater than +2 sigma.</i>
<b>R4UPPER</b>	<i>4 points out of the last 5 greater than +1 sigma.</i>
<b>R8UPPER</b>	<i>8 consecutive points above the center line.</i>
<b>R8LOWER</b>	<i>8 consecutive points below the center line.</i>
<b>R4LOWER</b>	<i>4 points out of the last 5 less than -1 sigma.</i>
<b>R2LOWER</b>	<i>2 points out of the last 3 less than -2 sigma.</i>
<b>LCL</b>	<i>Less than -3 sigma.</i>
<b>TRUP</b>	<i>6 consecutive points trending up.</i>
<b>TRDOWN</b>	<i>6 consecutive points trending down.</i>
<b>ALTERNATING</b>	<i>14 consecutive points alternating.</i>

## **ID Subcommand**

ID specifies a variable that identifies points in the table of rule violations. If this subcommand is omitted, the BY variable is used. Without the RULES subcommand, ID has no effect.

## **CAPSIGMA Subcommand**

This subcommand defines the capability sigma estimator, which is required in computing all the Process Capability Indices except  $\kappa$ . There are four options:

<b>RBAR</b>	<i>Mean sample range.</i> The estimated capability sigma is based on the mean of the sample group ranges.
<b>SBAR</b>	<i>Mean sample standard deviation.</i> The estimated capability sigma is based on the mean of the sample group standard deviations.
<b>MRBAR</b>	<i>Mean sample moving range.</i> The estimated capability sigma is based on the mean of the sample moving ranges. The span that is defined by the SPAN subcommand is used. (Recall that its passive default value is 2.)
<b>WITHIN</b>	<i>Sample within-group variance.</i> The estimated capability sigma is the square root of the sample within-group variance.

The validity of specification depends on the chart specification (i.e., /XR, /XS, /I, /IR, or /CPCHART).

Table 230-10  
Valid CAPSIGMA options by chart specification

<b>Chart Specification</b>	<b>Valid CAPSIGMA Options</b>
XR	RBAR (default) SBAR WITHIN
XS	RBAR SBAR (default) WITHIN
I	MRBAR (default)
IR	MRBAR (default)
CPCHART	RBAR SBAR WITHIN MRBAR

- When this subcommand is omitted or specified without a keyword by the user, the default conditional on the chart specification is implicitly assumed (see the table above). When the chart specification is /CPCHART, the default is RBAR when the data are organized in subgroups, and MRBAR otherwise.
- An invalid specification (e.g., /I and /CAPSIGMA = RBAR) causes a syntax error when the chart specification is /XR, /XS, /I, or /IR and an applicable STATISTICS keyword is specified. An invalid specification (for those same chart types) without an applicable STATISTICS keyword generates a syntax warning.
- When the chart specification is /CPCHART, an invalid specification (e.g., /CAPSIGMA = RBAR when the data are not organized in subgroups) causes a syntax error.

- When the chart specification is /XR, /XS, /I, or /IR, CAPSIGMA is silently ignored if the subcommand specification is valid (with respect to the chart specification) but an applicable STATISTICS keyword is not specified.
- CAPSIGMA is silently ignored if the chart specification is not /XR, /XS, /I, /IR, or /CPCHART.
- A duplicated subcommand name causes a syntax error.
- A duplicated keyword is silently ignored, but if two or more keywords are specified and they do not have identical meanings, a syntax error message is issued.

## ***SPAN Subcommand***

SPAN specifies the span from which the moving range for an individuals chart or process capability chart (when the data are specified in the form `var`) is calculated. The specification must be an integer value that is greater than 1. The default is 2. SPAN applies only to I, IR, and CPCHART chart specifications.

### ***Example***

```
SPCHART /IR=SUBSIZE /SPAN=5.
```

- The SPAN subcommand specifies that the moving ranges are computed from every five individual samples.

## ***CONFORM and NONCONFORM Subcommands***

Either CONFORM or NONCONFORM is required when you specify a status variable on the P or NP subcommand. You make that specification when data are organized so that each case is an inspection unit (see P and NP Subcommands on p. 1907).

- Either subcommand requires a value specification. The value can be numeric or string. String values must be enclosed within quotes.
- If CONFORM is specified, all values for the status variable (other than the specified value) are tabulated as nonconformities. If NONCONFORM is specified, only the specified value is tabulated as a nonconformity.
- CONFORM and NONCONFORM apply only to P and NP chart specifications.

## ***SIGMAS Subcommand***

SIGMAS allows you to define the sigma level for a control chart. The value specified on SIGMAS is used in calculating the upper and lower control limits on the chart. It also determines the placement of reference lines on a process capability chart (/CPCHART) when the REFERENCE subcommand includes the keyword SIGMAS. You can specify a number larger than 1 but less than or equal to 10. A larger SIGMAS value means a greater range between the upper and the lower control limits. The default is 3.

## ***MINSAMPLE Subcommand***

MINSAMPLE specifies the minimum sample size for X-bar, R, or s charts, as well as process capability charts when the data are organized in subgroups. When you specify XR, XS, or CPCHART on SPCHART, any subgroup with a size that is smaller than the size that is specified on MINSAMPLE is excluded from the chart and from all computations. If each case is a subgroup, there must be at least as many variables named as the number that is specified on MINSAMPLE. The default, as well as the minimum, is 2.

## ***LSL and USL Subcommand***

LSL and USL allow you to specify fixed lower and upper control limits. Fixed control limits are often used in manufacturing processes as designer-specified limits.

- For chart types other than CPCHART, these limits are displayed on the chart, along with the calculated limits. If you do not specify LSL and USL, no fixed control limits are displayed. However, if you want only the specified control limits, you must edit the chart in a chart window to suppress the calculated series.
- For a chart type of CPCHART (process capability chart), LSL and USL are required. The REFERENCE subcommand controls whether these limits are displayed on the process capability chart. By default, they are displayed.

### ***Example***

```
SPCHART /TEMPLATE='ctrl.sgt'  
/XS=SUBSIZE  
/USL=74.50  
/LSL=73.50.
```

- The USL and LSL subcommands specify the control limits according to the designing engineer. The center line is probably at 74.00.
- The specified upper and lower limits are displayed together with the control limits that are calculated from the observed standard deviation and the sigma value.

## ***TARGET Subcommand***

This subcommand defines the target value that is used in computing CpM and PpM requested by the STATISTICS subcommand. It is also the target value that is displayed by default on a process capability chart (/CPCHART). The value may be any real number that is less than or equal to the USL value and greater than or equal to the LSL value.

- This subcommand is silently ignored if no applicable STATISTICS keyword is specified or the chart type is not CPCHART.
- If the value is numeric but out of the valid range, we issue a warning and ignore the CPM and/or PPM keyword(s), if any, in the STATISTICS subcommand.

## **MISSING Subcommand**

MISSING controls the treatment of missing values.

- The default is NOREPORT and EXCLUDE.
- REPORT and NOREPORT are alternatives and apply only to category variables. REPORT and NOREPORT control whether categories (subgroups) with missing values are created. They do not apply to process capability charts (/CPCHART).
- INCLUDE and EXCLUDE are alternatives and apply to process variables.

<b>NOREPORT</b>	<i>Suppress missing-value categories. This setting is the default.</i>
<b>REPORT</b>	<i>Report and plot missing-value categories.</i>
<b>EXCLUDE</b>	<i>Exclude user-missing values. Both user-missing and system-missing values for the process variable are excluded from computation and plotting. This setting is the default.</i>
<b>INCLUDE</b>	<i>Include user-missing values. Only system-missing values for the process variable are excluded from computation and plotting.</i>

## **NORMAL Subcommand**

NORMAL controls which normal curves are superimposed on a process capability chart. It only applies when the chart specification is /CPCHART. Specify one of the following values:

<b>CAPABILITY</b>	<i>Superimposes a normal curve with mean equal to the process mean and standard deviation equal to the capability sigma. This is the default.</i>
<b>PERFORMANCE</b>	<i>Superimposes a normal curve with mean equal to the process mean and standard deviation equal to the process standard deviation.</i>
<b>BOTH</b>	<i>Superimposes both the CAPABILITY and PERFORMANCE normal curves.</i>
<b>NONE</b>	<i>Specifies that no normal curves will be displayed.</i>

## **REFERENCE Subcommand**

REFERENCE controls which vertical reference lines are displayed on a process capability chart. It only applies when the chart specification is /CPCHART.

<b>SL</b>	<i>Vertical reference lines are displayed at the specified values of LSL and USL as well as at the value of TARGET if specified. This is the default.</i>
<b>SIGMAS</b>	<i>Vertical reference lines are displayed at the process mean and at a specified number of capability sigmas above and below the mean. The number of capability sigmas is given by the value of the SIGMAS subcommand, which is 3 by default.</i>
<b>NONE</b>	<i>Specifies that no reference lines will be displayed. This is an alternative to SL and/or SIGMAS and cannot be used in combination with them.</i>

# SPECTRA

SPECTRA is available in the Forecasting option.

SPECTRA VARIABLES= series names

```
[/{CENTER NO**}
  {CENTER      }

[/{CROSS NO**}
  {CROSS       }

[/WINDOW={HAMMING** [( {5  } )]   }
  {          {span}          }
  {BARTLETT [(span)]          }
  {PARZEN [(span)]           }
  {TUKEY [(span)]            }
  {UNIT or DANIELL [(span)]  }
  {NONE                      }
  {w-p, ..., w0, ..., wp    }

[/PLOT= [P] [S] [CS] [QS] [PH] [A]
  [G] [K] [ALL] [NONE]
  [BY {FREQ  }]]
  {PERIOD}

[/SAVE = [FREQ (name)] [PER (name)] [SIN (name)]
  [COS (name)] [P (name)] [S (name)]
  [RC (name)] [IC (name)] [CS (name)]
  [QS (name)] [PH (name)] [A (name)]
  [G (name)] [K (name)]]

[/APPLY [= 'model name']]
```

**\*\*Default if the subcommand is omitted.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
SPECTRA VARIABLES = HSTARTS.
```

## Overview

SPECTRA plots the periodogram and spectral density function estimates for one or more series. You can also request bivariate spectral analysis. Moving averages, termed *windows*, can be used for smoothing the periodogram values to produce spectral densities.

### Options

**Output.** In addition to the periodogram, you can use the PLOT subcommand to produce a plot of the estimated spectral density. You can use the keyword BY on PLOT to suppress the display of the plot by frequency or the plot by period. To reduce the range of values that are displayed in the plots, you can center the data by using the CENTER subcommand.



**Cross-Spectral Analysis.** You can specify cross-spectral (bivariate) analysis with the `CROSS` subcommand and select which bivariate plots are produced by using `PLOT`.

**New Variables.** Variables that are computed by `SPECTRA` can be saved to the active dataset for use in subsequent analyses with the `SAVE` subcommand. `TSET MXNEWVAR` specifies the maximum number of new variables that can be generated by a procedure. The default is 60.

**Spectral Windows.** You can specify a spectral window and its span for calculation of the spectral density estimates.

### ***Basic Specification***

The basic specification is one or more series names.

- By default, `SPECTRA` plots the periodogram for each specified series. The periodogram is shown first by frequency and then by period. No new variables are saved by default.

### ***Subcommand Order***

- Subcommands can be specified in any order.

### ***Syntax Rules***

- `VARIABLES` can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each subcommand is executed.

### ***Operations***

- `SPECTRA` cannot process series with missing observations. (You can use the `RMV` command to replace missing values, use `TSET MISSING=INCLUDE` to include user-missing values, and use `USE` to ignore missing observations at the beginning or end of a series. See `RMV` and `USE` for more information.)
- If the number of observations in the series is odd, the first case is ignored.
- If the `SAVE` subcommand is specified, new variables are created for each specified series. For bivariate analyses, new variables are created for each series pair.
- `SPECTRA` requires memory both to compute variables and to build plots. Requesting fewer plots may enable you to analyze larger series.

### ***Limitations***

- A maximum of one `VARIABLES` subcommand is allowed. There is no limit on the number of series named on the list.

## ***Example***

```
SPECTRA VARIABLES = HSTARTS  
/CENTER  
/PLOT P S BY FREQ.
```

- This example produces a plot of the periodogram and spectral density estimate for series *HSTARTS*.
- `CENTER` adjusts the series to have a mean of 0.
- `PLOT` specifies that the periodogram (`P`) and the spectral density estimate (`S`) should be plotted against frequency (`BY FREQ`).

## ***VARIABLES Subcommand***

`VARIABLES` specifies the series names and is the only required subcommand.

- `VARIABLES` must be specified before the other subcommands.
- Each specified series is analyzed separately unless the `CROSS` subcommand is specified.
- The series must contain at least six cases.

### ***Example***

```
SPECTRA VARIABLES = VARX VARY.
```

- This command produces the default display for two series, *VARX* and *VARY*.

## ***CENTER Subcommand***

`CENTER` adjusts the series to have a mean of 0. This process reduces the range of values that are displayed in the plots.

- If `CENTER` is not specified, the ordinate of the first periodogram value is  $2n$  times the square of the mean of the series, where  $n$  is the number of cases.
- You can specify `CENTER NO` to suppress centering when applying a previous model with `APPLY`.

### ***Example***

```
SPECTRA VARIABLES = VARX VARY
/CENTER.
```

- This example produces the default display for *VARX* and *VARY*. The plots are based on the series after their means have been adjusted to 0.

## ***WINDOW Subcommand***

`WINDOW` specifies a spectral window to use when the periodogram is smoothed to obtain the spectral density estimate. If `WINDOW` is not specified, the Tukey-Hamming window with a span of 5 is used.

- The specification on `WINDOW` is a window name and a span in parentheses, or a sequence of user-specified weights.
- The window name can be any one of the keywords listed below.

- Only one window keyword is accepted. If more than one keyword is specified, the first keyword is used.
- The span is the number of periodogram values in the moving average and can be any integer. If an even number is specified, it is decreased by 1.
- Smoothing near the end of series is accomplished via reflection. For example, if the span is 5, the second periodogram value is smoothed by averaging the first, third, and fourth values and twice the second value.

The following data windows can be specified. Each formula defines the upper half of the window. The lower half is symmetric with the upper half. In all formulas,  $p$  is the integer part of the number of spans divided by 2,  $D_p$  is the Dirichlet kernel of order  $p$ , and  $F_p$  is the Fejer kernel of order  $p$  (Priestley, 1981).

**HAMMING**      *Tukey-Hamming window.* The weights are  

$$W_k = 0.54D_p(2\pi f_k) + 0.23D_p\left(2\pi f_k + \frac{\pi}{p}\right) + 0.23D_p\left(2\pi f_k - \frac{\pi}{p}\right)$$
 where  $k=0, \dots, p$ . This is the default.

**TUKEY**      *Tukey-Hanning window.* The weights are  

$$W_k = 0.5D_p(2\pi f_k) + 0.25D_p\left(2\pi f_k + \frac{\pi}{p}\right) + 0.25D_p\left(2\pi f_k - \frac{\pi}{p}\right)$$
 where  $k=0, \dots, p$ .

**PARZEN**      *Parzen window.* The weights are  

$$W_k = \frac{1}{p}(2 + \cos(2\pi f_k))(F_{p/2}(2\pi f_k))^2$$
 where  $k=0, \dots, p$ .

**BARTLETT**      *Bartlett window.* The weights are  

$$W_k = F_p(2\pi f_k)$$
 where  $k=0, \dots, p$ .

**UNIT**      *Equal-weight window.* The weights are  $w_k = 1$  where  $k=0, \dots, p$ . DANIELL is an alias for UNIT.

**NONE**      *No smoothing.* If NONE is specified, the spectral density estimate is the same as the periodogram.

**w**      *User-specified weights.*  $W_0$  is applied to the periodogram value that is being smoothed, and the weights on either side are applied to preceding and following values. If the number of weights is even, it is assumed that  $w_p$  is not supplied. The weight after the middle one is applied to the periodogram value being smoothed.  $W_0$  must be positive.

### Example

```
SPECTRA VARIABLES = VAR01
/WINDOW=TUKEY(3)
/PLOT=P S.
```

- In this example, the Tukey window weights with a span of 3 are used.
- The `PLOT` subcommand plots both the periodogram and the spectral density estimate, both by frequency and period.

## ***PLOT Subcommand***

`PLOT` specifies which plots are displayed.

- If `PLOT` is not specified, only the periodogram is plotted for each specified series. Each periodogram is shown both by frequency and by period.
- You can specify more than one plot keyword.
- Keywords can be specified in any order.
- The plot keywords `K`, `CS`, `QS`, `PH`, `A`, and `G` apply only to bivariate analyses. If the subcommand `CROSS` is not specified, these keywords are ignored.
- The period (horizontal) axis on a plot by period (`BY PERIOD`) is scaled in natural logarithms from 0.69 to  $\ln(n)$ , where  $n$  is the number of cases.
- The frequency (horizontal) axis on a plot by frequency (`BY FREQ`) is scaled from 0 to 0.5, expressing the frequency as a fraction of the length of the series.
- The periodogram and estimated spectrum (vertical axis) are scaled in natural logs.

The following plot keywords are available:

<b>P</b>	<i>Periodogram.</i> This setting is the default.
<b>S</b>	<i>Spectral density estimate.</i>
<b>K</b>	<i>Squared coherency.</i> Applies only to bivariate analyses.
<b>CS</b>	<i>Cospectral density estimate.</i> Applies only to bivariate analyses.
<b>QS</b>	<i>Quadrature spectrum estimate.</i> Applies only to bivariate analyses.
<b>PH</b>	<i>Phase spectrum.</i> Applies only to bivariate analyses.
<b>A</b>	<i>Cross amplitude.</i> Applies only to bivariate analyses.
<b>G</b>	<i>Gain.</i> Applies only to bivariate analyses.
<b>ALL</b>	<i>All plots.</i> For bivariate analyses, this setting includes all plots listed above. For univariate analyses, this setting includes the periodogram and the spectral density estimate.

## ***BY Keyword***

By default, `SPECTRA` displays both frequency and period plots. You can use `BY` to produce only frequency plots or only period plots.

- `BY FREQ` indicates that all plots are plotted by frequency only. Plots by period are not produced.
- `BY PERIOD` indicates that all plots are plotted by period only. Plots by frequency are not produced.

### ***Example***

```
SPECTRA VARIABLES = SER01
```

---

```
/PLOT=P S BY FREQ.
```

- This command plots both the periodogram and the spectral density estimate for *SERO1*. The plots are shown by frequency only.

## ***CROSS Subcommand***

CROSS is used to specify bivariate spectral analysis.

- When CROSS is specified, the first series named on the VARIABLES subcommand is the independent variable. All remaining variables are dependent.
- Each series after the first series is analyzed with the first series independently of other series that is named.
- Univariate analysis of each specified series is still performed.
- You can specify CROSS NO to turn off bivariate analysis when applying a previous model with APPLY.

### ***Example***

```
SPECTRA VARIABLES = VARX VARY VARZ
/CROSS.
```

- In this example, bivariate spectral analyses of series *VARX* with *VARY* and *VARX* with *VARZ* are requested in addition to the usual univariate analyses of *VARX*, *VARY*, and *VARZ*.

## ***SAVE Subcommand***

SAVE saves computed SPECTRA variables to the active dataset for later use. SPECTRA displays a list of the new variables and their labels, showing the type and source of those variables.

- You can specify any or all of the output keywords listed below.
- A name to be used for generating variable names must follow each output keyword. The name must be enclosed in parentheses.
- For each output keyword, one variable is created for each series named on SPECTRA and for each bivariate pair.
- The keywords RC, IC, CS, QS, PH, A, G, and K apply only to bivariate analyses. If CROSS is not specified, these keywords are ignored.
- SAVE specifications are not used when models are reapplied by using APPLY. They must be specified each time variables are to be saved.
- The output variables correspond to the Fourier frequencies. They do not correspond to the original series.
- Because each output variable has only  $(n/2 + 1)$  cases (where  $n$  is the number of cases), the values for the second half of the series are set to system-missing.
- Variable names are generated by adding  $\_n$  to the specified name, where  $n$  ranges from 1 to the number of series specified.

- For bivariate variables, the suffix is *\_n\_n*, where the *ns* indicate the two variables that are used in the analysis.
- The frequency (FREQ) and period (PER) variable names are constant across all series and do not have a numeric suffix.
- If the generated variable name is longer than the maximum variable name length, or if the specified name already exists, the variable is not saved.

The following output keywords are available:

<b>FREQ</b>	<i>Fourier frequencies.</i>
<b>PER</b>	<i>Fourier periods.</i>
<b>SIN</b>	<i>Value of a sine function at the Fourier frequencies.</i>
<b>COS</b>	<i>Value of a cosine function at the Fourier frequencies.</i>
<b>P</b>	<i>Periodogram values.</i>
<b>S</b>	<i>Spectral density estimate values.</i>
<b>RC</b>	<i>Real part values of the cross-periodogram. Applies only to bivariate analyses.</i>
<b>IC</b>	<i>Imaginary part values of the cross-periodogram. Applies only to bivariate analyses.</i>
<b>CS</b>	<i>Cospectral density estimate values. Applies only to bivariate analyses.</i>
<b>QS</b>	<i>Quadrature spectrum estimate values. Applies only to bivariate analyses.</i>
<b>PH</b>	<i>Phase spectrum estimate values. Applies only to bivariate analyses.</i>
<b>A</b>	<i>Cross-amplitude values. Applies only to bivariate analyses.</i>
<b>G</b>	<i>Gain values. Applies only to bivariate analyses.</i>
<b>K</b>	<i>Squared coherency values. Applies only to bivariate analyses.</i>

### **Example**

```
SPECTRA VARIABLES=STRIKES RUNS
  /SAVE= FREQ (FREQ) P (PGRAM) S (SPEC).
```

- This example creates five variables: *FREQ*, *PGRAM\_1*, *PGRAM\_2*, *SPEC\_1*, and *SPEC\_2*.

## **APPLY Subcommand**

APPLY allows you to use a previously defined SPECTRA model without having to repeat the specifications.

- The only specification on APPLY is the name of a previous model in quotation marks. If a model name is not specified, the model that was specified on the previous SPECTRA command is used. Model names are either the default *MOD\_n* names that are assigned by Forecasting or the names that are assigned on the MODEL NAME command.
- To change one or more model specifications, specify the subcommands of only those portions you want to change after the APPLY subcommand.
- If no series are specified on the command, the series that were originally specified with the model that is being reapplied are used.
- To change the series that are used with the model, enter new series names before or after the APPLY subcommand. If a variable name is specified before APPLY, the slash before the subcommand is required.

- The `SAVE` specifications from the previous model are *not* reused by `APPLY`. They must be specified each time that variables are to be saved.

### **Examples**

```
SPECTRA VARIABLES = VAR01
  /WINDOW=DANIELL (3)
  /CENTER
  /PLOT P S BY FREQ.
SPECTRA APPLY
  /PLOT P S.
```

- The first command plots both the periodogram and the spectral density estimate for *VAR01*. The plots are shown by frequency only.
- Because the `PLOT` subcommand is respecified, the second command produces plots by both frequency and period. All other specifications remain the same as in the first command.

### **References**

- Bloomfield, P. 1976. *Fourier analysis of time series*. New York: John Wiley and Sons.
- Fuller, W. A. 1976. *Introduction to statistical time series*. New York: John Wiley and Sons.
- Gottman, J. M. 1981. *Time-series analysis: A comprehensive introduction for social scientists*. Cambridge: Cambridge University Press.
- Priestley, M. B. 1981. *Spectral analysis and time series, volumes 1 and 2*. London: Academic Press.

# SPLIT FILE

```
SPLIT FILE          {OFF          }  
                   [{LAYERED }] {BY varlist}  
                   {SEPARATE}
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
SPLIT FILE BY SEX.
```

## **Overview**

`SPLIT FILE` splits the active dataset into subgroups that can be analyzed separately. These subgroups are sets of adjacent cases in the file that have the same values for the specified split variables. Each value of each split variable is considered a break group, and cases within a break group must be grouped together in the active dataset. If they are not grouped together, the `SORT CASES` command must be used before `SPLIT FILE` to sort cases in the proper order.

### **Basic Specification**

The basic specification is keyword `BY` followed by the variable or variables that define the split-file groups.

- By default, the split-file groups are compared within the same table(s).
- You can turn off split-file processing by using keyword `OFF`.

### **Syntax Rules**

- `SPLIT FILE` can specify both numeric and string split variables, including variables that are created by temporary transformations. `SPLIT FILE` cannot specify scratch or system variables.
- `SPLIT FILE` is in effect for all procedures in a session unless you limit it with a `TEMPORARY` command, turn it off, or override it with a new `SPLIT FILE` or `SORT CASES` command.

### **Operations**

- `SPLIT FILE` takes effect as soon as it is encountered in the command sequence. Therefore, pay special attention to the position of `SPLIT FILE` among commands. [For more information, see the topic Command Order on p. 41.](#)
- The file is processed sequentially. A change or break in values on any one of the split variables signals the end of one break group and the beginning of the next break group.



- AGGREGATE ignores the SPLIT FILE command. To split files by using AGGREGATE, name the variables that are used to split the file as break variables ahead of any other break variables on AGGREGATE. AGGREGATE still produces one file, but the aggregated cases are in the same order as the split-file groups.
- If SPLIT FILE is in effect when a procedure writes matrix materials, the program writes one set of matrix materials for every split group. If a procedure reads a file that contains multiple sets of matrix materials, the procedure automatically detects the presence of multiple sets.
- If SPLIT FILE names any variable that was defined by the NUMERIC command, the program prints page headings that indicate the split-file grouping.

### **Limitations**

- SPLIT FILE can specify or imply up to eight variables.
- Each eight bytes of a string variable counts as a variable toward the limit of eight variables. So a string variable with a defined width of greater than 64 bytes cannot be used as a split file variable.

## **LAYERED and SEPARATE Subcommands**

LAYERED and SEPARATE specify how split-file groups are displayed in the output.

- Only one of these subcommands can be specified. If neither subcommand is specified with the BY variable list, LAYERED is the default.
- LAYERED and SEPARATE do not apply to the text output.

**LAYERED**            *Display split-file groups in the same table in the outermost column.*  
**SEPARATE**         *Display split-file groups as separate tables.*

## **Examples**

### **Sorting and Splitting a File**

```
SORT CASES BY SEX.  
SPLIT FILE BY SEX.  
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.
```

- SORT CASES arranges cases in the file according to the values of variable *SEX*.
- SPLIT FILE splits the file according to the values of variable *SEX*, and FREQUENCIES generates separate median income tables for men and women.
- By default, the two groups (men and women) are compared in the same Frequency and Statistics tables.

### **Applying a Temporary Split File**

```
SORT CASES BY SEX.  
TEMPORARY.  
SPLIT FILE SEPARATE BY SEX.  
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.  
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.
```

- Because of the `TEMPORARY` command, `SPLIT FILE` applies to the first procedure only. Thus, the first `FREQUENCIES` procedure generates separate tables for men and women. The second `FREQUENCIES` procedure generates tables that include both sexes.

### ***Turning Off a Split File***

```
SORT CASES BY SEX.  
SPLIT FILE SEPARATE BY SEX.  
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.  
SPLIT FILE OFF.  
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.
```

- `SPLIT FILE` does not apply to the second `FREQUENCIES` procedure because it is turned off after the first `FREQUENCIES` procedure. This example produces the same results as the example above.

### ***Overriding a Previous Split File***

```
SORT CASES BY SEX RACE.  
SPLIT FILE BY SEX.  
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.  
SPLIT FILE BY SEX RACE.  
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.
```

- The first `SPLIT FILE` command applies to the first `FREQUENCIES` procedure. The second `SPLIT FILE` command overrides the first command and splits the file by sex and race. This split is in effect for the second `FREQUENCIES` procedure.

# STRING

```
STRING varlist (An) [/varlist...]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
STRING STATE1 (A2).
```

## **Overview**

STRING declares new string variables that can be used as target variables in data transformations.

### **Basic Specification**

The basic specification is the name of the new variables and, in parentheses, the variable format.

### **Syntax Rules**

- If keyword `TO` is used to create multiple string variables, the specified format applies to each variable that is named and implied by `TO`.
- To declare variables with different formats, separate each format group with a slash.
- `STRING` can be used within an input program to determine the order of string variables in the dictionary of the active dataset. When used for this purpose, `STRING` must precede `DATA LIST` in the input program. See the Examples for the `NUMERIC` command on p. 1396.
- `STRING` cannot be used to redefine an existing variable.
- String variables cannot have zero length; `A0` is an illegal format.
- All implementations of the program allow the `A` format. Other string formats may be available on some systems. In addition, the definition of a long string depends on your operating system. Use keyword `LOCAL` on the `INFO` command to obtain documentation for your operating system.

### **Operations**

- `STRING` takes effect as soon as it is encountered in the command sequence. Therefore, pay special attention to the position of `STRING` among commands. [For more information, see the topic Command Order on p. 41.](#)
- New string variables are initialized as blanks.
- Variables that are declared on `STRING` are added to the active dataset in the order in which they are specified. This order is not changed by the order in which the variables are used in the transformation language.

- To change the defined width of a string variable use the `ALTER TYPE` command. [For more information, see the topic ALTER TYPE on p. 175.](#)

## Examples

```
STRING STATE1 (A2).  
RECODE STATE ('IO'='IA') (ELSE=COPY) INTO STATE1.
```

- `STRING` declares variable `STATE1` with an `A2` format.
- `RECODE` specifies `STATE` as the source variable and specifies `STATE1` as the target variable. The original value `IO` is recoded to `IA`. Keywords `ELSE` and `COPY` copy all other state codes unchanged. Thus, `STATE` and `STATE1` are identical except for cases with the original value `IO`.

```
STRING V1 TO V6 (A8) / V7 V10 (A16).
```

- `STRING` declares variables `V1`, `V2`, `V3`, `V4`, `V5`, and `V6`, each with an `A8` format, and variables `V7` and `V10`, each with an `A16` format.

# ***SUBTITLE***

```
SUBTITLE [']text[']
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Example***

```
SUBTITLE "Children's Training Shoes Only".
```

## ***Overview***

`SUBTITLE` inserts a left-justified subtitle on the second line from the top of each page of the output. The default subtitle contains the installation name and information about the hardware and operating system.

## ***Basic Specification***

The only specification is the subtitle itself.

## ***Syntax Rules***

- The subtitle can include any characters. To specify a blank subtitle, enclose a blank within quotes.
- The subtitle can be up to 60 characters. Subtitles that are longer than 60 characters are truncated.
- The quotes enclosing the subtitle are optional; using them allows you to include single or double quotes in the subtitle.
- More than one `SUBTITLE` command is allowed in a single session.
- A subtitle cannot be placed between a procedure command and `BEGIN DATA-END DATA` or within data records when the data are inline.

## ***Operations***

- Each `SUBTITLE` command overrides the previous command and takes effect on the next output page.
- `SUBTITLE` is independent of `TITLE` and each command can be changed separately.
- The subtitle will not be displayed if `HEADER=NO` is specified on `SET`.

## Examples

### **Using Single and Double Quotes**

```
TITLE 'Running Shoe Study from Runner''s World Data'.  
SUBTITLE "Children's Training Shoes Only".
```

- The title is enclosed in single quotes, so the single quote (apostrophe) in *Runner's* must be specified as a two single quotes.
- The subtitle is enclosed in double quotes, so the apostrophe in *Children's* is simply specified as a single quote.

### **Suppressing the Default Subtitle**

```
TITLE 'Running Shoe Study from Runner''s World Data'.  
SUBTITLE ' '.
```

- This subtitle is specified as a blank, which suppresses the default subtitle.

# SUMMARIZE

SUMMARIZE is available in the Statistics Base option.

```
SUMMARIZE [TABLES=]{varlist} [BY varlist] [BY...] [/varlist...]
           {ALL      }

[/TITLE = 'string']      [/FOOTNOTE= 'string']

[/CELLS= [COUNT**] [MEAN ] [STDDEV]
          [MEDIAN] [GMEDIAN] [SEMEAN] [SUM ]
          [MIN] [MAX] [RANGE] [VARIANCE]
          [KURT] [SEKURT] [SKEW] [SESKEW]
          [FIRST] [LAST]
          [NPCT] [SPCT] [NPCT(var)] [SPCT(var)]
          [HARMONIC] [GEOMETRIC]
          [DEFAULT]
          [ALL] [NONE] ]

[/MISSING={ {EXCLUDE**} } [ {VARIABLE** } ]
          {INCLUDE } {TABLE }
          {DEPENDENT}

[/FORMAT={ {NOLIST** } } [ {CASENUM } ] [ {TOTAL**} ] [MISSING='string']
          {LIST [LIMIT=n]} {NOCASENUM } {NOTOTAL}
          {VALIDLIST }

[/STATISTICS={ANOVA} [ {LINEARITY} ] [NONE**] ]
              {ALL      }
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
SUMMARIZE SALES.
```

## Overview

SUMMARIZE produces univariate statistics for specified variables. You can break the variables into groups defined by one or more control (independent) variables. Another procedure that displays univariate statistics is FREQUENCIES.

### Options

**Cell Contents.** By default, SUMMARIZE displays means, standard deviations, and cell counts. You can also use the CELLS subcommand to display aggregated statistics, including sums, variances, median, range, kurtosis, skewness, and their standard error.

**Statistics.** In addition to the statistics that are displayed for each cell of the table, you can use the STATISTICS subcommand to obtain a one-way analysis of variance and test of linearity.

**Format.** By default, `SUMMARIZE` produces a Summary Report with a total category for each group that is defined by the control variables. You can use the `FORMAT` subcommand to request a Listing Report with or without case numbers. You can also remove the total category from each group. You can use the `TITLE` and `FOOTNOTE` subcommands to specify a title and a caption for the Summary or Listing Report.

### ***Basic Specification***

The basic specification is `TABLES` with a variable list. Each variable creates a category. The actual keyword `TABLES` can be omitted.

- The minimum specification is a dependent variable.
- By default, `SUMMARIZE` displays a Case Processing Summary table, showing the number and percentage of cases included, excluded, and their total. `SUMMARIZE` also displays a Summary Report, showing means, standard deviations, and number of cases for each category.

### ***Syntax Rules***

- Both numeric and string variables can be specified. String variables can be short or long.
- If there is more than one `TABLES` subcommand, `FORMAT=LIST` or `VALIDLIST` results in an error.
- String specifications for `TITLE` and `FOOTNOTE` must be enclosed in quotes and cannot exceed 255 bytes. When the specification breaks on multiple lines, enclose each line in quotes, and separate the specifications for each line by at least one blank.
- Each subcommand (except `TABLES`) can be specified only once. If a subcommand is specified multiple times, a warning results, and the last specification is used.
- Multiple `TABLES` subcommands are allowed, but multiple specifications use a lot of computer resources and time.
- There is no limit on the number of variables that you can specify on each `TABLES` subcommand.
- When a variable is specified more than once, only the first occurrence is honored. If the same variables are specified after different `BY` keywords, an error results.

### ***Limitations***

- Only 10 `BY` keywords can be specified.

### ***Operations***

- The data are processed sequentially. It is not necessary to sort the cases before processing. If a `BY` keyword is used, the output is always sorted.
- A Case Processing Summary table is always generated, showing the number and percentage of the cases included, excluded, and the total.
- For each combination of control variables specified after different `BY` keywords, `SUMMARIZE` produces a group in the Summary Report (depending on the specification on the `FORMAT` subcommand). By default, mean, standard deviation, and number of cases are displayed for each group and for the total.



- An ANOVA table and a Measure of Association table are produced if additional statistics are requested.

## Example

### Creating a Grouped Summary Report

```
SUMMARIZE
  /TABLES=value BY street
  /FORMAT=NOLIST TOTAL
  /TITLE='Home Sale Statistics' /FOOTNOTE 'Grouped by Street'
  /MISSING=VARIABLE
  /CELLS=COUNT MEAN MEDIAN MIN MAX .
```

- The procedure produces summaries of *value* by values of *street*.
- `FORMAT` specifies that no case listing should be produced but that summary statistics should be produced.
- `TITLE` produces the requested title for the summary report and `FOOTNOTE` produces the specified table footnote.
- `CELLS` requests the cell counts, means, medians, minimums, and maximums.
- All other options are set to their default values.

### Displaying Individual Cases

```
SUMMARIZE
  /TABLES=address value marktime
  /FORMAT=VALIDLIST NOCASENUM MISSING='Unavailable' TOTAL
  /TITLE='Case Listing' /FOOTNOTE 'Only homes selling for $335,000 or more'
  /MISSING=VARIABLE
  /CELLS=COUNT MEAN MEDIAN .
```

- The procedure produces summaries of *address*, *value*, and *marktime*.
- `FORMAT` specifies that the case listing report should contain only valid cases, without the case numbers, and with system-missing values displayed with the label *Unavailable*.
- `TITLE` produces the requested title for the summary report and `FOOTNOTE` produces the specified table footnote.
- `CELLS` requests the cell counts, means, and medians.
- All other options are set to their default values.

## TABLES Subcommand

`TABLES` specifies the dependent and control variables.

- You can specify multiple `TABLES` subcommands on a single `SUMMARIZE` command.
- For `FORMAT=LIST` or `VALIDLIST`, only one `TABLES` subcommand is allowed. Multiple dependent and control variables add more breaks to the Listing Report. Total statistics are displayed at the end for each combination that is defined by different values of the control variables.

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- For `FORMAT=NOLIST`, which is the default, each use of keyword `BY` adds a dimension to the requested table. Total statistics are displayed with each group.
- The order in which control variables are displayed is the same as the order in which they are specified on `TABLES`. The values of the first control variable that is defined for the table appear in the leftmost column of the table and change the most slowly in the definition of groups.
- Statistics are displayed for each dependent variable in the same report.
- More than one dependent variable can be specified in a table list, and more than one control variable can be specified in each dimension of a table list.

## ***TITLE and FOOTNOTE Subcommands***

`TITLE` and `FOOTNOTE` provide a title and a caption for the Summary or Listing Report.

- `TITLE` and `FOOTNOTE` are optional and can be placed anywhere.
- The specification on `TITLE` or `FOOTNOTE` is a string enclosed in quotes. To specify a multiple-line title or footnote, enclose each line in quotes and separate the specifications for each line by at least one blank.
- The string that you specify cannot exceed 255 characters.

## ***CELLS Subcommand***

By default, `SUMMARIZE` displays the means, standard deviations, and cell counts in each cell. Use `CELLS` to modify cell information.

- If `CELLS` is specified without keywords, `SUMMARIZE` displays the default statistics.
- If any keywords are specified on `CELLS`, only the requested information is displayed.
- `MEDIAN` and `GMEDIAN` use a lot of computer resources and time. Requesting these statistics (via these keywords or `ALL`) may slow down performance.

<b>DEFAULT</b>	<i>Means, standard deviations, and cell counts.</i> This setting is the default if <code>CELLS</code> is omitted.
<b>MEAN</b>	<i>Cell means.</i>
<b>STDDEV</b>	<i>Cell standard deviations.</i>
<b>COUNT</b>	<i>Cell counts.</i>
<b>MEDIAN</b>	<i>Cell median.</i>
<b>GMEDIAN</b>	<i>Grouped median.</i>
<b>SEMEAN</b>	<i>Standard error of cell mean.</i>
<b>SUM</b>	<i>Cell sums.</i>
<b>MIN</b>	<i>Cell minimum.</i>
<b>MAX</b>	<i>Cell maximum.</i>
<b>RANGE</b>	<i>Cell range.</i>
<b>VARIANCE</b>	<i>Variances.</i>
<b>KURT</b>	<i>Cell kurtosis.</i>
<b>SEKURT</b>	<i>Standard error of cell kurtosis.</i>
<b>SKEW</b>	<i>Cell skewness.</i>
<b>SESKEW</b>	<i>Standard error of cell skewness.</i>

<b>FIRST</b>	<i>First value.</i>
<b>LAST</b>	<i>Last value.</i>
<b>SPCT</b>	<i>Percentage of total sum.</i>
<b>NPCT</b>	<i>Percentage of total number of cases.</i>
<b>SPCT(var)</b>	<i>Percentage of total sum within specified variable. The specified variable must be one of the control variables.</i>
<b>NPCT(var)</b>	<i>Percentage of total number of cases within specified variable. The specified variable must be one of the control variables.</i>
<b>HARMONIC</b>	<i>Harmonic mean.</i>
<b>GEOMETRIC</b>	<i>Geometric mean.</i>
<b>ALL</b>	<i>All cell information.</i>

## ***MISSING Subcommand***

MISSING controls the treatment of cases with missing values. There are two groups of keywords.

- EXCLUDE, INCLUDE, and DEPENDENT specify the treatment of user-missing values. The default is EXCLUDE.

<b>EXCLUDE</b>	<i>All user-missing values are excluded. This setting is the default.</i>
<b>INCLUDE</b>	<i>User-missing values are treated as valid values.</i>
<b>DEPENDENT</b>	<i>User-missing values are considered missing in a dependent variable and valid in a grouping variable (variables that are specified after a BY keyword).</i>

- VARIABLE and TABLE specify how cases with missing values for dependent variables are excluded. The default is VARIABLE.
- Cases with missing values for any control variables are always excluded.

<b>VARIABLE</b>	<i>A case is excluded when all values for variables in that table are missing. This setting is the default.</i>
<b>TABLE</b>	<i>A case is excluded when any value is missing within a table.</i>

## ***FORMAT Subcommand***

FORMAT specifies whether you want a case listing for your report and whether you want case numbers displayed for the listing. FORMAT also determines whether your reports will display a total category for each group and how the reports will indicate missing values.

<b>NOLIST</b>	<i>Display a Summary Report without a case listing. This setting is the default.</i>
<b>LIST</b>	<i>Display a Listing Report. By default, LIST shows all cases. It can be followed by the optional keyword LIMIT, an equals sign, and a positive integer to limit the number of cases in the report. For example, /FORMAT LIST LIMIT=10 will limit the report to the first 10 cases.</i>
<b>VALIDLIST</b>	<i>Display a Listing Report showing only valid cases.</i>
<b>CASENUM</b>	<i>Display case numbers as a category in the Listing Reports. This setting is the default when FORMAT=LIST or VALIDLIST.</i>
<b>NOCASENUM</b>	<i>Do not display case numbers.</i>

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<b>TOTAL</b>	<i>Display the summary statistics for the total of each group with the label Total. This is the default.</i>
<b>NOTOTAL</b>	<i>Display the total category without a label.</i>
<b>MISSING='string'</b>	<i>Display system-missing values as a specified string.</i>

**STATISTICS Subcommand**

Use **STATISTICS** to request a one-way analysis of variance and a test of linearity for each table list.

- Statistics that are requested on **STATISTICS** are computed in addition to the statistics that are displayed in the Group Statistics table.
- If **STATISTICS** is specified without keywords, **SUMMARIZE** computes ANOVA.
- If two or more dimensions are specified, the second and subsequent dimensions are ignored in the analysis-of-variance table.

<b>ANOVA</b>	<i>Analysis of variance.</i> ANOVA displays a standard analysis-of-variance table and calculates eta and eta squared (displayed in the Measures of Association table). This setting is the default if <b>STATISTICS</b> is specified without keywords.
<b>LINEARITY</b>	<i>Test of linearity.</i> LINEARITY (alias ALL) displays additional statistics to the tables that are created by the ANOVA keyword—the sums of squares, degrees of freedom, and mean square associated with linear and nonlinear components, the <i>F</i> ratio, and the significance level for the ANOVA table and Pearson's <i>r</i> and <i>r</i> <sup>2</sup> for the Measures of Association table. LINEARITY is ignored if the control variable is a string.
<b>NONE</b>	<i>No additional statistics.</i> This setting is the default if <b>STATISTICS</b> is omitted.

**Example**

```
SUMMARIZE TABLES=INCOME BY SEX BY RACE
  /STATISTICS=ANOVA.
```

- **SUMMARIZE** produces a Group Statistics table of *INCOME* by *RACE* within *SEX* and computes an analysis of variance only for *INCOME* by *SEX*.

# SURVIVAL

SURVIVAL is available in the Advanced Statistics option.

```
SURVIVAL TABLE=survival varlist
                [BY varlist (min, max)...][BY varlist (min, max)...]

/INTERVAL=THRU n BY a [THRU m BY b ...]

/STATUS=status variable({min, max}) FOR {ALL      }
                {value  } {survival varlist}
[/STATUS=...]

[/PLOTS  ({ALL      })={ALL      } BY {ALL      }      BY {ALL      }]
        {LOGSURV } {survival varlis} {varlist}      {varlist}
        {SURVIVAL}
        {HAZARD  }
        {DENSITY }
        {OMS     }

[/PRINT={TABLE**}]
        {NOTABLE}

[/COMPARE={ALL**      } BY {ALL**      }      BY {ALL**      }]
        {survival varlist} {varlist}      {varlist}

[/CALCULATE=[{EXACT**      } [PAIRWISE] [COMPARE] ]
        {CONDITIONAL}
        {APPROXIMATE}]

[/MISSING={GROUPWISE**} [INCLUDE] ]
        {LISTWISE      }

[/WRITE=[{NONE**}] ]
        {TABLES}
        {BOTH      }
```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
SURVIVAL TABLE=MOSFREE BY TREATMNT(1,3)
/STATUS = PRISON (1) FOR MOSFREE
/INTERVAL=THRU 24 BY 3.
```

## Overview

SURVIVAL produces actuarial life tables, plots, and related statistics for examining the length of time to the occurrence of an event, often known as **survival time**. Cases can be classified into groups for separate analyses and comparisons. Time intervals can be calculated with the IBM® SPSS® Statistics date- and time-conversion functions—for example, `CTIME.DAYS` or `YRMODE`. For a closely related alternative nonparametric analysis of survival times using the product-limit Kaplan-Meier estimator, see the `KM` command. For an analysis of survival times with covariates, including time-dependent covariates, see the `COXREG` command.

### **Options**

**Life Tables.** You can list the variables to be used in the analysis, including any control variables on the `TABLE` subcommand. You can also suppress the life tables in the output with the `PRINT` subcommand.

**Intervals.** `SURVIVAL` reports the percentage alive at various times after the initial event. You can select the time points for reporting with the `INTERVAL` subcommand.

**Plots.** You can plot the survival functions for all cases or separately for various subgroups with the `PLOTS` subcommand.

**Comparisons.** When control variables are listed on the `TABLE` subcommand, you can compare groups based on the Wilcoxon (Gehan) statistic using the `COMPARE` subcommand. You can request pairwise or approximate comparisons with the `CALCULATE` subcommand.

**Writing a File.** You can write the life tables, including the labeling information, to a file with the `WRITE` subcommand.

### **Basic Specification**

- The basic specification requires three subcommands: `TABLE`, `INTERVAL`, and `STATUS`. `TABLE` identifies at least one survival variable from the active dataset, `INTERVAL` divides the time period into intervals, and `STATUS` names a variable that indicates whether the event occurred.
- The basic specification prints one or more life tables, depending on the number of survival and control variables specified.

### **Subcommand Order**

- `TABLE` must be first.
- Remaining subcommands can be named in any order.

### **Syntax Rules**

- Only one `TABLE` subcommand can be specified, but multiple survival variables can be named. A survival variable cannot be specified as a control variable on any subcommands.
- Only one `INTERVAL` subcommand can be in effect on a `SURVIVAL` command. The interval specifications apply to all of the survival variables listed on `TABLE`. If multiple `INTERVAL` subcommands are used, the last specification supersedes all previous ones.
- Only one status variable can be listed on each `STATUS` subcommand. To specify multiple status variables, use multiple `STATUS` subcommands.
- You can specify multiple control variables on one `BY` keyword. Use a second `BY` keyword to specify second-order control variables to interact with the first-order control variables.
- All variables, including survival variables, control variables, and status variables, must be numeric. `SURVIVAL` does not process string variables.

**Operations**

- SURVIVAL computes time intervals according to specified interval widths, calculates the survival functions for each interval, and builds one life table for each group of survival variables. The life table is displayed unless explicitly suppressed.
- When the PLOTS subcommand is specified, SURVIVAL plots the survival functions for all cases or separately for various groups.
- When the COMPARE subcommand is specified, SURVIVAL compares survival-time distributions of different groups based on the Wilcoxon (Gehan) statistic.

**Limitations**

- Maximum 20 survival variables.
- Maximum 100 control variables total on the first- and second-order control-variable lists combined.
- Maximum 20 THRU and BY specifications on INTERVAL.
- Maximum 35 values can appear on a plot.

**Examples**

```
SURVIVAL TABLE=MOSFREE BY TREATMNT(1,3)
  /STATUS = PRISON (1) FOR MOSFREE
  /INTERVAL = THRU 24 BY 3.
```

- The survival analysis is used to examine the length of time between release from prison and return to prison for prisoners in three treatment programs. The variable *MOSFREE* is the length of time in months a prisoner stayed out of prison. The variable *TREATMNT* indicates the treatment group for each case.
- A value of 1 on the variable *PRISON* indicates a terminal outcome—that is, cases coded as 1 have returned to prison. Cases with other non-negative values for *PRISON* have not returned. Because we don't know their final outcome, such cases are called censored.
- Life tables are produced for each of the three subgroups. *INTERVAL* specifies that the survival experience be described every three months for the first two years.

**Using Life Tables to Examine Customer Time to Churn**

```
SURVIVAL
  TABLE=tenure BY custcat(1 4)
  /INTERVAL=THRU 60 BY 3
  /STATUS=churn(1)
  /PRINT=TABLE
  /PLOTS ( SURVIVAL )=tenure BY custcat
  /COMPARE=tenure BY custcat /CALCULATE PAIRWISE .
```

- The survival analysis is used to examine the number of months before a customer churns. The variable *tenure* is the length of time in months a person was a customer. The variable *custcat* indicates the customer category for each case.
- A value of 1 on the variable *churn* indicates a terminal outcome—that is, cases coded as 1 have churned. Cases with other non-negative values for *churn* have not returned. Because we don't know their final outcome, such cases are called censored.

- Life tables are produced for each of the four subgroups. The `INTERVAL` subcommand specifies that the survival experience be described every three months for the first five years.
- The `PLOTS` subcommand produces plots of the cumulative survival distribution for *tenure* for the four values of *custcat*.
- The `COMPARE` and `CALCULATE` subcommands produce pairwise comparisons of the survival curves for *tenure* by the values of *custcat*.

## TABLE Subcommand

`TABLE` identifies the survival and control variables to be included in the analysis.

- The minimum specification is one or more survival variables.
- To specify one or more first-order control (or factor) variables, use the keyword `BY` followed by the control variable(s). First-order control variables are processed in sequence. For example, `BY A(1,3) B(1,2)` results in five groups (A=1, A=2, A=3, B=1, and B=2).
- You can specify one or more second-order control variables following a second `BY` keyword. Separate life tables are generated for each combination of values of the first-order and second-order controls. For example, `BY A(1,3) BY B(1,2)` results in six groups (A=1 B=1, A=1 B=2, A=2 B=1, A=2 B=2, A=3 B=1, and A=3 B=2).
- Each control variable must be followed by a value range in parentheses. These values must be integers separated by a comma or a blank. Non-integer values in the data are truncated, and the case is assigned to a subgroup based on the integer portion of its value on the variable. To specify only one value for a control variable, use the same value for the minimum and maximum.
- To generate life tables for all cases combined, as well as for control variables, use `COMPUTE` to create a variable that has the same value for all cases. With this variable as a control, tables for the entire set of cases, as well as for the control variables, will be produced.

### Example

```
SURVIVAL TABLE = MOSFREE BY TREATMNT(1,3) BY RACE(1,2)
/STATUS = PRISON(1)
/INTERVAL = THRU 24 BY 3.
```

- *MOSFREE* is the survival variable, and *TREATMNT* is the first-order control variable. The second `BY` defines *RACE* as a second-order control group having a value of 1 or 2.
- Six life tables with the median survival time are produced, one for each pair of values for the two control variables.

## INTERVAL Subcommand

`INTERVAL` determines the period of time to be examined and how the time will be grouped for the analysis. The interval specifications apply to all of the survival variables listed on `TABLE`.

- `SURVIVAL` always uses 0 as the starting point for the first interval. Do not specify the 0. The `INTERVAL` specification *must* begin with the keyword `THRU`.



- Specify the terminal value of the time period after the keyword `THRU`. The final interval includes any observations that exceed the specified terminal value.
- The grouping increment, which follows the keyword `BY`, must be in the same units as the survival variable.
- The period to be examined can be divided into intervals of varying lengths by repeating the `THRU` and `BY` keywords. The period must be divided in ascending order. If the time period is not a multiple of the increment, the endpoint of the period is adjusted upward to the next even multiple of the grouping increment.
- When the period is divided into intervals of varying lengths by repeating the `THRU` and `BY` specifications, the adjustment of one period to produce even intervals changes the starting point of subsequent periods. If the upward adjustment of one period completely overlaps the next period, no adjustment is made and the procedure terminates with an error.

### **Example**

```
SURVIVAL TABLE = MOSFREE BY TREATMNT(1,3)
  /STATUS = PRISON(1) FOR MOSFREE
  /INTERVAL = THRU 12 BY 1 THRU 24 BY 3.
```

- `INTERVAL` produces life tables computed from 0 to 12 months at one-month intervals and from 13 to 24 months at three-month intervals.

### **Example**

```
SURVIVAL ONSSURV BY TREATMNT (1,3)
  /STATUS = OUTCOME (3,4) FOR ONSSURV
  /INTERVAL = THRU 50 BY 6.
```

- On the `INTERVAL` subcommand, the value following `BY` (6) does not divide evenly into the period to which it applies (50). Thus, the endpoint of the period is adjusted upward to the next even multiple of the `BY` value, resulting in a period of 54 with 9 intervals of 6 units each.

### **Example**

```
SURVIVAL ONSSURV BY TREATMNT (1,3)
  /STATUS = OUTCOME (3,4) FOR ONSSURV
  /INTERVAL = THRU 50 BY 6 THRU 100 BY 10 THRU 200 BY 20.
```

- Multiple `THRU` and `BY` specifications are used on the `INTERVAL` subcommand to divide the period of time under examination into intervals of different lengths.
- The first `THRU` and `BY` specifications are adjusted to produce even intervals as in the previous example. As a result, the following `THRU` and `BY` specifications are automatically readjusted to generate 5 intervals of 10 units (through 104), followed by 5 intervals of 20 units (through 204).

## **STATUS Subcommand**

To determine whether the terminal event has occurred for a particular observation, `SURVIVAL` checks the value of a status variable. `STATUS` lists the status variable associated with each survival variable and the codes that indicate that a terminal event occurred.

- Specify a status variable followed by a value range enclosed in parentheses. The value range identifies the codes that indicate that the terminal event has taken place. All cases with non-negative times that do not have a code in the value range are classified as **censored cases**, which are cases for which the terminal event has not yet occurred.
- If the status variable does not apply to all the survival variables, specify `FOR` and the name of the survival variable(s) to which the status variable applies.
- Each survival variable on `TABLE` must have an associated status variable identified by a `STATUS` subcommand.
- Only one status variable can be listed on each `STATUS` subcommand. To specify multiple status variables, use multiple `STATUS` subcommands.
- If `FOR` is omitted on the `STATUS` specification, the status-variable specification applies to all of the survival variables not named on another `STATUS` subcommand.
- If more than one `STATUS` subcommand omits the keyword `FOR`, the final `STATUS` subcommand without `FOR` applies to all survival variables not specified by `FOR` on other `STATUS` subcommands. No warning is printed.

**Example**

```
SURVIVAL ONSSURV BY TREATMNT (1,3)
  /INTERVAL = THRU 50 BY 5, THRU 100 BY 10
  /STATUS = OUTCOME (3,4) FOR ONSSURV.
```

- `STATUS` specifies that a code of 3 or 4 on *OUTCOME* means that the terminal event for the survival variable *ONSSURV* occurred.

**Example**

```
SURVIVAL TABLE = NOARREST MOSFREE BY TREATMNT(1,3)
  /STATUS = ARREST (1) FOR NOARREST
  /STATUS = PRISON (1)
  /INTERVAL=THRU 24 BY 3.
```

- `STATUS` defines the terminal event for *NOARREST* as a value of 1 for *ARREST*. Any other value for *ARREST* is considered censored.
- The second `STATUS` subcommand defines the value of 1 for *PRISON* as the terminal event. The keyword `FOR` is omitted. Thus, the status-variable specification applies to *MOSFREE*, which is the only survival variable not named on another `STATUS` subcommand.

**PLOTS Subcommand**

`PLOTS` produces plots of the cumulative survival distribution, the hazard function, and the probability density function. The `PLOTS` subcommand can plot only the survival functions generated by the `TABLE` subcommand; `PLOTS` cannot eliminate control variables.

- When specified by itself, the `PLOTS` subcommand produces all available plots for each survival variable. Points on each plot are identified by values of the first-order control variables. If second-order controls are used, a separate plot is generated for every value of the second-order control variables.

- To request specific plots, specify, in parentheses following PLOTS, any combination of the keywords defined below.
- Optionally, generate plots for only a subset of the requested life tables. Use the same syntax as used on the TABLE subcommand for specifying survival and control variables, omitting the value ranges. Each survival variable named on PLOTS must have as many control levels as were specified for that variable on TABLE. However, only one control variable needs to be present for each level. If a required control level is missing on the PLOTS specification, the default BY ALL is used for that level. The keyword ALL can be used to refer to an entire set of survival or control variables.
- To determine the number of plots that will be produced, multiply the number of functions plotted by the number of survival variables times the number of first-order controls times the number of distinct values represented in all of the second-order controls.

<b>ALL</b>	<i>Plot all available functions. ALL is the default if PLOTS is used without specifications.</i>
<b>LOGSURV</b>	<i>Plot the cumulative survival distribution on a logarithmic scale.</i>
<b>SURVIVAL</b>	<i>Plot the cumulative survival distribution on a linear scale.</i>
<b>HAZARD</b>	<i>Plot the hazard function.</i>
<b>DENSITY</b>	<i>Plot the density function.</i>
<b>OMS</b>	<i>Plot the one-minus-survival function.</i>

### Example

```
SURVIVAL TABLE = NOARREST MOSFREE BY TREATMNT(1,3)
  /STATUS = ARREST (1) FOR NOARREST
  /STATUS = PRISON (1) FOR MOSFREE
  /INTERVAL = THRU 24 BY 3
  /PLOTS (SURVIVAL,HAZARD) = MOSFREE.
```

- Separate life tables are produced for each of the survival variables (*NOARREST* and *MOSFREE*) for each of the three values of the control variable *TREATMNT*.
- PLOTS produces plots of the cumulative survival distribution and the hazard rate for *MOSFREE* for the three values of *TREATMNT* (even though *TREATMNT* is not included on the PLOTS specification).
- Because plots are requested only for the survival variable *MOSFREE*, no plots are generated for the variable *NOARREST*.

## PRINT Subcommand

By default, SURVIVAL prints life tables. PRINT can be used to suppress the life tables.

<b>TABLE</b>	<i>Print the life tables. This is the default.</i>
<b>NOTABLE</b>	<i>Suppress the life tables. Only plots and comparisons are printed. The WRITE subcommand, which is used to write the life tables to a file, can be used when NOTABLE is in effect.</i>

### Example

```
SURVIVAL TABLE = MOSFREE BY TREATMNT(1,3)
```

```

/STATUS = PRISON (1) FOR MOSFREE
/INTERVAL = THRU 24 BY 3
/PLOTS (ALL)
/PRINT = NOTABLE.

```

- `PRINT NOTABLE` suppresses the printing of life tables.

## **COMPARE Subcommand**

`COMPARE` compares the survival experience of subgroups defined by the control variables. At least one first-order control variable is required for calculating comparisons.

- When specified by itself, the `COMPARE` subcommand produces comparisons using the `TABLE` variable list.
- Alternatively, specify the survival and control variables for the comparisons. Use the same syntax as used on the `TABLE` subcommand for specifying survival and control variables, omitting the value ranges. Only variables that appear on the `TABLE` subcommand can be listed on `COMPARE`, and their role as survival, first-order, and second-order control variables cannot be altered. The keyword `TO` can be used to refer to a group of variables, and the keyword `ALL` can be used to refer to an entire set of survival or control variables.
- By default, `COMPARE` calculates exact comparisons between subgroups. Use the `CALCULATE` subcommand to obtain pairwise comparisons or approximate comparisons.

### **Example**

```

SURVIVAL TABLE = MOSFREE BY TREATMNT(1,3)
/STATUS = PRISON (1) FOR MOSFREE
/INTERVAL = THRU 24 BY 3
/COMPARE.

```

- `COMPARE` computes the Wilcoxon (Gehan) statistic, degrees of freedom, and observed significance level for the hypothesis that the three survival curves based on the values of `TREATMNT` are identical.

### **Example**

```

SURVIVAL TABLE=ONSSURV,RECSURV BY TREATMNT(1,3)
/STATUS = RECURSIT(1,9) FOR RECSURV
/STATUS = STATUS(3,4) FOR ONSSURV
/INTERVAL = THRU 50 BY 5 THRU 100 BY 10
/COMPARE = ONSSURV BY TREATMNT.

```

- `COMPARE` requests a comparison of `ONSSURV` by `TREATMNT`. No comparison is made of `RECSURV` by `TREATMNT`.

## **CALCULATE Subcommand**

`CALCULATE` controls the comparisons of survival for subgroups specified on the `COMPARE` subcommand.

- The minimum specification is the subcommand keyword by itself. `EXACT` is the default.

- Only one of the keywords `EXACT`, `APPROXIMATE`, and `CONDITIONAL` can be specified. If more than one keyword is used, only one is in effect. The order of precedence is `APPROXIMATE`, `CONDITIONAL`, and `EXACT`.
- The keywords `PAIRWISE` and `COMPARE` can be used with any of the `EXACT`, `APPROXIMATE`, or `CONDITIONAL` keywords.
- If `CALCULATE` is used without the `COMPARE` subcommand, `CALCULATE` is ignored. However, if the keyword `COMPARE` is specified on `CALCULATE` and the `COMPARE` subcommand is omitted, an error message is generated.
- Data can be entered into `SURVIVAL` for each individual case or aggregated for all cases in an interval. The way in which data are entered determines whether an exact or an approximate comparison is most appropriate. For more information, see the topic [Using Aggregated Data on p. 1952](#).

<b>EXACT</b>	<i>Calculate exact comparisons.</i> This is the default. You can obtain exact comparisons based on the survival experience of each observation with individual data. While this method is the most accurate, it requires that all of the data be in memory simultaneously. Thus, exact comparisons may be impractical for large samples. It is also inappropriate when individual data are not available and data aggregated by interval must be used.
<b>APPROXIMATE</b>	<i>Calculate approximate comparisons only.</i> Approximate comparisons are appropriate for aggregated data. The approximate-comparison approach assumes that all events occur at the midpoint of the interval. With exact comparisons, some of these midpoint ties can be resolved. However, if interval widths are not too great, the difference between exact and approximate comparisons should be small.
<b>CONDITIONAL</b>	<i>Calculate approximate comparisons if memory is insufficient.</i> Approximate comparisons are produced only if there is insufficient memory available for exact comparisons.
<b>PAIRWISE</b>	<i>Perform pairwise comparisons.</i> Comparisons of all pairs of values of the first-order control variable are produced along with the overall comparison.
<b>COMPARE</b>	<i>Produce comparisons only.</i> Survival tables specified on the <code>TABLE</code> subcommand are not computed, and requests for plots are ignored. This allows all available workspace to be used for comparisons. The <code>WRITE</code> subcommand cannot be used when this specification is in effect.

### Example

```
SURVIVAL TABLE = MOSFREE BY TREATMNT(1,3)
  /STATUS = PRISON (1) FOR MOSFREE
  /INTERVAL = THRU 24 BY 3
  /COMPARE /CALCULATE = PAIRWISE.
```

- `PAIRWISE` on `CALCULATE` computes the Wilcoxon (Gehan) statistic, degrees of freedom, and observed significance levels for each pair of values of `TREATMNT`, as well as for an overall comparison of survival across all three `TREATMNT` subgroups: group 1 with group 2, group 1 with group 3, and group 2 with group 3.
- All comparisons are exact comparisons.

### Example

```
SURVIVAL TABLE = MOSFREE BY TREATMNT(1,3)
  /STATUS = PRISON (1) FOR MOSFREE
```

```
/INTERVAL = THRU 24 BY 3
/COMPARE /CALCULATE = APPROXIMATE COMPARE.
```

- APPROXIMATE on CALCULATE computes the Wilcoxon (Gehan) statistic, degrees of freedom, and probability for the overall comparison of survival across all three *TREATMNT* subgroups using the approximate method.
- Because the keyword COMPARE is specified on CALCULATE, survival tables are not computed.

## Using Aggregated Data

When aggregated survival information is available, the number of censored and uncensored cases at each time point must be entered. Up to two records can be entered for each interval, one for censored cases and one for uncensored cases. The number of cases included on each record is used as the weight factor. If control variables are used, there will be up to two records (one for censored and one for uncensored cases) for each value of the control variable in each interval. These records must contain the value of the control variable and the number of cases that belong in the particular category as well as values for survival time and status.

### Example

```
DATA LIST / SURVEVAR 1-2 STATVAR 4 SEX 6 COUNT 8.
VALUE LABELS  STATVAR 1 'DECEASED' 2 'ALIVE'
              /SEX 1 'FEMALE' 2 'MALE'.

BEGIN DATA
1 1 1 6
1 1 1 1
1 2 2 2
1 1 2 1
2 2 1 1
2 1 1 2
2 2 2 1
2 1 2 3
. . .
END DATA.
WEIGHT COUNT.
SURVIVAL TABLE = SURVEVAR BY SEX (1,2)
/INTERVAL = THRU 10 BY 1
/STATUS = STATVAR (1) FOR SURVEVAR.
```

- This example reads aggregated data and performs a SURVIVAL analysis when a control variable with two values is used.
- The first data record has a code of 1 on the status variable *STATVAR*, indicating that it is an uncensored case, and a code of 1 on *SEX*, the control variable. The number of cases for this interval is 6, the value of the variable *COUNT*. Intervals with weights of 0 do not have to be included.
- *COUNT* is not used in SURVIVAL but is the weight variable. In this example, each interval requires four records to provide all of the data for each *SURVEVAR* interval.

## MISSING Subcommand

MISSING controls missing-value treatments. The default is GROUPWISE.

- Negative values on the survival variables are automatically treated as missing data. In addition, cases outside of the value range on a control variable are excluded.
- `GROUPWISE` and `LISTWISE` are mutually exclusive. However, each can be used with `INCLUDE`.

<b>GROUPWISE</b>	<i>Exclude missing values groupwise.</i> Cases with missing values on a variable are excluded from any calculation involving that variable. This is the default.
<b>LISTWISE</b>	<i>Exclude missing values listwise.</i> Cases missing on any variables named on <code>TABLE</code> are excluded from the analysis.
<b>INCLUDE</b>	<i>Include user-missing values.</i> User-missing values are included in the analysis.

## WRITE Subcommand

`WRITE` writes data in the survival tables to a file. This file can be used for further analyses or to produce graphics displays.

- When `WRITE` is omitted, the default is `NONE`. No output file is created.
- When `WRITE` is used, a `PROCEDURE OUTPUT` command must precede the `SURVIVAL` command. The `OUTFILE` subcommand on `PROCEDURE OUTPUT` specifies the output file.
- When `WRITE` is specified without a keyword, the default is `TABLES`.

<b>NONE</b>	<i>Do not write procedure output to a file.</i> This is the default when <code>WRITE</code> is omitted.
<b>TABLES</b>	<i>Write survival-table data records.</i> All survival-table statistics are written to a file.
<b>BOTH</b>	<i>Write out survival-table data and label records.</i> Variable names, variable labels, and value labels are written out along with the survival table statistics.

## Format

`WRITE` writes five types of records. The keyword `TABLES` writes record types 30, 31, and 40. The keyword `BOTH` writes record types 10, 20, 30, 31, and 40. The format of each record type is described in the following tables.

Table 236-1  
Record type 10, produced only by keyword `BOTH`

Columns	Content	Format
1–2	Record type (10)	F2.0
3–7	Table number	F5.0
8–15	Name of survival variable	A8
16–55	Variable label of survival variable	A40
56	Number of <code>BY</code> 's (0, 1, or 2)	F1.0
57–60	Number of rows in current survival table	F4.0

- One type-10 record is produced for each life table.

- Column 56 specifies the number of orders of control variables (0, 1, or 2) that have been applied to the life table.
- Columns 57–60 specify the number of rows in the life table. This number is the number of intervals in the analysis that show subjects entering; intervals in which no subjects enter are not noted in the life tables.

Table 236-2

*Record type 20, produced by keyword BOTH*

Columns	Content	Format
1–2	Record type (20)	F2.0
3–7	Table number	F5.0
8–15	Name of control variable	A8
16–55	Variable label of control variable	A40
56–60	Value of control variable	F5.0
61–80	Value label for this value	A20

- One type-20 record is produced for each control variable in each life table.
- If only first-order controls have been placed in the survival analysis, one type-20 record will be produced for each table. If second-order controls have also been applied, two type-20 records will be produced per table.

Table 236-3

*Record type 30, produced by both keywords TABLES and BOTH*

Columns	Content	Format
1–2	Record type (30)	F2.0
3–7	Table number	F5.0
8–13	Beginning of interval	F6.2
14–21	Number entering interval	F8.2
22–29	Number withdrawn in interval	F8.2
30–37	Number exposed to risk	F8.2
38–45	Number of terminal events	F8.2

- Information on record type 30 continues on record type 31. Each pair of type-30 and type-31 records contains the information from one line of the life table.

Table 236-4

*Record type 31, continuation of record type 30*

Columns	Content	Format
1–2	Record type (31)	F2.0
3–7	Table number	F5.0
8–15	Proportion terminating	F8.6
16–23	Proportion surviving	F8.6
24–31	Cumulative proportion surviving	F8.6
32–39	Probability density	F8.6
40–47	Hazard rate	F8.6
48–54	S.E. of cumulative proportion surviving	F7.4
55–61	S.E. of probability density	F7.4
62–68	S.E. of hazard rate	F7.4



- Record type 31 is a continuation of record type 30.
- As many type-30 and type-31 record pairs are output for a table as it has lines (this number is noted in columns 57–60 of the type-10 record for the table).

Table 236-5

Record type 40, produced by both keywords TABLES and BOTH

Columns	Content	Format
1–2	Record type (40)	F2.0

- Type-40 records indicate the completion of the series of records for one life table.

## Record Order

The SURVIVAL output file contains records for each of the life tables specified on the TABLE subcommand. All records for a given table are produced together in sequence. The records for the life tables are produced in the same order as the tables themselves. All life tables for the first survival variable are written first. The values of the first- and second-order control variables rotate, with the values of the first-order controls changing more rapidly.

### Example

```
PROCEDURE OUTPUT OUTFILE = SURVTBL.
SURVIVAL TABLE = MOSFREE BY TREATMNT(1,3)
  /STATUS = PRISON (1) FOR MOSFREE
  /INTERVAL = THRU 24 BY 3
  /WRITE = BOTH.
```

- WRITE generates a procedure output file called *SURVTBL*, containing life tables, variable names and labels, and value labels stored as record types 10, 20, 30, 31, and 40.

# ***SYSFILE INFO***

```
SYSFILE INFO [FILE=] 'file specification'
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## ***Example***

```
SYSFILE INFO FILE='PERSNL.SAV'.
```

## ***Overview***

`SYSFILE INFO` displays complete dictionary information for all variables in an external IBM® SPSS® Statistics data file. You do not have to retrieve the file with `GET` to use `SYSFILE INFO`. If the file has already been retrieved, use `DISPLAY DICTIONARY` to display dictionary information.

### ***Basic Specification***

The basic specification is the command keyword and a complete file specification that is enclosed in quotes.

### ***Syntax Rules***

- Only one file specification is allowed per command. To display dictionary information for more than one SPSS Statistics data file, use multiple `SYSFILE INFO` commands.
- The file extension, if there is one, must be specified, even if it is the default.
- The subcommand keyword `FILE` is optional. When `FILE` is specified, the equals sign is required.

### ***Operations***

- No procedure is needed to execute `SYSFILE INFO`, because `SYSFILE INFO` obtains information from the dictionary alone.
- `SYSFILE INFO` displays file and variable information. File information includes number of variables and cases in the file; file label; file documents; defined multiple response sets; variable sets; and information that is used by other applications (such as IBM® SPSS® Modeler, Data Entry for Windows, and TextSmart). Variable information includes the variable name; label; sequential position in the file; print and write format; missing values; and value labels for each variable in the specified file.

# TDISPLAY

```
TDISPLAY [{ALL          }]  
          {model names }  
          {command names}  
  
[/TYPE={MODEL**}]  
          {COMMAND}
```

**\*\***Default if the subcommand is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
TDISPLAY.
```

## Overview

TDISPLAY displays information about currently active models. These models are automatically generated by a number of procedures for use with the APPLY subcommand.

## Options

If models are specified on TDISPLAY, information about only those models is displayed. You can use the TYPE subcommand to control whether models are specified by model name or by the name of the procedure that generated them.

## Basic Specification

The basic specification is simply the command keyword TDISPLAY.

- By default, TDISPLAY produces a list of all currently active models. The list includes the model names, the commands that created each model, model labels (if specified), and creation dates and times.

## Syntax Rules

- To display information about a subset of active models, specify those models after TDISPLAY.
- Models can be specified by using individual model names or the names of the procedures that created them. To use procedure names, you must specify the TYPE subcommand with the keyword COMMAND.
- Model names are either the default MOD\_n names or the names that are assigned with MODEL NAME.
- If procedure names are specified, all models that are created by those procedures are displayed.
- Model names and procedure names cannot be mixed on the same TDISPLAY command.

- You can specify the keyword `ALL` after `TDISPLAY` to display all models that are currently active. This setting is the default.

### **Operations**

- Only currently active models are displayed.
- The following procedures can generate models that can be displayed with the `TDISPLAY` command: `AREG`, `ARIMA`, `EXSMOOTH`, `SEASON`, and `SPECTRA` in the Forecasting add-on module; `ACF`, `CASEPLOT`, `CCF`, `CURVEFIT`, `PACF`, `PLOT`, and `TSLOT` in the Core system; and `WLS` and `2SLS` in the Regression add-on module.

## **TYPE Subcommand**

`TYPE` indicates whether models are specified by model name or procedure name.

- One keyword, `MODEL` or `COMMAND`, can be specified after `TYPE`.
- `MODEL` is the default and indicates that models are specified as model names.
- `COMMAND` specifies that models are specified by procedure name.
- `TYPE` has no effect if model names or command names are not listed after `TDISPLAY`.
- If more than one `TYPE` subcommand is specified, only the last subcommand is used.
- The `TYPE` specification applies only to the current `TDISPLAY` command.

### **Example**

```
TDISPLAY ACF CURVEFIT  
/TYPE=COMMAND.
```

- This command displays all currently active models that were created by procedures `ACF` and `CURVEFIT`.

# TEMPORARY

TEMPORARY

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## Overview

TEMPORARY signals the beginning of temporary transformations that are in effect only for the next procedure. With the exception of variables created or modified by the next procedure (for example, residual values saved as variables by REGRESSION), numeric or string variables created while the TEMPORARY command is in effect are temporary variables, and any modifications made to existing variables while the TEMPORARY command is in effect are also temporary. Any variables created or modified by the next procedure are permanent.

With TEMPORARY, you can perform separate analyses for subgroups in the data and then repeat the analysis for the file as a whole. You can also use TEMPORARY to transform data for one analysis but not for other subsequent analyses.

TEMPORARY can be applied to the following commands:

- Transformation commands COMPUTE, RECODE, IF, and COUNT, and the DO REPEAT utility.
- The LOOP and DO IF structures.
- Format commands PRINT FORMATS, WRITE FORMATS, and FORMATS.
- Data selection commands SELECT IF, SAMPLE, FILTER, and WEIGHT.
- Variable declarations NUMERIC, STRING, and VECTOR.
- Labeling commands VARIABLE LABELS and VALUE LABELS, and the MISSING VALUES command.
- SPLIT FILE.
- XSAVE.

### Basic specification

The only specification is the keyword TEMPORARY. There are no additional specifications.

### Operations

- Once TEMPORARY is specified, you cannot refer to previously existing scratch variables.
- Temporary transformations apply to the next command that reads the data. Once the data are read, the temporary transformations are no longer in effect.
- The XSAVE command leaves temporary transformations in effect. SAVE, however, reads the data and turns temporary transformations off after the file is written. (See “Examples.”)

- TEMPORARY cannot be used with SORT CASES, MATCH FILES, ADD FILES, or COMPUTE with a LAG function. If any of these commands follows TEMPORARY in the command sequence, there must be an intervening procedure or command that reads the data to first execute the TEMPORARY command.
- TEMPORARY cannot be used within the DO IF-END IF or LOOP-END LOOP structures.

If any variables have an unknown measurement level, the measurement level for those variables remains unknown after a procedure that follows a TEMPORARY command. (In the absence of the TEMPORARY command, the data pass performed by the procedure would assign default measurement levels to all variables with an unknown measurement level.)

## Examples

### Basic example

```
SORT CASES BY SEX.
TEMPORARY.
SPLIT FILE BY SEX.
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.
FREQUENCIES VARS=INCOME /STATISTICS=MEDIAN.
```

- SPLIT FILE applies to the first FREQUENCIES procedure, which generates separate median income tables for men and women.
- SPLIT FILE is not in effect for the second FREQUENCIES procedure, which generates a single median income table that includes both men and women.

### Temporary transformations for a new variable

```
DATA LIST FILE=HUBDATA RECORDS=3
  /1 #MOBIRTH #DABIRTH #YRBIRTH 6-11 DEPT88 19.
COMPUTE AGE=($JDATE - YRMODA(#YRBIRTH,#MOBIRTH,#DABIRTH))/365.25.
VARIABLE LABELS AGE 'EMPLOYEE'S AGE'
  DEPT88 'DEPARTMENT CODE IN 1988'.

TEMPORARY.
RECODE AGE (LO THRU 20=1)(20 THRU 25=2)(25 THRU 30=3)(30 THRU 35=4)
  (35 THRU 40=5)(40 THRU 45=6)(45 THRU 50=7)(50 THRU 55=8)
  (55 THRU 60=9)(60 THRU 65=10)(65 THRU HI=11).
VARIABLE LABELS AGE 'EMPLOYEE AGE CATEGORIES'.
VALUE LABELS AGE 1 'Up to 20' 2 '20 to 25' 3 '25 to 30' 4 '30 to 35'
  5 '35 to 40' 6 '40 to 45' 7 '45 to 50' 8 '50 to 55'
  9 '55 to 60' 10 '60 to 65' 11 '65 and older'.

FREQUENCIES VARIABLES=AGE.
MEANS AGE BY DEPT88.
```

- COMPUTE creates variable AGE from the dates in the data.
- FREQUENCIES uses the temporary version of variable AGE with temporary variable and value labels.
- MEANS uses the unrecoded values of AGE and the permanent variable label.

**Using XSAVE with TEMPORARY**

```
GET FILE='/data/hub.sav'.
TEMPORARY.
RECODE DEPT85 TO DEPT88 (1,2=1) (3,4=2) (ELSE=9).
VALUE LABELS DEPT85 TO DEPT88 1 'MANAGEMENT'
                                2 'OPERATIONS'
                                3 'UNKNOWN'.
XSAVE OUTFILE='/temp/hubtemp.sav'.
CROSSTABS DEPT85 TO DEPT88 BY JOBCAT.
```

- Both the saved data file and the CROSSTABS output will reflect the temporary recoding and labeling of the department variables.
- If XSAVE is replaced with SAVE, the saved data file will reflect the temporary recoding and labeling but the CROSSTABS output will not.

# TIME PROGRAM

```
TIME PROGRAM.  
[commands to compute time dependent covariates]
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## Overview

TIME PROGRAM is required to define time-dependent covariates. These are covariates whose values change during the course of the study.

### Basic Specification

The basic specification is the TIME PROGRAM command followed by one or more commands that define time-dependent covariates.

### Syntax Rules

- TIME PROGRAM and the transformations that define the time-dependent covariate(s) must precede the command that analyzes the time-dependent covariate(s).
- A time-dependent covariate is a function of the current time, which is represented by the special variable  $T_.$
- The active dataset must not have a variable named  $T_.$  If it does, rename the variable. Otherwise, you will trigger an error.
- $T_.$  cannot be specified as a covariate. Any other variable in the TIME PROGRAM can be specified on the covariate list.
- For every time-dependent covariate, values are generated for each valid case for all uncensored times in the same stratum that occur before the observed time. If no STRATA subcommand is specified, all cases are considered to belong to one stratum.
- If any function defined by the time program results in a missing value for a case that has no missing values for any other variable used in the procedure, this causes an error.

## Example

```
TIME PROGRAM.  
COMPUTE Z=AGE + T_.  
  
COXREG VARIABLES = SURVIVAL WITH Z  
/STATUS SURVSTA EVENT (1).
```

- TIME PROGRAM defines the time-dependent covariate  $Z$  as the current age.  $Z$  is then specified as a covariate in COXREG.



# TITLE

```
TITLE [']text[']
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
TITLE "Running Shoe Study from Runner's World Data".
```

## **Overview**

TITLE inserts a left-justified title on the top line of each page of output. The default title indicates the version of the system being used.

### **Basic Specification**

The only specification is the title.

### **Syntax Rules**

- The title can include any characters. To specify a blank title, enclose a blank in quotes (“ ”).
- The title can be up to 60 bytes long. Titles longer than 60 bytes are truncated.
- The quotes enclosing the title are optional; using them allows you to include apostrophes or quotation marks in the title. [For more information, see the topic String Values in Command Specifications on p. 40.](#)
- More than one TITLE command is allowed in a single session.
- A title cannot be placed between a procedure command and BEGIN DATA-END DATA or within data records when the data are inline.

### **Operations**

- The title is displayed as part of the output heading, which also includes the date and page number. If HEADER=NO is specified on SET, the heading, including the title and subtitle, will not be displayed.
- Each TITLE command overrides the previous one and takes effect on the next output page.
- Only the title portion of the heading changes. The date and page number are still displayed.
- TITLE is independent of SUBTITLE, and each can be changed separately.

## **Examples**

```
TITLE My Title.  
TITLE "Fred's Title".  
TITLE 'Fred''s "Quoted" Title'.
```

- The first title doesn't need to be enclosed in quotes because it doesn't contain any quotes or apostrophes.
- The second title is enclosed in double quotes because it contains an apostrophe (single quote).
- The last title is enclosed in single quotes since it contains double quotes, and two single quotes are used to specify the apostrophe.

# TMS BEGIN

```
TMS BEGIN
  /DESTINATION OUTFILE='file specification'
```

## Release History

Release 15.0

- Command introduced.

Release 16.0

- Added support for new string functions `CHAR.CONCAT`, `CHAR.LENGTH`, and `CHAR.SUBSTR` within TMS blocks.

## Example

```
TMS BEGIN
  /DESTINATION OUTFILE='/mydir/mytransformations.xml'.
  COMPUTE modelvar=ln(var).
TMS END.
```

## Overview

The `TMS BEGIN` command indicates the beginning of a block of transformations to be exported to a file in PMML format—specifically, PMML 3.1 with IBM® SPSS® Statistics extensions. The exported transformations can be merged (using `TMS MERGE`) with a PMML model file in order to include any transformations of the raw data required to create the variables used in the model. The merged PMML file can be used to score data with scoring functions. (For more information on scoring functions, see [Scoring expressions](#).) When you score data with a model file containing transformations, the transformations will be applied to the data before scoring. The transformations are carried out as part of the internal scoring process and have no effect on the active dataset. The `TMS END` command is used to end the block.

From the point of view of the active dataset, transformations in a `TMS BEGIN`–`TMS END` block are no different than transformations outside a TMS block—they are pending and will be executed with the next command that reads the active dataset. Source variables used in transformation expressions in TMS blocks must exist in the active dataset. In a typical scenario, you would run `TMS BEGIN`–`TMS END` command syntax on the raw dataset used to build the associated model, assuming it contains the source variables used in any transformation expressions.

Commands within a TMS block are categorized as belonging to one of four types: exported, tolerated, invalid, and those causing an implicit `TMS END`.

**Exported.** These are commands that produce export to PMML. Commands in this category include the supported transformations—for example, `RECODE`. Many transformations are expressed with functions included on a `COMPUTE` command. Export to PMML is supported for a limited set

of such functions. The set of commands that produce export to PMML, as well as the set of supported functions, are provided at the end of this topic.

**Tolerated.** These are commands that can be included in a TMS BEGIN–TMS END block, are executed in normal fashion, but do not produce export to PMML. Examples include NUMERIC, STRING, USE, INCLUDE, and INSERT. In the case of INCLUDE or INSERT, supported transformations from the inserted file will be exported to PMML.

**Invalid.** These are commands that cannot be processed within a TMS block and will give rise to an error. The list of invalid commands is provided at the end of this topic.

**Implicit TMS END.** These commands are executed in normal fashion but cause an implicit end to the current block, which then causes the output PMML file to be written. Transformations (in the current block) occurring after one of these commands are not exported to the output PMML file. Commands in this category include all of the statistical procedures—for example, ANOVA and REGRESSION—as well as the EXECUTE command. ADD FILES and MATCH FILES belong to this category as well as any command that causes a different dataset to become active—for example, DATASET ACTIVATE, GET, and GET DATA.

### ***Basic Specification***

The basic specification for TMS BEGIN is the command name followed by a DESTINATION subcommand that contains an OUTFILE specification.

### ***Syntax Rules***

- The DESTINATION subcommand is required.
- An error occurs if the DESTINATION subcommand is specified more than once.
- An error occurs if the OUTFILE keyword is specified more than once.
- Equal signs (=) shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.

### ***Operations***

- Once a TMS BEGIN command is executed, it remains in effect until the end of the session, until explicitly ended by a TMS END command, or until a command that implicitly ends a TMS block—for example, any statistical procedure—is encountered.
- TMS commands (for example, TMS MERGE) are not allowed within a TMS block.
- TMS BEGIN verifies that the path specified for the destination file is valid. If the path is invalid, the commands in the block will be processed in the normal manner but no PMML will be exported.
- The following conditions will cause the destination file to be written: a command that causes an implicit end of the block is executed; the TMS block ends explicitly with TMS END; the current session ends before the block ends (either explicitly or implicitly). The destination file is unavailable to other commands until after it is written. TMS END has no effect if the block is ended implicitly.

- An error occurs if a transformation in the block is not supported. The command containing the invalid transformation is not executed. This type of error does not cause an implicit end to the TMS block and has no effect on when the destination file is written.
- The order of the transformations in the destination file may not be the order in which they were specified in the TMS block.
- Transformations specified in a TMS block are processed in the normal manner—that is, pending and to be executed with the next command that reads the active dataset.

### ***Exported Commands***

The following commands produce export to PMML.

ADD VALUE LABELS  
COMPUTE  
DO REPEAT - END REPEAT  
MISSING VALUES  
RECODE  
VALUE LABELS  
VARIABLE LABELS  
VARIABLE LEVEL

- For COMPUTE, the target variable must be a variable that has not already been assigned a value or otherwise used in the associated TMS block. For example, the following is invalid:  

```
COMPUTE var = var + 1.
```
- For RECODE, all input and output keywords are supported except CONVERT. In addition, you must always recode INTO a variable that has not already been assigned a value or otherwise used in the associated TMS block. When a recoding transformation is carried out on a dataset to be scored, the results depend on whether the target variable exists in the dataset. If the target variable already exists, cases with values (of the source variable) not mentioned in the value specifications for the recode are not changed. If the target variable does not already exist, cases with values not mentioned are assigned the system-missing value.

### ***Invalid Commands***

The following commands are invalid inside a TMS block.

AGGREGATE  
BREAK  
COUNT  
DO IF - END IF  
ELSE  
ELSE IF  
END CASE  
END FILE  
FILE TYPE - END FILE TYPE  
FILTER

IF  
INPUT PROGRAM - END INPUT PROGRAM  
KEYED DATA LIST  
LEAVE  
LOOP - END LOOP  
N OF CASES  
POINT  
RECORD TYPE  
REPEATING DATA  
REREAD  
SAMPLE  
SELECT IF  
VECTOR  
WEIGHT  
WRITE

### ***Supported Functions, Operations, and System Variables***

The following functions, operations, and system variables are supported for use with the COMPUTE command in TMS blocks. In a few cases—for example, CHAR.SUBSTR—multiple forms (differing in the number of parameters) exist, but only a single form is supported. The parameters available in the supported form are shown.

+ Addition  
- Subtraction  
\* Multiplication  
/ Division  
\*\* Exponentiation  
ABS  
CDF.NORMAL  
CDFNORM  
CHAR.CONCAT  
CHAR.LENGTH  
CHAR.SUBSTR(strexp, pos)  
CONCAT  
EXP  
IDF.NORMAL  
LG10  
LENGTH  
LN  
LOWER  
LTRIM(strexp)  
MAX  
MEAN  
MIN  
MOD

PDF.NORMAL  
 RND  
 RTRIM(strexp)  
 SD  
 SQRT  
 SUBSTR(strexp,pos)  
 SUM  
 SYSMIS  
 TRUNC  
 UPCASE  
 VALUE  
 VARIANCE

### Notes

- The relational operators EQ, NE, LT, LE, GT, and GE are not supported.
- Within a TMS block it is recommended to use CHAR.CONCAT, CHAR.LENGTH, and CHAR.SUBSTR in place of CONCAT, LENGTH, and SUBSTR.
- String arguments for the MAX and MIN functions are not supported for scoring.

## EXAMPLES

```
TMS BEGIN
  /DESTINATION OUTFILE='/mydir/mytransformations.xml'.
  COMPUTE modelvar=ln(var).
TMS END.
```

- TMS BEGIN marks the beginning of a block of transformation commands that will be evaluated for export to PMML.
- The variable *var* is transformed with a log function to create the variable *modelvar*, which is to be used in a model. The details of this transformation will be included as PMML in the file */mydir/mytransformations.xml*.
- TMS END marks the end of the block and causes the output file to be written, but has no effect on the state of the transformations contained in the block. In the present example, the transformation to create *modelvar* is still pending after the completion of the block.

### Tolerated Commands in a TMS Block

```
TMS BEGIN
  /DESTINATION OUTFILE='/mydir/mytransformations.xml'.
  STRING new_strvar (A1).
  RECODE strvar ('A','B','C'='A') ('D','E','F'='B') (ELSE=' ')
  INTO new_strvar.
TMS END.
```

- The STRING command is used to create a new string variable that is to be the target of a recode transformation of an existing string variable. STRING is a tolerated command so it is executed in the normal manner, without generating an error or implicitly ending the TMS block.

**Commands that Cause an Implicit TMS END**

```
TMS BEGIN
  /DESTINATION OUTFILE='/mydir/mytransformations.xml'.
  RECODE numvar1 (0=1) (1=0) (ELSE=SYSMIS) INTO new_numvar1.
  FREQUENCIES new_numvar1.
  RECODE numvar2 (1 THRU 5=1) (6 THRU 10=2) (11 THRU HI=3) (ELSE=0)
    INTO new_numvar2.
TMS END.
```

- The `FREQUENCIES` command causes an implicit end to the `TMS` block, causing the output PMML file to be written. The output file only contains the recode transformation for *numvar1*—that is, the transformations prior to the command causing the implicit `TMS END`.
- Although the `FREQUENCIES` command implicitly ends the block, the recode transformation for *numvar2* is processed in the normal manner and remains pending until the next command that reads the active dataset.
- Although we have included it in this example, the `TMS END` command is not needed when there is a command that implicitly ends the `TMS` block. In this example, execution of the `TMS END` command generates a warning indicating that there is no current `TMS` command in effect.

**Using EXECUTE in a TMS Block**

```
TMS BEGIN
  /DESTINATION OUTFILE='/mydir/mytransformations.xml'.
  RECODE numvar1 (0=1) (1=0) (ELSE=SYSMIS) INTO new_numvar1.
  RECODE numvar2 (1 THRU 5=1) (6 THRU 10=2) (11 THRU HI=3) (ELSE=0)
    INTO new_numvar2.
EXECUTE.
```

- You can include `EXECUTE` in a `TMS` block. In its usual manner, it will cause all pending transformations to be executed. In addition, it causes an implicit end to the block which then causes the output PMML file to be written.
- In the current example, we have omitted the `TMS END` command since the block is implicitly ended by the `EXECUTE` command.

**DESTINATION Subcommand**

The `DESTINATION` subcommand is required. It must include the `OUTFILE` keyword followed by an equals sign (=) and a file specification in quotes or a previously defined file handle defined with the `FILE HANDLE` command. It is recommended to include the extension `.xml` in the file specification.



# TMS END

```
TMS END
  [/PRINT RESULTS={SUMMARY**}]
  {NONE}
```

\*\* Default if the subcommand or keyword is omitted.

## Release History

Release 15.0

- Command introduced.

## Example

```
TMS BEGIN
  /DESTINATION OUTFILE='/mydir/mytransformation.xml'.
  COMPUTE mymodelvar=ln(var).
TMS END
  /PRINT RESULTS=NONE.
```

## Overview

The `TMS END` command is used to indicate the end of a block of transformations to be exported to a file in PMML format—specifically, PMML 3.1 with IBM® SPSS® Statistics extensions. The exported transformations can be merged (using `TMS MERGE`) with a PMML model file in order to include any transformations of the raw data required to create the variables used in the model. The merged PMML file can be used to score data with scoring functions. (For more information on scoring functions, see [Scoring expressions](#).) When you score data with a model file containing transformations, the transformations will be applied to the data before scoring. The transformations are carried out as part of the internal scoring process and have no effect on the active dataset. The `TMS BEGIN` command is used to indicate the beginning of the block.

## Basic Specification

The basic specification for `TMS END` is the command name. For examples of `TMS BEGIN`–`TMS END` blocks, see the `TMS BEGIN` command.

## Syntax Rules

- An error occurs if the `PRINT` subcommand is specified more than once.
- An error occurs if the `RESULTS` keyword is specified more than once.
- Equal signs (=) shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.

**Operations**

- TMS END will generate a warning if there is no TMS BEGIN command in effect. This condition occurs in the obvious case that there is no prior TMS BEGIN command, but will also occur if there is a TMS BEGIN command followed by a command that implicitly ends a TMS block—for instance, any statistical procedure or the EXECUTE command—followed by TMS END.
- TMS END causes the destination file to be written, unless the associated TMS block ended implicitly (see previous bullet), thus causing the file to be written before execution of TMS END. For more information, see the [Examples](#) for the TMS BEGIN command.
- TMS END does not cause transformations in the associated TMS block to be executed. In normal fashion, they are pending and will be executed with the next command that reads the active dataset.

**PRINT Subcommand**

The PRINT subcommand is optional and specifies whether or not to print a table summarizing the derived variables defined by the transformations in the block.

- |                |   |
|----------------|---|
| <b>SUMMARY</b> | <i>Print a summary of the derived variables. This is the default.</i> |
| <b>NONE</b>    | <i>Do not print a summary.</i>  |

# TMS IMPORT

```
TMS IMPORT
  /INFILE TRANSFORMATIONS='filespec'
    [MODE={FORWARD** (ROLES={UPDATE**})          }]
      {LEAVE      }
      {BACK (PREDICTED=fieldname
        SUFFIX={'_backtransformed'}
          {suffixname      } ) }
  [/SAVE TRANSFORMED={YES**}]
    {NO      }
  [/OUTFILE SYNTAX='filespec'].
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 18

- Command introduced.

## **Example**

```
TMS IMPORT
  /INFILE TRANSFORMATIONS='file.xml'.
```

## **Overview**

TMS IMPORT converts transformations contained in an XML file created by ADP (transformations files created by TMS BEGIN-END are not currently “reversible”) into command syntax and applies the converted transformations to the active dataset or saves them to a syntax file.

## **Basic Specification**

The basic specification is the TMS IMPORT command and an INFILE subcommand with a transformations file specification.

## **Syntax Rules**

- The INFILE subcommand is required. All other subcommands are optional.
- Each subcommand can be specified only once.
- Subcommands may be used in any order.
- An error occurs if a keyword is specified more than once within a subcommand.
- Equal signs (=) shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.

## Examples

```
TMS IMPORT
  /INFILE TRANSFORMATIONS='file.xml'.
```

- TMS IMPORT reads the file *file.xml*, converts the transformations from XML to command syntax, and applies them to the active dataset.

### Using Backtransformations

```
TMS IMPORT
  /INFILE TRANSFORMATIONS='workingDirectory/car_sales_transformations.xml'
  MODE=BACK (PREDICTED=PRE_2 SUFFIX='_backtransformed').
EXECUTE.
```

- The TMS IMPORT command reads the transformations in *car\_sales\_transformations.xml* and applies the backtransformation to *PRE\_2*.
- The new field containing the backtransformed values is named *PRE\_2\_backtransformed*.
- The EXECUTE command causes the transformations to be processed. When using this as part of a longer stream of syntax, you may be able to remove the EXECUTE command to save some processing time.

## INFILE Subcommand

The INFILE subcommand specifies the location of the XML file containing the transformations.

### TRANSFORMATIONS Keyword

The TRANSFORMATIONS keyword specifies the location of the XML file containing the transformations. Specify a filename in quotation marks. The filename must be specified in full.

### MODE Keyword

The MODE keyword specifies the direction of the transformations.

#### FORWARD(ROLES=UPDATE|LEAVE)

Convert the “forward” transformations from XML to command syntax. These are the transformations from the original fields into fields that are used in modeling. ROLES determines whether to update field roles based upon transformations in the file. UPDATE will set the role of any “obsolete” field to None. LEAVE will leave the roles of existing fields as they currently are. [For more information, see the topic VARIABLE ROLE on p. 2133.](#)

FORWARD is the default mode.

#### BACK(PREDICTED=fieldname SUFFIX=suffixname)

Convert the backtransformation of target values into command syntax and apply to the field specified on the PREDICTED keyword. The backtransformed values are saved to a field with the name on the PREDICTED keyword and suffix on the SUFFIX keyword. *\_backtransformed* is the default suffix.

### Example

```
TMS IMPORT
  /INFILE TRANSFORMATIONS='file.xml'.
REGRESSION VARIABLES=target_transformed input_transformed
  /DEPENDENT=target_transformed
  /METHOD=ENTER input_transformed
  /SAVE=PRED(predicted) SDBETA(BETA) ICIN.
TMS IMPORT
  /INFILE TRANSFORMATIONS='file.xml' MODE=BACK(PREDICTED=predicted).
```

- The first `TMS IMPORT` reads the file *file.xml*, converts the transformations from XML to command syntax, and applies them to the active dataset. The fields *target\_transformed* and *input\_transformed* result.
- The `REGRESSION` command builds a model for *target\_transformed* using *input\_transformed*, and adds *predicted* to the active dataset, containing the predicted values.
- The second `TMS IMPORT` reads the file *file.xml*, converts the backtransformations from XML to command syntax, and applies them to the active dataset. The field *predicted\_backtransformed* results, and contains predicted values in the original scale of the target field.

## SAVE Subcommand

The `SAVE` subcommand adds the prepared fields to the active dataset.

### TRANSFORMED

*Transformed data.* The transformed fields are added to the active dataset. The values are not computed immediately; instead the transformations are stored, pending execution with the next command that reads the dataset.

## OUTFILE Subcommand

The `OUTFILE` subcommand specifies the location of the file to which command syntax should be saved.

### SYNTAX=filename

*Syntax file.* Specify a filename in quotation marks. The filename must be specified in full; no extension is supplied. It is recommended to include the extension *.sps* in the file specification.

# TMS MERGE

```
TMS MERGE
  /DESTINATION OUTFILE='file specification'
  /TRANSFORMATIONS INFILE='file specification'
  /MODEL INFILE='file specification'
  [/PRINT RESULTS={SUMMARY**}]
  {NONE}
```

\*\* Default if the subcommand or keyword is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 15.0

- Command introduced.

## **Example**

```
TMS MERGE
  /DESTINATION OUTFILE=' /mydir/mymergedmodel.xml '
  /TRANSFORMATIONS INFILE=' /mydir/mytransformations.xml '
  /MODEL INFILE=' /mydir/mymodel.xml '.
```

## **Overview**

The TMS MERGE command is used to merge a PMML file containing transformations with a PMML model file containing variables whose derivations are expressed by those transformations. The merged PMML file can be used to score data with scoring functions. (For more information on scoring functions, see [Scoring expressions](#).) When you score data with a model file containing transformations, the transformations will be applied to the data before scoring. The transformations are carried out as part of the internal scoring process and have no effect on the active dataset. Transformations are exported as PMML using the TMS BEGIN and TMS END commands. [For more information, see the topic TMS BEGIN on p. 1965.](#)

## **Basic Specification**

The basic specification is the command name followed by DESTINATION, TRANSFORMATIONS, and MODEL subcommands that contain file specifications.

## **Syntax Rules**

- Each subcommand can be specified only once.
- Subcommands may be used in any order.
- An error occurs if a keyword is specified more than once within a subcommand.

- Equal signs (=) shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.

### **Operations**

- TMS MERGE has no effect on the active dataset. It relies completely on the information contained in the transformations and model files.
- TMS MERGE verifies that the paths specified for the input files and destination file are valid. The destination file is unavailable to other commands until the TMS MERGE command is ended.
- An error occurs if properties of derived variables in the transformations file do not match the properties for those same variables in the model file—for example, conflicting definitions of missing values.

### **Limitations**

- TMS MERGE cannot be used to merge transformations with an ensemble model.
- TMS MERGE cannot be used to merge transformations with a model that contains splits (separate model information for each split group).

## **TRANSFORMATIONS, MODEL, and DESTINATION Subcommands**

The TRANSFORMATIONS, MODEL, and DESTINATION subcommands specify the file containing the transformations (created from a TMS BEGIN–TMS END block), the file containing the model specifications, and the destination for the merged file. All three of these subcommands are required.

- File specifications should be enclosed in quotation marks.
- The filename must be specified in full. No extension is supplied.
- It is recommended to include the extension *.xml* in the specification for the destination file.

## **PRINT Subcommand**

The PRINT subcommand is optional and specifies whether or not to print tables summarizing the variables in the merged XML file.

<b>SUMMARY</b>	<i>Print summary tables.</i> This is the default.
<b>NONE</b>	<i>Do not print summary tables.</i>

# TREE

TREE is available in the Decision Trees option.

*Note:* Square brackets used in the TREE syntax chart are required parts of the syntax and are not used to indicate optional elements. Equals signs (=) used in the syntax chart are required elements. All subcommands are optional.

```
TREE dependent_variable [level]
  BY variable [level] variable [level]...
  FORCE = variable [level]

  /TREE DISPLAY ={TOPDOWN** } NODES = {STATISTICS**}
    {LEFTTORIGHT}      {CHART      }
    {RIGHTTOLEFT}     {BOTH       }
    {NONE              }
  BRANCHSTATISTICS ={YES**} NODEDEFS = {YES**}
    {NO }              {NO }
  SCALE={AUTO**      }
    {percent value}

  /DEPCATEGORIES USEVALUES=[VALID** value, value...MISSING]
    TARGET=[value value...]

  /PRINT MODELSUMMARY** CLASSIFICATION** RISK** CPS** IMPORTANCE SURROGATES
    TREETABLE CATEGORYSPCS NONE

  /GAIN SUMMARYTABLE = {YES**} CATEGORYTABLE = {YES**}
    {NO }              {NO }
  TYPE = [NODE** PTile] SORT = {DESCENDING**}
    {ASCENDING }
  INCREMENT = {10** } CUMULATIVE = {YES**}
    {value}          {NO }

  /PLOT GAIN RESPONSE INDEX MEAN PROFIT ROI IMPORTANCE INCREMENT = {10 }
    {value}

  /RULES NODES = {TERMINAL** } SYNTAX = {INTERNAL** }
    {ALL }          {SQL }
    {TOPN(value) }  {GENERIC }
    {TOPPCT(value) }
    {MININDEX(value)}
  TYPE = {SCORING**} SURROGATES = {EXCLUDE**} LABELS = {YES**}
    {SELECTION}      {INCLUDE }      {NO }
  OUTFILE = 'filespec'

  /SAVE NODEID(varname) PREDVAL(varname) PREDPROB(rootname) ASSIGNMENT(varname)

  /METHOD TYPE = {CHAID**      }
    {EXHAUSTIVECHAID}
    {CRT }
    {QUEST }
  MAXSURROGATES = {AUTO**} PRUNE = {NONE** }
    {value }      {SE({1 })}
    {value}

  /GROWTHLIMIT MAXDEPTH = {AUTO**} MINPARENTSIZE = {100**}
    {value }      {value}

  MINCHILDSIZE = {50** }
    {value}

  /VALIDATION TYPE = {NONE**      } OUTPUT = {BOTHSAMPLES}
    {SPLITSAMPLE({50 })}      {TESTSAMPLE }
    {percent }
    {varname}
    {CROSSVALIDATION({10 })}
```



```

                                {value}

/CHAID  ALPHASPLIT = {.05**}  ALPHAMERGE = {.05**}
        {value}              {value}
        SPLITMERGED = {NO**}  CHISQUARE = {PEARSON**}
        {YES }           {LR }
        CONVERGE = {.001**}  MAXITERATIONS = {100**}
        {value }          {value}
        ADJUST = {BONFERRONI**}
        {NONE }
        INTERVALS = {10** }
        {value }
        {varlist(value) varlist(value) ...}

/CRT IMPURITY = {GINI** } MINIMPROVEMENT = {.0001**}
               {TWOING } {value }
               {ORDEREDTWOING}

/QUEST ALPHASPLIT = {.05**}
        {value}

/COSTS {EQUAL** }
       {CUSTOM = actcat, predcat [value] actcat, predcat [value] ...}

/PRIORS {FROMDATA** } ADJUST = {NO**}
        {EQUAL } {YES }
        {CUSTOM = cat [value] cat [value] ...}

/SCORES {EQUALINCREMENTS** }
        {CUSTOM = cat [value] cat [value] ...}

/PROFITS CUSTOM = cat [revenue, expense] cat [revenue, expense] ...

/INFLUENCE varname

/OUTFILE TRAININGMODEL = 'filespec' TESTMODEL = 'filespec'

/MISSING NOMINALMISSING = {MISSING**}
                       {VALID }

/TARGETRESPONSE RATE = {NONE }
                       {value }

```

\*\* Default if subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- Command introduced.

Release 18.0

- TARGETRESPONSE subcommand introduced.

### **Example**

TREE risk BY income age creditscore employment.

## Overview

The `TREE` procedure creates a tree-based model. It classifies cases into groups or predicts values of a dependent variable based on values of predictor variables. The procedure provides validation tools for exploratory and confirmatory classification analysis.

### Options

**Model.** You can specify the dependent (target) variable and one or more independent (predictor) variables. Optionally you can force one independent variable into the model as the first variable.

**Growing Method.** Four growing algorithms are available: CHAID (the default), Exhaustive CHAID, CRT, and QUEST. Each performs a type of recursive splitting. First, all predictors are examined to find the one that gives the best classification or prediction by splitting the sample into subgroups (nodes). The process is applied recursively, dividing the subgroups into smaller and smaller groups. It stops when one or more stopping criteria are met.

The four growing methods have different performance characteristics and features:

- CHAID chooses predictors that have the strongest interaction with the dependent variable. Predictor categories are merged if they are not significantly different with respect to the dependent variable (Kass, 1980).
- Exhaustive CHAID is a modification of CHAID that examines all possible splits for each predictor (Biggs et al., 1991).
- CRT is a family of methods that maximizes within-node homogeneity (Breiman et al., 1984).
- QUEST trees are computed rapidly, but the method is available only if the dependent variable is nominal. (Loh and Shih, 1997).

**Stopping Criteria.** You can set parameters that limit the size of the tree and control the minimum number of cases in each node.

**Validation.** You can assess how well your tree structure generalizes to a larger sample. Split-sample partitioning and cross-validation are supported. Partitioning divides your data into a training sample, from which the tree is grown, and a testing sample, on which the tree is tested. Cross-validation involves dividing the sample into a number of smaller samples. Trees are generated excluding the data from each subsample in turn. For each tree, misclassification risk is estimated using data for the subsample that was excluded in generating it. A cross-validated risk estimate is calculated as the average risk across trees.

**Output.** Default output includes a tree diagram and risk statistics. Classification accuracy is reported if the dependent variable is categorical. Optionally, you can obtain charts of gain- and profit-related measures as well as classification rules that can be used to select or score new cases. You can also save the model's predictions to the active dataset, including assigned segment (node), predicted class/value, and predicted probability.

### Basic Specification

- The basic specification is a dependent variable and one or more independent variables.

### **Operations**

- The tree is grown until one or more stopping criteria are met. The default growing method is CHAID.
- The type of model depends on the measurement level of the dependent variable. If the dependent variable is scale (continuous), a prediction model is computed. If it is categorical (nominal or ordinal), a classification model is generated.
- Measurement level determines allowable combinations of predictor values within a node. For ordinal and scale predictors, only adjacent categories/values may occur in a node. There are no restrictions on grouping of nominal categories.
- TREE honors the SET SEED value if split-sample model validation is requested.
- SPLIT FILE is ignored by the TREE procedure.
- If a WEIGHT variable is defined the weights are treated as replication weights. Fractional weights are rounded.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### **Syntax Rules**

- The minimum specification is a dependent variable, the keyword BY and one or more independent variables.
- All subcommands are optional.
- Only a single instance of each subcommand is allowed.
- A keyword may be specified only once within a subcommand.
- Equals signs (=) shown in the syntax chart are required.
- Subcommand names and keywords must be spelled in full.
- Subcommands may be used in any order.

### **Limitations**

- SPLIT FILE is ignored by the TREE procedure.
- CHAID and Exhaustive CHAID: A categorical dependent variable may not have more than 126 categories. If the dependent variable categorical, then the limit for a categorical predictor is also 126 categories.
- CRT: A nominal predictor may not have more than 25 categories.
- QUEST: If a predictor is nominal, then the limit for the dependent variable (which must be nominal) is 127 categories. A nominal predictor may not have more than 25 categories.

### **Examples**

```
TREE risk BY income age creditscore employment.
```

- A tree model is computed that estimates credit risk using an individual's income, age, credit score, and employment category as predictor variables.
- The default method, CHAID, is used to grow the tree.
- Since measurement level is not specified, it is obtained from the data dictionary for each model variable. If no measurement levels have been defined, numeric variables are treated as scale and string variables are treated as nominal.

```
TREE risk [o] BY income [o] age [s] creditscore [s] employment [n]
/METHOD TYPE=CRT
/VALIDATION TYPE=SPLITSAMPLE
/SAVE NODEID PREDVAL.
```

- A tree model is computed that estimates credit risk using an individual's income, age, credit score, and employment category as predictor variables.
- Age and credit score will be treated as scale variables, risk and income as ordinal, and employment category as nominal.
- The CRT method, which performs binary splits, is used to grow the tree.
- Split-sample validation is requested. By default, 50% of cases are assigned to the training sample. Remaining cases are used to validate the tree.
- Two variables are saved to the active dataset: node (segment) identifier and predicted value.

## Model Variables

The command name `TREE` is followed by the dependent variable, the keyword `BY`, and one or more independent variables.

### Dependent Variable

The dependent (target) variable must be the first specification on the `TREE` command.

- Only one dependent variable can be specified.
- The dependent variable cannot be the `WEIGHT` variable.

### Independent Variables

You can specify one or more independent (predictor) variables.

- `ALL` and `TO` can be used in the independent variable list. `ALL` represents all variables in the active dataset. `TO` refers to a range of variables in the active dataset.
- In general, if the independent variable list refers to the `WEIGHT` variable or a variable that is used elsewhere in the tree specification (e.g., dependent variable or cell weight variable), the variable is not used as an independent variable. The exception is that scale variables used both in the independent variable list and the `INTERVAL` subcommand are not dropped from the independent variable list.
- Repeated instances of the same independent variable are filtered out of the list. For example, `a b a c a` is equivalent to `a b c`. Assuming that the active dataset contains variables *a*, *b*, and *c*, `a ALL` is also equivalent to `a b c`.

## Measurement Level

Optionally, a measurement level can be specified in square brackets following any model variable. The measurement level can be defined as scale ([S]), ordinal ([O]), or nominal ([N]). Categorical ([C]) is accepted as a synonym for nominal.

- If a measurement level is specified, it temporarily overrides the setting recorded in the data dictionary. If a measurement level is not specified for a variable, the dictionary setting is used.
- If a measurement level is not specified and no setting is recorded in the data dictionary, a numeric variable is treated as scale and a string variable is treated as nominal.
- If a string variable is defined as scale, a warning is issued and the variable is treated as nominal.
- A measurement level applies only to the variable that immediately precedes it in the variable list. For example, `age income [S]` and `age TO income [S]` assign the scale level of measurement to *income* only.
- The keyword `ALL` followed by a measurement level specification will apply that measurement level to all independent variables.
- If a variable appears more than once in the independent variable list and the measurement level varies across instances, the measurement level specified for the last instance of the variable is used. For example, if `ALL [S] z [N]` is specified, *z* is treated as a nominal predictor.

## FORCE Keyword

- By default, the procedure evaluates each predictor variable for inclusion in the model. The optional `FORCE` keyword forces an independent variable into the model as the first split variable (the first variable used to split the root node into child nodes).
- The `FORCE` keyword must be followed by an equals sign (=) and a single variable.
- A measurement level may be specified in square brackets following the variable name. If the variable is also included in the independent variable list, the measurement level specified for the `FORCE` variable is used. If the variable has a specified measurement level in the independent list but no explicit measurement level on the `FORCE` specification, the defined measurement level for the variable is used.
- The variable cannot be the dependent variable or the `WEIGHT` variable.

### Example

```
TREE risk [o] BY income [o] age [s] creditscore [s] FORCE = gender [n].
```

- The independent variables *income*, *age*, and *creditscore* may or may not be included in the model, depending on how strongly they interact with the dependent variable *risk*.
- The independent variable *gender* will be included in the model regardless of the level of interaction with the dependent variable, and the root node will first be split into nodes for male and female before any other independent variables are considered.

## ***DEPCATEGORIES Subcommand***

For categorical dependent variables, the `DEPCATEGORIES` subcommand controls the categories of the dependent variable included in the model and/or specifies target categories.

- By default, all valid categories are used in the analysis and user-missing values are excluded.
- There is no default target category.
- `DEPCATEGORIES` applies only to categorical dependent variables. If the dependent variable is scale, the subcommand is ignored and a warning is issued.

### ***Example***

```
TREE risk [n] BY income [o] age [s] creditscore [s] employment [n]
  /DEPCATEGORIES USEVALUES=[VALID 99] TARGET=[1].
```

### ***USEVALUES keyword***

`USEVALUES` controls the categories of the dependent variable included in the model.

- The keyword is followed by an equals sign (=) and a list of values enclosed in square brackets, as in: `USEVALUES=[1 2 3]`.
- At least two values must be specified.
- Any cases with dependent variable values not included in the list are excluded from the analysis.
- Values can be string or numeric but must be consistent with the data type of the dependent variable. String and date values must be quoted. Date values must be consistent with the variable's print format.
- If the dependent variable is nominal, the list can include user-missing values. If the dependent variable is ordinal, any user-missing values in the list are excluded and a warning is issued.
- The keywords `VALID` and `MISSING` can be used to specify all valid values and all user-missing values respectively. For example, `USEVALUES=[VALID MISSING]` will include all cases with valid or user-missing values for the dependent variable.
- A warning is issued if a specified category does not exist in the data or in the training sample if split-sample validation is in effect. [For more information, see the topic VALIDATION Subcommand on p. 1994.](#)

### ***TARGET Keyword***

The `TARGET` keyword specifies categories of the dependent variable that are of primary interest in the analysis. For example, if you are trying to predict mortality, "death" would be defined as the target category. If you are trying to identify individuals with high or medium credit risk, you would define both "high credit risk" and "medium credit risk" as target categories.

- The keyword is followed by an equals sign (=) and a list of values enclosed in square brackets, as in: `TARGET=[1]`.
- There is no default target category. If not specified, some classification rule options and gains-related output are not available.

- Values can be string or numeric but must be consistent with the data type of the dependent variable. String and date values must be quoted. Date values must be consistent with the variable's print format.
- If USEVALUES is also specified, the target categories must also be included (either implicitly or explicitly) in the USEVALUES list.
- If a target category is user-missing and the dependent variable is not nominal, the target category is ignored and a warning is issued.

## TREE Subcommand

The TREE subcommand allows you to specify the appearance of the tree or suppress it altogether. You can control its orientation, node contents, and the labeling of splits and nodes. Each keyword is followed by an equals sign (=) and the value for that keyword.

### Example

```
TREE risk BY income age creditscore
  /TREE DISPLAY=LEFTTORIGHT NODES=BOTH BRANCHSTATISTICS=NO.
```

### DISPLAY Keyword

The DISPLAY keyword controls the orientation the tree and can also be used to suppress the display of the tree. The following alternatives are available:

<b>TOPDOWN</b>	<i>Tree grows top down.</i> The root node appears at the top of the diagram. The tree grows downward. This is the default.
<b>LEFTTORIGHT</b>	<i>Tree grows from left to right.</i>
<b>RIGHTTOLEFT</b>	<i>Tree grows from right to left.</i>
<b>NONE</b>	<i>Tree diagram is not generated.</i>

### NODES Keyword

The NODES keyword controls the contents of tree nodes.

<b>STATISTICS</b>	<i>Nodes display summary statistics.</i> This is the default. For a categorical dependent variable, percentages and counts are shown for each category as well as the total count and percentage. (Total percentage is the percentage of cases in the sample assigned to the node. If you specify prior probabilities, the total percentage is adjusted to account for those priors.) For a scale dependent variable, nodes display mean, standard deviation, number of cases, and predicted value.
<b>CHART</b>	<i>Nodes display a summary chart.</i> For a categorical dependent variable, nodes display a bar chart of percentages for each category. For a scale dependent variable a histogram is shown that summarizes the distribution of the dependent variable.
<b>BOTH</b>	<i>Statistics and chart are shown.</i>

**BRANCHSTATISTICS Keyword**

The `BRANCHSTATISTICS` keyword controls whether branch (split) statistics are shown in the tree.

- For CHAID and Exhaustive CHAID, branch statistics include  $F$  value (for scale dependent variables) or chi-square value (for categorical dependent variables), as well as the significance value and degrees of freedom.
- For CRT, the improvement value is shown.
- For QUEST,  $F$ , the significance value, and degrees of freedom are shown for scale and ordinal independent variables. For nominal independent variables, chi-square, the significance value, and degrees of freedom are shown.

**YES**                    *Branch statistics are displayed. This is the default.*

**NO**                    *Branch statistics are not displayed.*

**NODEDEFS Keyword**

The `NODEDEFS` keyword controls whether descriptive labels are shown for each node. Node labels show the independent variable (predictor) values assigned to each node.

**YES**                    *Node labels are displayed. This is the default.*

**NO**                    *Node labels are not displayed.*

**SCALE Keyword**

The `SCALE` keyword controls the initial scale of the tree diagram.

**AUTO**                    *Adjust the scale according to the number of nodes and levels of tree growth. The scale is reduced automatically for large trees. This is the default.*

**value**                    *User-specified percentage value. The value can be between 5 and 200, inclusive.*

**PRINT Subcommand**

The `PRINT` subcommand controls the display of optional tabular output. If the `PRINT` subcommand is specified, only the output explicitly requested will be produced. The subcommand name is followed by an equals sign (=) and one or more of the following options:

**MODELSUMMARY**    *Model summary.* The summary includes the method used, variables included in the model, and variables specified but not included in the model. Shown by default if subcommand omitted.

**IMPORTANCE**        *Predictor importance.* Ranks each independent (predictor) variable according to its importance to the model. `IMPORTANCE` can be specified only for the CRT method; otherwise it is ignored, and a warning is issued.

**SURROGATES**        *Surrogate predictors.* Lists surrogate predictors for each split in the tree. Each node is split based on an independent (predictor) variable; for cases in which the value for that predictor is missing, other predictors having high associations with the original predictor are used for classification. These alternative predictors are called surrogates. `SURROGATES` can be specified only for the CRT and QUEST methods. If surrogates are not computed, this keyword is ignored, and a warning is issued.



<b>CLASSIFICATION</b>	<i>Classification table.</i> Shows the number of cases classified correctly and incorrectly for each category of the dependent variable. A warning is issued if CLASSIFICATION is specified for a scale dependent variable. Shown by default for categorical dependent variables if subcommand omitted.
<b>RISK</b>	<i>Risk estimate and its standard error.</i> A measure of the tree's predictive accuracy. For categorical dependent variables, the risk estimate is the proportion of cases correctly classified after adjustment for prior probabilities and misclassification costs. For scale dependent variables, the risk estimate is within-node variance. Shown by default if subcommand omitted.
<b>CATEGORYSPEC</b>	<i>Cost, prior probability, score, and profit values used in the analysis.</i> This keyword is ignored if the dependent variable is a scale variable.
<b>TREETABLE</b>	<i>Tree model in tabular format.</i> Summary information for each node in the tree, including parent node number, branch statistics, independent variable value(s) for the node, mean and standard deviation for scale dependent variables or counts and percents for categorical dependent variables.
<b>CPS</b>	<i>Case processing summary.</i> Shown by default if subcommand omitted.
<b>NONE</b>	<i>None of the output controlled by the PRINT subcommand is generated.</i> An error occurs if any other PRINT keyword is specified with NONE.

## **GAIN Subcommand**

The GAIN subcommand controls the display of tables that summarize gain and index values and other measures that indicate the relative performance of nodes in the tree. You can obtain tables that summarize results by node, percentile group, or both. Each keyword is followed by an equals sign (=) and the value for that keyword.

### **Example**

```
TREE risk BY income age creditscore
  /GAIN SUMMARYTABLE=YES TYPE=[NODE, PTILE].
```

### **SUMMARYTABLE Keyword**

The SUMMARYTABLE keyword controls whether a summary table of index values is shown.

- For scale dependent variables, the table includes node number, number of cases, and mean value of the dependent variable.
- For categorical dependent variables, the table includes node number, number of cases, average profit, and ROI (return on investment) values.
- For categorical dependent variables, the summary table is not displayed unless profits have been specified on the PROFITS subcommand.

**YES**     *Display summary table.* This is the default for scale dependent variables if the subcommand is omitted or SUMMARYTABLE=NO is not included on the subcommand.

**NO**     *The summary table is not displayed.*

### **CATEGORYTABLE Keyword**

The CATEGORYTABLE keyword displays gain and index values for target categories of the dependent variable.

- The dependent variable must be categorical.
- Target categories must be specified on the DEPCATEGORIES subcommand.
- A separate table is generated for each target category.
- If the dependent variable is scale or no target categories are specified, the keyword is ignored and a warning is issued.

**YES**     *Display a table of gain statistics for each target category of the dependent variable. The table includes percentage gain, response percent, and index percent (lift) by node or percentile group. This is the default.*

**NO**     *No category table is shown.*

### **TYPE Keyword**

The **TYPE** keyword specifies the type of gain table(s) that are produced. The keyword is followed by an equals sign (=) and one or both of the following types enclosed in square brackets:

**NODE**     *Terminal nodes. Rows of the table represent terminal nodes. This is the default.*

**PTILE**     *Percentile. Each row of the table summarizes a percentile group.*

### **SORT Keyword**

The **SORT** keyword specifies how rows are sorted in all gain tables.

**DESCENDING**     *Rows are sorted in descending order of index value. This is the default.*

**ASCENDING**     *Rows are sorted in ascending order of index value.*

### **INCREMENT Keyword**

The **INCREMENT** keyword controls the width of percentile groupings for a percentile gain table (**TYPE**= [PTILE]).

- The default increment is 10, which gives decile intervals.
- You can specify one of the following percentage increments: 1, 2, 5, 10, 20, or 25. If a different value is specified, a warning is issued and the default increment is used.
- **INCREMENT** is ignored for node-by-node gain tables.

### **CUMULATIVE Keyword**

The **CUMULATIVE** keyword specifies whether node-by-node tables (**TYPE**= [NODE]) include cumulative statistics. It is ignored for percentile gain tables, which are always cumulative.

**YES**     *Cumulative statistics are displayed. This is the default.*

**NO**     *Cumulative statistics are not displayed.*

## **PLOT Subcommand**

The PLOT subcommand generates charts that may help you evaluate your tree model. The subcommand name is followed by an equals sign (=) and any or all of the following chart types:

<b>GAIN</b>	<i>Line chart of percentage gain.</i> The chart shows percentage gain for each target category of the dependent variable. GAIN is available only for categorical dependent variables with target categories defined on the DEPCATEGORIES subcommand.
<b>RESPONSE</b>	<i>Line chart of response.</i> The chart shows percentage of responses for each target category of the dependent variable. RESPONSE is available only for categorical dependent variables with target categories defined on the DEPCATEGORIES subcommand.
<b>INDEX</b>	<i>Line chart of index percent.</i> The chart shows index percent (lift). INDEX is available only for categorical dependent variables with target categories defined on the DEPCATEGORIES subcommand.
<b>MEAN</b>	<i>Line chart of mean value for scale dependent variables.</i> If the dependent variable is categorical, this keyword is ignored and a warning is issued.
<b>PROFIT</b>	<i>Line chart of average profit for categorical dependent variables.</i> If profits are not defined or the dependent variable is scale, this keyword is ignored and a warning is issued. <a href="#">For more information, see the topic PROFITS Subcommand on p. 2002.</a>
<b>ROI</b>	<i>Line chart of ROI (return on investment).</i> ROI is computed as the ratio of profits to expenses. If profits are not defined or the dependent variable is scale, this keyword is ignored and a warning is issued.
<b>IMPORTANCE</b>	<i>Bar chart of model importance by predictor.</i> IMPORTANCE is available for CRT models only; otherwise the keyword is ignored and a warning is issued.

### **Example**

```
TREE risk [o] BY income age creditscore
/METHOD TYPE=CRT
/DEPCATEGORIES TARGET=[1]
/PLOT=IMPORTANCE GAIN INDEX INCREMENT=5.
```

### **INCREMENT Keyword**

The INCREMENT keyword controls the width of percentile group charts.

- The keyword is followed by an equals sign (=) and one of the following values: 1, 2, 5, 10, 20, or 25. If a different value is specified a warning is issued and the default increment is used.
- INCREMENT is ignored for the predictor importance chart.

## **RULES Subcommand**

The RULES subcommand generates syntax that can be used to select or classify cases based on values of independent (predictor) variables.

- You can generate rules for all nodes, all terminal nodes, the top *n* terminal nodes, terminal nodes that correspond to the top *n* percent of cases, or nodes with index values that meet or exceed a cutoff value.
- Rules are available in three different forms: internal command syntax format, SQL, and generic (plain English pseudocode).

- You can specify an external destination file for the rules.
- Each keyword is followed by an equals sign (=) and the value for that keyword.

### **Example**

```
TREE risk [o] BY income age creditscore
/RULES NODES=TERMINAL SYNTAX=INTERNAL TYPE=SCORING
OUTFILE='/jobfiles/treescores.sps'.
```

### **NODES Keyword**

The NODES keyword specifies the scope of generated rules. Specify one of the following alternatives:

- TERMINAL**     *Generates rules for each terminal node.* This is the default.  
**ALL**             *Generates rules for all nodes.* Rules are shown for all parent and terminal nodes.

For categorical dependent variables with defined target categories, the following additional alternatives are available:

- TOPN(value)**             *Generates rules for the top n terminal nodes based on index values.* The number must be a positive integer, enclosed in parentheses. If the number exceeds the number of nodes in the tree, a warning is issued and rules are generated for all terminal nodes.
- TOPPCT(value)**           *Generates rules for terminal nodes for the top n percent of cases based on index values.* The percent value must be a positive number greater than zero and less than 100, enclosed in parentheses.
- MININDEX(value)**         *Generates rules for all terminal nodes with an index value greater than or equal to the specified value.* The value be a positive number, enclosed in parentheses.

### **SYNTAX Keyword**

The SYNTAX keyword specifies the syntax of generated rules. It determines the form of the selection rules in both output displayed in the Viewer and selection rules saved to an external file.

- INTERNAL**     *Command syntax language.* Rules are expressed as a set of commands that define a filter condition that can be used to select subsets of cases (this is the default) or as COMPUTE statements that can be used to score cases (with TYPE=SCORING).
- SQL**             *SQL.* Standard SQL rules are generated to select/extract records from a database or assign values to those records. The generated SQL rules do not include any table names or other data source information.
- GENERIC**       *Plain English pseudocode.* Rules are expressed as a set of logical “if..then” statements that describe the model’s classifications or predictions for each node.

### **TYPE Keyword**

The `TYPE` keyword specifies the type of SQL or internal command syntax rules to generate. It is ignored if generic rules are requested.

- SCORING**      *Scoring of cases.* The rules can be used to assign the model's predictions to cases that meet node membership criteria. A separate rule is generated for each node within the scope specified on the `NODES` keyword. This is the default.
- SELECTION**    *Selection of cases.* The rules can be used to select cases that meet node membership criteria. For internal command syntax and SQL rules, a single rule is generated to select all cases within the scope specified on the `NODES` keyword.
- Note:* For internal command syntax or SQL rules with `NODES=TERMINAL` and `NODES=ALL`, `TYPE=SELECTION` will produce a rule that effectively selects every case included in the analysis.

### **SURROGATES Keyword**

For `CRT` and `QUEST`, the `SURROGATES` keyword controls whether rules use surrogate predictors to classify cases that have missing predictor values. The keyword is ignored if the method is not `CRT` or `QUEST` or if generic rules are requested.

Rules that include surrogates can be quite complex. In general, if you just want to derive conceptual information about your tree, exclude surrogates. If some cases have incomplete predictor data and you want rules that mimic your tree, include surrogates.

- INCLUDE**      *Include surrogates.* Surrogate predictors are used in generated rules. This is the default.
- EXCLUDE**      *Exclude surrogates.* Rules exclude surrogate predictors.

### **LABELS Keyword**

The `LABELS` keyword specifies whether value and variable labels are used in generic decision rules.

- By default, any defined value and variable labels are used. When labels aren't available, values and variable names are used.
- `LABELS` is ignored for SQL and internal command syntax rules.

- YES**          *Any defined value and variable labels are used in generic rules.* This is the default.
- NO**          *Values and variable names are used instead of labels.*

### **OUTFILE Keyword**

`OUTFILE` writes the rules to an external text file.

- The keyword is followed by an equals sign (=) and a file specification enclosed in quotes.
- If the file specification includes a path, an error will result if the specified directory/folder location doesn't exist.
- For command syntax, the file can be used as a command syntax file in both interactive and batch modes.

- For SQL syntax, the generated SQL does not include any table names or other data source information.
- OUTFILE is ignored if NODES=NONE, and a warning is issued.

## SAVE Subcommand

The SAVE subcommand writes optional model variables to the active dataset.

- Specify one or more model variables.
- Optionally, you can specify new variable names in parentheses after each model variable name. The names must be unique, valid variable names.
- If new names are not specified, or if specified names are not valid variable names, unique variable names are automatically generated using the model variable name with a suffix.

<b>NODEID</b>	<i>Node identifier.</i> The terminal node to which a case is assigned. The value is the tree node number.
<b>PREDVAL</b>	<i>Predicted value.</i> The class or value predicted by the model.
<b>PREDPROB</b>	<i>Predicted probability.</i> The probability associated with the model's prediction. One variable is saved for each category of the dependent variable. If you specify an optional variable name, it is used as a root for generated variable names. PREDPROB is ignored if the dependent variable is a scale variable. Probabilities are saved only for the first 126 categories of the dependent variable when the CRT or QUEST algorithms are used.
<b>ASSIGNMENT</b>	<i>Sample assignment.</i> The variable indicates whether a case was used in the training or testing sample. Cases in the training sample have a value of 1, and cases in the testing sample have a value of 0. ASSIGNMENT is ignored if split-sample validation is not specified.

### Example

```
TREE risk [o] BY income age creditscore
  /SAVE NODEID(node_number) PREDVAL(predicted_value).
```

## METHOD Subcommand

The METHOD subcommand specifies the growing method and optional parameters. Each keyword is followed by an equals sign (=) and the value for that keyword.

### Example

```
TREE risk [o] BY income age creditscore
  /METHOD TYPE=CRT MAXSURROGATES=2 PRUNE=SE(0).
```

**TYPE Keyword**

TYPE specifies the growing method. For CRT and QUEST, splits are always binary. CHAID and Exhaustive CHAID allow multiway splits.

<b>CHAID</b>	<i>Chi-squared Automatic Interaction Detection.</i> At each step, CHAID chooses the independent (predictor) variable that has the strongest interaction with the dependent variable. Categories of each predictor are merged if they are not significantly different with respect to the dependent variable. This is the default method.
<b>EXHAUSTIVECHAID</b>	<i>Exhaustive CHAID.</i> A modification of CHAID that examines all possible ways of merging predictor categories.
<b>CRT</b>	<i>Classification and Regression Trees.</i> CRT splits the data into segments that are as homogeneous as possible with respect to the dependent variable.
<b>QUEST</b>	<i>Quick, Unbiased, Efficient Statistical Tree.</i> A method that is fast and avoids other methods' bias in favor of predictors with many categories. QUEST can be specified only if the dependent variable is nominal. An error occurs if the dependent variable is ordinal or scale.

**MAXSURROGATES Keyword**

CRT and QUEST can use surrogates for independent (predictor) variables. For cases in which the value for that predictor is missing, other predictors having high associations with the original predictor are used for classification. These alternative predictors are called surrogates. The MAXSURROGATES keyword specifies the maximum number of surrogate predictors to compute.

If the growing method is CHAID or EXHAUSTIVECHAID, this keyword is ignored and a warning is issued.

<b>AUTO</b>	<i>The maximum is the number of independent variables minus one.</i> This is the default.
<b>value</b>	<i>User-specified value.</i> The value must be a non-negative integer that is less than the number of independent variables in the model. If you don't want to use surrogates in the model, specify MAXSURROGATES=0. If the value equals or exceeds the number of independent variables, the setting is ignored and a warning is issued.

**PRUNE Keyword**

For CRT and QUEST, the tree can be automatically pruned. Pruning can help avoid creating a tree that overfits the data. If you request pruning, the tree is grown until stopping criteria are met. Then it is trimmed automatically according to the specified criterion.

If the growing method is CHAID or EXHAUSTIVECHAID, PRUNE is ignored and a warning is issued.

<b>NONE</b>	The tree is not pruned. This is the default.
<b>SE(value)</b>	<i>Prune tree using standard error criterion.</i> The procedure prunes down to the smallest subtree with a risk value within a specified number of standard errors of that of the subtree with the minimum risk. You can specify the number of standard errors in parentheses. The default is 1. The value must be nonnegative. To obtain the subtree with the minimum risk, specify 0.

## **GROWTHLIMIT Subcommand**

The GROWTHLIMIT subcommand specifies stopping criteria that limit the size of the tree. Each keyword is followed by an equals sign (=) and the value for that keyword.

### **Example**

```
TREE risk [o] BY income age creditscore
/METHOD TYPE=CRT
/GROWTHLIMIT MAXDEPTH=4 MINCHILDSIZE=10.
```

### **MAXDEPTH Keyword**

MAXDEPTH specifies the maximum number of levels of growth beneath the root node. You can change the maximum depth to adjust the size of the tree.

**AUTO**        *Three levels for CHAID and Exhaustive CHAID, five levels for CRT and QUEST. This is the default.*

**value**        *User-specified value. The value must be a positive integer.*

### **MINPARENTSIZE Keyword**

MINPARENTSIZE specifies the minimum number of cases required to split a node. Nodes with fewer cases are not split. You can use this setting to avoid splitting nodes that have few cases.

- The size value must be a positive integer. The default is 100.
- MINPARENTSIZE must be greater than MINCHILDSIZE.

### **MINCHILDSIZE Keyword**

MINCHILDSIZE specifies the minimum number of cases in any child node. A node will not be split if any of the resulting child nodes would have fewer cases than the specified value.

- The size value must be a positive integer. The default is 50.
- MINCHILDSIZE must be less than MINPARENTSIZE.

### **MINIMPROVEMENT Keyword**

For CRT, you can use MINIMPROVEMENT to specify the minimum decrease in impurity. The CRT growing method attempts to maximize within-node homogeneity. In other words, a terminal node in which all cases have the same value for the dependent variable is a homogeneous, “pure” node. A node is not split if impurity would decrease less than the specified value.

The improvement value must be positive value. The default is 0.0001. As the value increases, the number of nodes in the tree tends to decrease.

## **VALIDATION Subcommand**

The VALIDATION subcommand allows you to assess how well your tree structure generalizes to a larger population.



- Split-sample validation and cross-validation are available.
- By default, validation is not performed.
- If you want to be able to reproduce validated results later, use `SET SEED` before the `TREE` procedure to initial the random number seed.

Each keyword in the subcommand is followed by an equals sign (=) and the value for that keyword.

### Example

```
TREE risk [o] BY income age creditscore
  /VALIDATION TYPE=SPLITSAMPLE(25) OUTPUT=TESTSAMPLE.
```

### TYPE Keyword

<b>NONE</b>	<i>The tree model is not validated.</i> This is the default.
<b>SPLITSAMPLE(percent)</b>	<i>Split-sample validation.</i> The model is generated using a training sample and tested on a hold-out sample. The value or variable specified in parentheses determines the training sample size. Enter a percent value greater than 0 and less than 100, or a numeric variable, the values of which determine how cases are assigned to the training or testing samples: cases with a value of 1 for the variable are assigned to the training sample, and all other cases are assigned to the testing sample. The variable cannot be the dependent variable, weight variable, influence variable or a forced independent variable. <a href="#">For more information, see the topic INFLUENCE Subcommand on p. 2003.</a> <i>Note:</i> Split-sample validation should be used with caution on small data files (data files with a small number of cases). Small training sample sizes may yield poor models since there may not be enough cases in some categories to adequately grow the tree.
<b>CROSSVALIDATION(value)</b>	<i>Crossvalidate the tree model.</i> The sample is divided into a number of subsamples, or <b>folds</b> . Tree models are then generated excluding the data from each subsample in turn. The first tree is based on all of the cases <i>except</i> those in the first sample fold, the second tree is based on all of the cases <i>except</i> those in the second sample fold, and so on. For each tree, misclassification risk is estimated by applying the tree to the subsample excluded in generating it. Specify a positive integer between 2 and 25 in parentheses. The higher the value, the fewer the number of cases excluded for each tree model. Crossvalidation produces a single, final tree model. The cross-validated risk estimate for the final tree is calculated as the average of the risks for all of the trees. <code>CROSSVALIDATION</code> is ignored with a warning if <code>FORCE</code> is also specified.

### OUTPUT Keyword

With split-sample validation, the `OUTPUT` keyword controls the output generated. This setting is ignored if `SPLITSAMPLE` is not specified.

<b>BOTHSAMPLES</b>	<i>Output is produced for training and test samples.</i> This is the default. Choose this option if you want to compare results for each partition.
<b>TESTSAMPLE</b>	<i>Output is produced for the test sample only.</i> Choose this option if you want validated results only.



**CONVERGE Keyword**

For nominal and ordinal dependent variables, the `CONVERGE` keyword specifies the convergence value for estimation of the CHAID model.

- Specify a value greater than zero and less than 1.
- The default value is 0.05.
- If the dependent variable is nominal or scale, this keyword is ignored, and a warning is issued.

**MAXITERATIONS Keyword**

For nominal and ordinal dependent variables, the `MAXITERATIONS` keyword specifies the maximum number of iterations for estimation of the CHAID model.

- Specify a positive integer value.
- The default value is 100.
- If the dependent variable is nominal or scale, this keyword is ignored, and a warning is issued.

**ADJUST Keyword**

The `ADJUST` keyword specifies how to adjust significance values for multiple comparisons.

- BONFERRONI**     *Significance values are adjusted using the Bonferroni method. This is the default.*  
**NONE**             *Significance values are not adjusted.*

**INTERVALS Keyword**

In CHAID analysis, scale independent (predictor) variables are always banded into discrete groups (for example, 0-10, 11-20, 21-30, and so on) prior to analysis. You can use the `INTERVALS` keyword to control the number of discrete intervals for scale predictors.

- By default, each scale predictor is divided into 10 intervals that have approximately equal numbers of cases.
- The `INTERVALS` keyword is ignored if the model contains no scale independent variables.

**value**             *The specified value applies to all scale predictors. For example: `INTERVALS=5`. The value must be a positive integer less than or equal to 64.*

**varlist (value)**     *The specified value applies to the preceding variable list only. Specify a list of variables followed by the number of intervals in parentheses. Multiple lists can be specified. For example: `INTERVALS=age income (10) creditscore (5)`. The value must be a positive integer less than or equal to 64.*

For the `varlist (value)` form:

- If a variable in the list is not a scale variable and/or the variable is not specified as an independent variable, the interval specification is ignored for that variable, and a warning is issued.
- If a variable appears in more than one list, the last specification is used.
- For any scale variables not include in the list(s), the default number of intervals (10) is used.

## ***CRT Subcommand***

The **CRT** subcommand sets parameters for the CRT method. For other tree growing methods, the subcommand is ignored, and a warning is issued.

Each keyword in the subcommand is followed by an equals sign (=) and the value for that keyword.

### ***Example***

```
TREE risk [o] BY income age creditscore  
/METHOD TYPE=CRT  
/CRT IMPURITY=TWOING MINIMPROVEMENT=.001.
```

### ***IMPURITY Keyword***

For categorical dependent variables, the **IMPURITY** keyword specifies the impurity measure used.

<b>GINI</b>	Gini is based on squared probabilities of membership for each category of the dependent variable. It reaches its minimum (zero) when all cases in a node fall into a single category. This is the default.
<b>TWOING</b>	Twoing groups categories of the dependent variable into the two best subclasses and computes the variance of the binary variable indicating the subclass to which each case belongs.
<b>ORDEREDTWOING</b>	A variant of twoing for ordinal dependent variables in which only adjacent classes can be combined. This method is available only for ordinal dependent variables only.

For scale dependent variables the least squared deviation (LSD) measure of impurity is used. It is computed as the within-node variance, adjusted for any frequency weights or influence values.

### ***MINIMPROVEMENT Keyword***

**MINIMPROVEMENT** specifies the minimum decrease in impurity for CRT trees. A node is not split if impurity would decrease less than the specified value.

- The default value is 0.0001.
- Specify a positive value. Large improvement values tend to result in smaller trees.

## ***QUEST Subcommand***

The **ALPHASPLIT** keyword on the **QUEST** subcommand sets the alpha criterion for selection of independent (predictor) variables. Specify a value greater than 0 and less than 1. The default is 0.05.

### ***Example***

```
TREE risk [n] BY income age creditscore  
/METHOD TYPE=QUEST  
/QUEST ALPHASPLIT=.01.
```

## ***COSTS Subcommand***

For categorical dependent variables, the `COSTS` subcommand allows you to specify penalties for misclassifying cases.

- Costs are used in node assignment and risk estimation for all trees.
- They also affect the tree-growing process for all trees except `QUEST` trees and `CRT` trees that use a twoing measure. To include cost information in the tree-growing process for such trees, use adjusted priors. For more information, see the topic `PRIORS` Subcommand on p. 2000.
- The `COST` subcommand is ignored if the dependent variable is scale, and a warning is issued.

The subcommand name is followed by an equals sign (=) and one of the following alternatives:

- EQUAL**      *Misclassification costs are equal across categories of the dependent variable. Cost is set to 1 for all possible misclassifications. This is the default.*
- CUSTOM**     *User-specified costs. Use this to specify costs for one or more combinations of actual and predicted categories. Any combination for which a cost is not specified defaults to 1.*

### ***Custom Costs***

For each combination of dependent variable categories:

- Specify the actual category, predicted category, and misclassification cost, in that order.
- The cost value must be enclosed in square brackets.
- At least one cost specification must be provided, but you don't have to specify all possible combinations. Any combination for which a cost is not specified defaults to 1.
- If multiple cost specifications are provided for a particular combination of predicted and actual categories, the last one is used.
- Costs specified for categories that do not exist in the data are ignored, and a warning is issued.
- Cost values must be non-negative.
- Category values must be consistent with the data type of the dependent variable. String and date values must be quoted. Date values must be consistent with the variable's print format.
- Correct classifications, where the predicted category and the actual category match, are always assigned a cost of zero. A warning is issued if a nonzero cost is specified for a correct classification. For example, `COSTS CUSTOM=1 1 [3]` would be ignored and the cost would be set to 0.

#### ***Example***

```
TREE risk [0] BY age income employment
  /COSTS CUSTOM=3 2 [2]
                3 1 [8].
```

- Assuming that the dependent variable is coded 1=*low*, 2=*medium*, and 3=*high*, the cost of misclassifying a high-risk individual as medium risk is 2.
- The cost of misclassifying a high-risk individual as low risk is 8.

- All other misclassifications, such as classifying a medium-risk individual as high risk, are assigned the default cost, 1.
- Correct classifications such as classifying a high-risk individual as high risk are always assigned a cost of 0.

## **PRIORS Subcommand**

For CRT and QUEST trees with categorical dependent variables, the `PRIORS` subcommand allows you to specify prior probabilities of group membership. **Prior probabilities** are estimates of the overall relative frequency for each target category of the dependent variable prior to knowing anything about the values of the independent (predictor) variables. Using prior probabilities helps correct any tree growth caused by data in the sample that is not representative of the entire population.

- If the growing method is CHAID or Exhaustive CHAID, this subcommand is ignored and a warning is issued.
- If the dependent variable is scale, this subcommand is ignored and a warning is issued.

- FROMDATA** *Obtain priors from the training sample.* Use this setting if the distribution of groups in the training sample is representative of the population distribution. This is the default. If you have not specified a training sample using the `VALIDATION` subcommand, then the distribution of values in the entire data file is used.
- EQUAL** *Equal priors across categories.* Use this if categories of the dependent variable are represented equally in the population. For example, if there are four categories, approximately 25% of the cases are in each category.
- CUSTOM** *User-specified prior probabilities.* For each category of the dependent variable, specify the category followed by a non-negative prior probability value enclosed in square brackets.

### **Custom Prior Probabilities**

- Prior probabilities must be specified for all values of the dependent variable included in the analysis (either all non-missing values found in the data or all values defined on the `DEPCATEGORIES` subcommand).
- The values must be non-negative.
- The specified category values must be consistent with the data type of the dependent variable. String and date values must be quoted. Date values must be consistent with the variable's print format.
- If you specify the same category more than once, the last one is used.
- A warning is issued if you specify a prior probability for a category that does not exist in the data or in the training sample if split-sample validation is in effect. [For more information, see the topic VALIDATION Subcommand on p. 1994.](#)
- Prior probability values are “normalized” to relative proportions before the tree is grown.

### **Example**

```
TREE risk [o] BY age income employment
/METHOD TYPE=CRT
/PRIORS CUSTOM= 1 [30] 2 [75] 3 [45].
```

The prior probabilities of 30, 75, and 45 are normalized to proportions of 0.2, 0.5, and 0.3, respectively.

### **ADJUST Keyword**

The ADJUST keyword specifies whether prior probabilities are adjusted to take into account misclassification costs.

- NO**            *Priors are unadjusted.* This is the default.  
**YES**           *Priors are adjusted to take into account misclassification costs.*

## **SCORES Subcommand**

For CHAID or Exhaustive CHAID trees with ordinal dependent variables, the SCORES subcommand allows you to specify scores that are used to scale the dependent variable. By default, scores are assigned in sorted value order: the first category is assigned a score of 1, the second category is assigned a score of 2, and so forth.

- If the growing method is CRT or QUEST, this subcommand is ignored, and a warning is issued.
- If the dependent variable is nominal or scale, this subcommand is ignored, and a warning is issued.

The subcommand name is followed by an equals sign (=) and one of the two following alternatives:

- EQUALINCREMENTS**            *Scores are ordinal ranks.* The lowest category of the dependent variable is assigned a score of 1, the next highest category is assigned a score of 2, and so on. This is the default.
- CUSTOM**                        *User-specified scores.* For each category of the dependent variable, specify the category value followed by a score in square brackets.

### **Custom Scores**

- Scores must be numeric and unique. You cannot assign the same score to more than one category.
- Scores must be specified for all values of the dependent variable included in the analysis (either all non-missing values found in the data or all values defined on the DEPCATEGORIES subcommand).
- Category values must be consistent with the data type of the dependent variable. String and date values must be quoted. Date values must be consistent with the variable's print format.
- If you specify multiple score values for the same category value, the last one is used.
- A warning is issued if you specify a score for a category that does not exist in the data or in the training sample if split-sample validation is in effect. [For more information, see the topic VALIDATION Subcommand on p. 1994.](#)

### **Example**

```
TREE income [o] BY age employment gender region
  /SCORES CUSTOM=1 [1] 2 [2.5] 3 [8].
```

- By default, the CHAID method is used to grow the tree.
- `CUSTOM` contains one score specification for each category of the ordinal dependent variable.
- Category 1 is assigned a score of 1; category 2 is assigned a score of 2.5; and category 3 is assigned a score of 8. In other words, scores are used to rescale income so there is a smaller gap between categories 1 and 2 than between 2 and 3.

## ***PROFITS Subcommand***

For categorical dependent variables, the `PROFITS` subcommand allows you to assign revenue and expense values to levels of the dependent variable. Revenues and expenses are used in tables and charts that summarize expected profits and return on investment (ROI). Profit is computed as revenue minus expense.

If the dependent variable is scale, this subcommand is ignored and warning is issued.

- The subcommand name is followed by an equals sign (=), the keyword `CUSTOM`, and a list of dependent variable values, each followed by a pair of values, representing revenues and expenses respectively, enclosed in square brackets.
- Revenue and expense values must be numeric and must be specified for all values of the dependent variable included in the analysis (either all non-missing values found in the data or all values defined on the `DEPCATEGORIES` subcommand).
- Category values must be consistent with the data type of the dependent variable. String and date values must be quoted. Date values must be consistent with the variable's print format.
- If you specify the same category value more than once, the expense and revenue values for the last one are used.
- A warning is issued if you specify a category of the dependent variable that does not exist in the data or in the training sample if split-sample validation is in effect. [For more information, see the topic VALIDATION Subcommand on p. 1994.](#)

### ***Example***

```
TREE income [o] BY age employment gender region
  /PROFITS CUSTOM=1 [.1 .25]
                2 [.5 .25]
                3 [15 .25].
```

- Expected revenue and expense values are defined for each income category.
- Revenues vary across groups. For group 1, the lowest income group, expected revenue is 0.10 per individual. For group 2, the middle income group, expected revenue is 0.50 per individual. For group 3, the high income group, expected revenue is 15 per individual.
- The matrix specifies a fixed expense for each group. For each group, expected expense is 0.25 per person.



## ***INFLUENCE Subcommand***

The `INFLUENCE` subcommand defines an optional influence variable that defines how much influence a case has on the tree-growing process. Cases with lower influence values have less influence, cases with higher values have more.

- The variable must be numeric.
- The dependent variable cannot be used as the influence variable.
- The weight variable (see the `WEIGHT` command) cannot be used as the influence variable.
- For `QUEST`, this subcommand is ignored and a warning is issued.

## ***OUTFILE Subcommand***

The `OUTFILE` subcommand writes the tree model to an XML file.

<code>TRAININGMODEL='filespec'</code>	<i>Writes the model to the specified file.</i> For split-sample validated trees, this is the model for the training sample.
<code>TESTMODEL='filespec'</code>	<i>Writes the model for the test sample to the specified file.</i> Available only with <code>/VALIDATION TYPE=SPLITSAMPLE</code> . Ignored for unvalidated and cross-validated models.

### ***Example***

```
TREE income [o] BY age employment gender region  
/OUTFILE TRAININGMODEL='/myfolder/model.xml'.
```

## ***MISSING Subcommand***

The `MISSING` subcommand controls the handling of nominal, user-missing, independent (predictor) variable values.

- Handling of ordinal and scale user-missing independent variable values varies between growing methods.
- By default, if the values of all independent variables are system- or user-missing, the case is excluded.
- Handling of nominal dependent variables is controlled by the `DEPCATEGORIES` subcommand.
- For the dependent variable, cases with system-missing or user-missing ordinal or scale values are always excluded.

**NOMINALMISSING Keyword**

The `NOMINALMISSING` keyword controls the treatment of user-missing values of nominal independent (predictor) variables.

- MISSING**      *Handling of user-missing values of nominal predictor variables depends on the growing method. This is the default.*
- VALID**        *Treat user-missing values of nominal independent variables as valid values.*  
User-missing values of nominal independent variables are treated as ordinary values in tree growing and classification.

**Method-Dependent Rules**

If some, but not all independent variable values are system-or user-missing:

- For CHAID and Exhaustive CHAID, system- and user-missing scale and ordinal independent variable values are included in the tree-growing process as a “floating” category that is allowed to merge with other categories in tree nodes. By default, nominal user-missing values are also handled in this fashion.
- For CRT and QUEST, cases with missing independent variable values are excluded from the tree-growing process but are classified using surrogate predictors, if surrogates are included in the method. By default, nominal user-missing values are also handled in this fashion.

**Weight Variables**

If the analysis uses frequency weights or an influence variable, all cases with zero, negative, system-missing, or user-missing values of either variable are excluded from the analysis.

**TARGETRESPONSE Subcommand**

The `TARGETRESPONSE` subcommand displays a description of the attributes of each terminal node and the percentage of cases in that node with the specified target value. The output includes a table of percentages and cumulative percentages of the target value in each terminal node and a chart of cumulative response rates.

The `TARGETRESPONSE` subcommand is only executed under the following conditions:

- The subcommand must be explicitly specified. There is no default setting.
- The target variable must be nominal or ordinal.
- Target values must be specified on the `DEPCATEGORIES` subcommand. [For more information, see the topic DEPCATEGORIES Subcommand on p. 1984.](#)

The results are sorted in descending order of percentages. The table can be color-coded based on a specified cumulative minimum percentage, with rows at or above the minimum colored green and rows below the minimum colored red. The minimum value is also displayed as a reference line on the chart. If more than one target value is specified, a separate table and chart are displayed for each target value.

You must specify one of the following alternatives:

- RATE=NONE**                      *No minimum percentage.* The table is not color-coded, and there is no reference line on the chart.
- RATE=value**                      *Use the specified percentage value as the minimum.* The value must be greater than 0 and less than 100.

***Special Considerations***

- The terminal node attributes displayed in the table represent a general description. They are not a substitute for precise rules. [For more information, see the topic RULES Subcommand on p. 1989.](#)
- For the CHAID and Exhaustive CHAID methods, the attributes displayed in the table do not include any missing data conditions, and terminal nodes that represent the missing categories of a predictor are not included in the output. [For more information, see the topic MISSING Subcommand on p. 2003.](#)
- For the CRT and QUEST methods, surrogate values are not included in the list of attribute conditions. [For more information, see the topic METHOD Subcommand on p. 1992.](#)

# TSAPPLY

TSAPPLY is available in the Forecasting option.

*Note:* Square brackets that are shown in the TSAPPLY syntax chart are not used to indicate optional elements. Where indicated, they are required parts of syntax when a list of values is specified, but they may be omitted when a single value is specified. Equals signs (=) that are used in the syntax chart are required elements. The MODEL subcommand is required. All other subcommands are optional.

TSAPPLY

## *Global subcommands:*

```
/MODELSUMMARY PRINT=[MODELFIT** RESIDACF RESIDPACF NONE]
                PLOT=[SRSQUARE RSQUARE RMSE MAPE MAE MAXAPE
                    MAXAE NORMBIC RESIDACF RESIDPACF]

/MODELSTATISTICS DISPLAY={YES**}
                  {NO }
                MODELFIT=[SRSQUARE** RSQUARE RMSE MAPE MAE
                    MAXAPE MAXAE NORMBIC]

/MODELDETAILS PRINT=[PARAMETERS RESIDACF RESIDPACF FORECASTS]
                PLOT=[RESIDACF RESIDPACF]

/SERIESPLOT OBSERVED FORECAST FIT FORECASTCI FITCI

/OUTPUTFILTER DISPLAY={ALLMODELS**
                    {[BESTFIT({N=integer }) WORSTFIT({N=integer })]}
                    {PCT=percent} {PCT=percent}}
                MODELFIT={SRSQUARE**
                    {RSQUARE }
                    {RMSE }
                    {MAPE }
                    {MAE }
                    {MAXAPE }
                    {MAXAE }
                    {NORMBIC }

/SAVE PREDICTED(rootname)
      LCL(rootname)
      UCL(rootname)
      NRESIDUAL(rootname)

/AUXILIARY REESTIMATE={NO**} CILEVEL={95** } MAXACFLAGS={24** }
              {YES } {number} {integer}
          SEASONLENGTH=integer

/MISSING USERMISSING={EXCLUDE**}
                    {INCLUDE }
```

## *MODEL (local subcommand):*

```
/MODEL FILE=file {DROP=['ModelID' 'ModelID' ...]}
          {KEEP=['ModelID' 'ModelID' ...]}
          OUTFILE=file
```

**\*\*Default if the subcommand or keyword is omitted.**

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 14.0

- Command introduced.

### **Example**

```
PREDICT THRU YEAR 2006 MONTH 6.  
TSAPPLY  
  /MODELDETAILS PRINT=FORECASTS  
  /MODEL FILE='/models/models.xml'.
```

## **Overview**

The TSAPPLY procedure loads existing time series models from an external file and applies them to data. Models are generated by using the TSMODEL procedure. You can use TSAPPLY to obtain forecasts for series for which new or revised data are available.

### **Options**

**Models.** By default, parameter estimates saved in the model file are used to produce forecasts. Optionally, model parameters can be reestimated. Reestimation of model parameters has no effect on the model structure. For example, an ARIMA(1,0,1) model will remain so, but the autoregressive and moving-average parameters will be reestimated by using the data in the active dataset. Outliers, if any, are always taken from the model file.

**Output.** TSAPPLY offers the same output as the TSMODEL procedure. However, autocorrelations and partial autocorrelations are available only if model parameters are reestimated.

**Saved Variables.** You can save forecasts and their confidence intervals to the active dataset. Fit values and noise residuals are available if model parameters are reestimated.

**Missing Values.** You can control whether user-missing values are treated as valid or invalid values.

### **Basic Specification**

- The basic specification is a MODEL subcommand that specifies a model file. By default, all models in the file are loaded.
- Default output includes a summary of the distribution of goodness of fit across loaded models and a table of stationary *R*-square, Ljung-Box *Q*, and number of outliers by model.

### **Syntax Rules**

- The following subcommands are global: MODELSUMMARY, MODELSTATISTICS, MODELDETAILS, SERIESPLOT, OUTPUTFILTER, SAVE, AUXILIARY, and MISSING.
- Each global subcommand is optional, may be used only once, and must appear before any MODEL subcommand.

- The `MODEL` subcommand must be used at least once.
- An error occurs if a keyword or attribute is specified more than once within a subcommand.
- All subcommands other than `MODEL` are optional.
- Subcommand names and keywords must be spelled in full.
- Equals signs (=) that are shown in the syntax chart are required.
- An error is issued if any subcommand is empty.

### ***Operations***

- Models are applied to variables in the active dataset with the same names as the variables that are specified in the model.
- `TSAPPLY` honors time intervals and periodicity specified with the `DATE` command.
- The procedure verifies that date variables that are implied by the current date specification exist and that their values coincide with the date specification within the estimation period. If the data are filtered, values of date variables are verified within the filtered subset of data.
- If you choose to reestimate model parameters, reestimation happens within the currently active `USE` period. If the model has outliers, the outlier periods must be inside the effective `USE` period. The `USE` period is ignored if model parameters are not reestimated.
- The `PREDICT` command defines the end of the forecast period for tables, charts, and model variables that are produced by `TSAPPLY`.
- `TSAPPLY` does not honor the following commands: `MODEL NAME`, `SAVE MODEL`, and `READ MODEL`.
- `TSAPPLY` does not honor `TSET`. It does, however, provide options for missing-value handling, setting the width of confidence intervals, and setting the maximum number of lags displayed for autocorrelations.
- The `TDISPLAY` command does not display models that are saved by using `TSAPPLY`.

### ***Limitations***

- An error occurs if `SPLIT FILE` is in effect.
- `WEIGHT` is ignored with a warning.

## ***Examples***

This section provides simple examples that are designed to get you started with the `TSAPPLY` procedure. Further examples that are specific to each subcommand are provided in the subcommand topics.

### ***Forecasting***

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /MODEL FILE='/models/models.xml' .
```

- The `PREDICT` command is used to specify the forecast period for the `TSAPPLY` procedure.

- The `MODELDETAILS` subcommand specifies that the output includes a table containing the forecasts.
- The `MODEL` subcommand specifies that all models in the file `/models/models.xml` will be applied to the active dataset. Models are applied to variables in the active dataset with the same names as the variables that are specified in the model.
- Forecasts are created by using the model parameters from `models.xml`, without reestimating any of those parameters. This behavior is the default behavior.

### **Forecasting with Reestimated Model Parameters**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /AUXILIARY REESTIMATE=YES
  /MODEL FILE='/models/models.xml'
  OUTFILE='/models/models_updated.xml'.
```

- The `AUXILIARY` subcommand specifies that model parameters for all models in `/models/models.xml` will be reestimated prior to forecasting. Model parameters are reestimated by using the data in the active dataset.
- Reestimation of model parameters has no effect on the model structure. For example, an `ARIMA(1,0,1)` model will remain so, but the autoregressive and moving-average parameters will be reestimated.
- The `OUTFILE` keyword specifies that the reestimated models will be saved to the file `/models/models_updated.xml`.

### **Applying a Selected Subset of Models**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /MODEL FILE='/models/models.xml' KEEP=['CustomArima_1'].
```

- The `KEEP` keyword specifies that only the model named `CustomArima_1` will be applied to the active dataset. All other models in `/models/models.xml` will be excluded.

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /MODEL FILE='/models/models.xml' DROP=['Expert_1' 'Expert_2'].
```

- The `DROP` keyword specifies that the models named `Expert_1` and `Expert_2` will be excluded. All other models in `/models/models.xml` will be applied to the active dataset.

## Goodness-of-Fit Measures

The following model goodness-of-fit measures are available in TSAPPLY:

<b>SRSQUARE</b>	<b>Stationary R-squared.</b> A measure that compares the stationary part of the model to a simple mean model. This measure is preferable to ordinary R-squared when there is a trend or seasonal pattern. Stationary R-squared can be negative with a range of negative infinity to 1. Negative values mean that the model under consideration is worse than the baseline model. Positive values mean that the model under consideration is better than the baseline model.
<b>RSQUARE</b>	<b>R-squared.</b> An estimate of the proportion of the total variation in the series that is explained by the model. This measure is most useful when the series is stationary. R-squared can be negative with a range of negative infinity to 1. Negative values mean that the model under consideration is worse than the baseline model. Positive values mean that the model under consideration is better than the baseline model.
<b>RMSE</b>	<b>RMSE.</b> Root Mean Square Error. The square root of mean square error. A measure of how much a dependent series varies from its model-predicted level, expressed in the same units as the dependent series.
<b>MAPE</b>	<b>MAPE.</b> Mean Absolute Percentage Error. A measure of how much a dependent series varies from its model-predicted level. It is independent of the units used and can therefore be used to compare series with different units.
<b>MAE</b>	<b>MAE.</b> Mean absolute error. Measures how much the series varies from its model-predicted level. MAE is reported in the original series units.
<b>MAXAPE</b>	<b>MaxAPE.</b> Maximum Absolute Percentage Error. The largest forecasted error, expressed as a percentage. This measure is useful for imagining a worst-case scenario for your forecasts.
<b>MAXAE</b>	<b>MaxAE.</b> Maximum Absolute Error. The largest forecasted error, expressed in the same units as the dependent series. Like MaxAPE, it is useful for imagining the worst-case scenario for your forecasts. Maximum absolute error and maximum absolute percentage error may occur at different series points—for example, when the absolute error for a large series value is slightly larger than the absolute error for a small series value. In that case, the maximum absolute error will occur at the larger series value and the maximum absolute percentage error will occur at the smaller series value.
<b>NORMBIC</b>	<b>Normalized BIC.</b> Normalized Bayesian Information Criterion. A general measure of the overall fit of a model that attempts to account for model complexity. It is a score based upon the mean square error and includes a penalty for the number of parameters in the model and the length of the series. The penalty removes the advantage of models with more parameters, making the statistic easy to compare across different models for the same series.

## MODELSUMMARY Subcommand

The MODELSUMMARY subcommand controls the display of tables and charts that summarize goodness of fit, residual autocorrelations, and residual partial autocorrelations across estimated models. Each keyword is followed by an equals sign (=) and one or more of the available options enclosed in square brackets.

### Example

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /AUXILIARY REESTIMATE=YES
```



```
/MODELSUMMARY PLOT=[SRSQUARE MAXAPE]
/MODEL FILE=' /models/models.xml ' .
```

- The output includes two histograms: one histogram for stationary  $R$ -squared and one histogram for the maximum absolute percentage error. Each histogram consists of results across all models.

### **PRINT Keyword**

The PRINT keyword controls the display of model summary tables.

<b>MODELFIT</b>	<i>Goodness of fit.</i> Table of summary statistics and percentiles for stationary $R$ -square, $R$ -square, root mean square error, mean absolute percentage error, mean absolute error, maximum absolute percentage error, maximum absolute error, and normalized Bayesian Information Criterion. Note that if models are not reestimated, model fit statistics are loaded from the model file. This table is displayed by default.
<b>RESIDACF</b>	<i>Residual autocorrelation function.</i> Table of summary statistics and percentiles for autocorrelations of the residuals. RESIDACF is ignored with a warning if REESTIMATE=NO.
<b>RESIDPACF</b>	<i>Residual partial autocorrelation function.</i> Table of summary statistics and percentiles for partial autocorrelations of the residuals. RESIDPACF is ignored with a warning if REESTIMATE=NO.
<b>NONE</b>	<i>No tables are shown.</i> An error occurs if NONE is used in combination with any other PRINT option.

### **PLOT Keyword**

The PLOT keyword controls the display of model summary charts. By default no charts are shown. (For detailed definitions of the following terms, see [Goodness-of-Fit Measures](#) on p. 2010.)

<b>SRSQUARE</b>	<i>Histogram of Stationary R-Square.</i>
<b>RSQUARE</b>	<i>Histogram of R-Square.</i>
<b>RMSE</b>	<i>Histogram of Root Mean Square Error.</i>
<b>MAPE</b>	<i>Histogram of Mean Absolute Percentage Error.</i>
<b>MAE</b>	<i>Histogram of Mean Absolute Error.</i>
<b>MAXAPE</b>	<i>Histogram of Maximum Absolute Percentage Error.</i>
<b>MAXAE</b>	<i>Histogram of Maximum Absolute Error.</i>
<b>NORMBIC</b>	<i>Histogram of Normalized Bayesian Information Criterion (BIC).</i>
<b>RESIDACF</b>	<i>Boxplot of Residual Autocorrelation Function by Lag.</i> RESIDACF is ignored with a warning if REESTIMATE=NO.
<b>RESIDPACF</b>	<i>Boxplot of Residual Partial Autocorrelation Function by Lag.</i> RESIDPACF is ignored with a warning if REESTIMATE=NO.

## **MODELSTATISTICS Subcommand**

The MODELSTATISTICS subcommand controls display of a table listing all models along with chosen goodness-of-fit statistics. Note that if models are not reestimated, model fit statistics are loaded from the model file.

**Example**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /AUXILIARY REESTIMATE=YES
  /MODELSTATISTICS DISPLAY=YES MODELFIT=RSQUARE
  /MODEL FILE=' /models/models.xml '.
```

- The output includes a table displaying the value of *R*-squared for each model.

**DISPLAY Keyword**

The DISPLAY keyword controls whether the model statistics table is shown.

<b>YES</b>	<i>Model statistics table is shown.</i> Table of model goodness of fit, Ljung-Box Q statistic, and number of outliers by model. This setting is the default. The Q statistic measures the degree of pattern in the residuals. Large values of Q in relation to its degrees of freedom indicate that model residuals are not randomly distributed.
<b>NO</b>	<i>Model statistics table is not shown.</i>

**MODELFIT Keyword**

The MODELFIT keyword controls which fit statistics are shown in the model statistics table, and this keyword is ignored if DISPLAY=NO. The keyword is followed by an equals sign (=) and one or more of the following options enclosed in square brackets. (For detailed definitions of the following terms, see [Goodness-of-Fit Measures](#) on p. 2010.)

<b>SRSQUARE</b>	<i>Stationary R-square.</i> This setting is the default.
<b>RSQUARE</b>	<i>R-Square.</i>
<b>RMSE</b>	<i>Root Mean Square Error.</i>
<b>MAPE</b>	<i>Mean Absolute Percentage Error.</i>
<b>MAE</b>	<i>Mean Absolute Error.</i>
<b>MAXAPE</b>	<i>Maximum Absolute Percentage Error.</i>
<b>MAXAE</b>	<i>Maximum Absolute Error.</i>
<b>NORMBIC</b>	<i>Normalized Bayesian Information Criterion (BIC).</i>

**MODELDETAILS Subcommand**

The MODELDETAILS subcommand controls the display of model parameters, model forecasts, and autocorrelations of noise residuals for individual models. Each keyword is followed by an equals sign (=) and one or more of the available options enclosed in square brackets.

**Example**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /AUXILIARY REESTIMATE=YES
  /MODELDETAILS PRINT=[PARAMETERS FORECASTS]
```

```
/MODEL FILE=' /models/models.xml ' .
```

- The output includes tables displaying the model parameters and forecasts for each model.

### **PRINT Keyword**

The PRINT keyword controls display of tables for individual models. By default, no tables are shown.

<b>PARAMETERS</b>	<i>Model parameter estimates.</i> Shows separate tables for ARIMA and exponential smoothing models. If outliers exist, parameter estimates for outliers are also displayed. Note that if models are not reestimated, parameter estimates are loaded from the model file.
<b>RESIDACF</b>	<i>Residual autocorrelation function.</i> Shows residual autocorrelations by lag. RESIDACF is ignored with a warning if REESTIMATE=NO.
<b>RESIDPACF</b>	<i>Residual partial autocorrelation function.</i> Shows residual partial autocorrelations by lag. RESIDPACF is ignored with a warning if REESTIMATE=NO.
<b>FORECASTS</b>	<i>Forecasts and confidence intervals.</i> Shows model forecasts and confidence intervals.

### **PLOT Keyword**

The PLOT keyword controls display of charts for individual models. By default no charts are shown. PLOT is ignored with a warning if REESTIMATE=NO.

<b>RESIDACF</b>	<i>Residual autocorrelation function.</i> Shows residual autocorrelations by lag.
<b>RESIDPACF</b>	<i>Residual partial autocorrelation function.</i> Shows residual partial autocorrelations by lag.

## **SERIESPLOT Subcommand**

The SERIESPLOT subcommand allows you to obtain plots of predicted values, observed values, and confidence intervals for each model. By default, no plots are shown. The subcommand is followed by one or more of the following keywords:

<b>OBSERVED</b>	<i>Displays all observed values of the dependent series.</i> OBSERVED is ignored with a warning if REESTIMATE=NO.
<b>FORECAST</b>	<i>Displays model-predicted values within the forecast period.</i>
<b>FIT</b>	<i>Displays model-predicted values within the estimation period.</i> FIT is ignored with a warning if REESTIMATE=NO.
<b>FORECASTCI</b>	<i>Displays upper and lower confidence limits within the forecast period.</i>
<b>FITCI</b>	<i>Displays upper and lower confidence limits within the estimation period.</i> FITCI is ignored with a warning if REESTIMATE=NO.

### **Example**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /AUXILIARY REESTIMATE=YES
  /SERIESPLOT OBSERVED FORECAST
```

```
/MODEL FILE=' /models/models.xml ' .
```

- The output includes plots of the model-predicted values in the forecast period, as well as all observed values, for each model.

## ***OUTPUTFILTER Subcommand***

The OUTPUTFILTER subcommand controls which models are included in the model statistics table (MODELSTATISTICS subcommand), detailed model output (MODELDETAILS subcommand), and time series plots (SERIESPLOT subcommand).

- By default, all models are included in the output. OUTPUTFILTER can be used to display only the best-fitting or worst-fitting models (or both). If you request both, two sets of output are displayed.
- If model parameters are not reestimated, models are selected based on goodness-of-fit statistics that are loaded from the model file.
- OUTPUTFILTER is ignored if no output is requested on the MODELSTATISTICS, MODELDETAILS, or SERIESPLOT subcommands.
- OUTPUTFILTER has no effect on output from the MODELSUMMARY subcommand.

### ***Example***

```
PREDICT THRU YEAR 2006 MONTH 6.  
TSAPPLY  
  /MODELDETAILS PRINT=FORECASTS  
  /OUTPUTFILTER DISPLAY=[BESTFIT(N=5) WORSTFIT(PCT=10)]  
  /MODEL FILE=' /models/models.xml ' .
```

- The output consists of two sets of results: those results for the 5 best-fitting models and those results for models with fit values in the bottom 10%. The stationary *R*-squared value is used as the goodness-of-fit measure.

### ***DISPLAY Keyword***

The DISPLAY keyword controls which models are included in MODELSTATISTICS, MODELDETAILS, and SERIESPLOT output.

- By default, all models are displayed (`DISPLAY=ALLMODELS`).
- To display only those models with extreme goodness-of-fit values, specify `BESTFIT`, `WORSTFIT`, or both, in square brackets. One or both keywords must be specified.

<b>ALLMODELS</b>	<i>All models are displayed.</i> All models that could be loaded ( <code>REESTIMATE=NO</code> ) or computed ( <code>REESTIMATE=YES</code> ) are shown. This setting is the default.
<b>BESTFIT</b>	<i>Models having the highest fit values are shown.</i> To display the models with the <code>N</code> highest fit values, specify <code>BESTFIT</code> followed by <code>N=count</code> in parentheses—for example, <code>BESTFIT(N=5)</code> . The count must be a positive integer. If the count exceeds the number of estimated models, all models are shown. To display the models with fit values in the top <code>N%</code> , specify <code>PCT=percent</code> in parentheses. For example, <code>BESTFIT(PCT=5)</code> displays models with fit values in the top 5% among estimated models. The percentage value must be greater than zero and less than 100.
<b>WORSTFIT</b>	<i>Models having the lowest fit values are shown.</i> To display the models with the <code>N</code> lowest fit values, specify <code>WORSTFIT</code> followed by <code>N=count</code> in parentheses—for example, <code>WORSTFIT(N=5)</code> . The count must be a positive integer. If the count exceeds the number of models estimated, all models are shown. To display the models with fit values in the lowest <code>N%</code> , specify <code>PCT=percent</code> in parentheses. For example, <code>WORSTFIT(PCT=5)</code> displays models with fit values in the bottom 5% among estimated models. The percentage value must be greater than zero and less than 100.

### **MODELFIT Keyword**

The `MODELFIT` keyword specifies the fit measure that is used to filter models. `MODELFIT` is ignored if `DISPLAY=ALLMODELS`.

Specify one of the following options. (For detailed definitions of the following terms, see [Goodness-of-Fit Measures](#) on p. 2010.)

<b>SRSQUARE</b>	<i>Stationary R-Square.</i> This setting is the default.
<b>RSQUARE</b>	<i>R-Square.</i>
<b>RMSE</b>	<i>Root Mean Square Error.</i>
<b>MAPE</b>	<i>Mean Absolute Percentage Error.</i>
<b>MAE</b>	<i>Mean Absolute Error.</i>
<b>MAXAPE</b>	<i>Maximum Absolute Percentage Error.</i>
<b>MAXAE</b>	<i>Maximum Absolute Error.</i>
<b>NORMBIC</b>	<i>Normalized Bayesian Information Criterion (BIC).</i>

## **SAVE Subcommand**

The `SAVE` subcommand is used to save new variables representing predicted values, residuals, and confidence intervals to the active dataset. By default, no new variables are saved to the active dataset.

- Specify one or more keywords, each keyword followed by an optional rootname to be used as the prefix for new variable names. Enclose the root name in parentheses. Each keyword gives rise to one new variable for each dependent variable.
- The root name, if specified, must conform to the rules for valid variable names.
- If no root name is specified, `TSAPPLY` uses a default name.

- The full variable name is the concatenation of the root name, the name of the associated dependent variable, and a model identifier. The variable name is extended if necessary to avoid variable naming conflicts.
- The PREDICT command controls whether new cases are added to the file when new variables are saved. New cases are added if PREDICT specifies a forecast period that extends beyond the length of the dependent-variable series.
- SAVE is ignored with a warning if temporary transformations are in effect.

**Example**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /AUXILIARY REESTIMATE=YES
  /SAVE PREDICTED(ApplyPred) NRESIDUAL
  /MODEL FILE=' /models/models.xml'.
```

- Two new variables are created for each of the models: one variable contains the model predictions, and the other variable contains the noise residuals.
- The root name (prefix) for the variables containing model predictions is *ApplyPred*.
- The root name for the variables containing noise residuals is *NResidual* (the default).

<b>PREDICTED(rootname)</b>	<i>Model fit and forecast values.</i> The default root name is <i>Predicted</i> . If REESTIMATE=NO, only forecasts are produced.
<b>LCL(rootname)</b>	<i>Lower confidence limits.</i> The default root name is <i>LCL</i> . If REESTIMATE=NO, lower confidence limits are only produced for forecasts.
<b>UCL(rootname)</b>	<i>Upper confidence limits.</i> The default root name is <i>UCL</i> . If REESTIMATE=NO, upper confidence limits are only produced for forecasts.
<b>NRESIDUAL(rootname)</b>	<i>Noise residuals.</i> The default rootname is <i>NResidual</i> . When transformations of the dependent variable are performed (for example, natural log), these residuals are the residuals for the transformed series. NRESIDUAL is ignored with a warning if REESTIMATE=NO.

**AUXILIARY Subcommand**

The AUXILIARY subcommand is used to specify whether model parameters are reestimated, set the confidence interval level for forecasts, set the maximum number of lags for autocorrelation and partial autocorrelation plots and tables, and specify the season length.

**Example**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=[PARAMETERS FORECASTS]
  PLOT=[RESIDACF RESIDPACF]
  /AUXILIARY REESTIMATE=YES MAXACFLAGS=36
  /MODEL FILE=' /models/models.xml'.
```

**REESTIMATE Keyword**

REESTIMATE controls whether model parameters are reestimated. If REESTIMATE=NO (the default), parameter estimates that are recorded in the model file are used to produce forecasts. If REESTIMATE=YES, forecasts reflect reestimated model parameters. Outliers, if any, are always taken from the model file.

Note that if REESTIMATE=NO, any parameter estimates or goodness-of-fit statistics that are displayed in output are loaded from the model file and reflect the data that were used when each model was developed (or last updated). And if you use the OUTPUTFILTER subcommand, models will be selected on the basis of goodness-of-fit statistics that are recorded in the model file. In addition, forecasts will not take into account historical data—for either dependent or independent variables—in the active dataset. You must set REESTIMATE=YES if you want historical data to impact the forecasts. Finally, forecasts do not take into account values of the dependent series in the forecast period—but the forecasts do take into account values of independent variables in the forecast period. If you have more current values of the dependent series and want them to be included in the forecasts, you need to reestimate, adjusting the estimation period to include these values.

**CILEVEL Keyword**

The CILEVEL keyword sets the confidence level.

- Specify a positive number that is less than 100.
- The default is 95.

**MAXACFLAGS Keyword**

The MAXACFLAGS keyword sets the maximum number of lags displayed in residual autocorrelation and partial autocorrelation tables and plots. MAXACFLAGS is ignored if REESTIMATE=NO.

- Specify a positive integer.
- The default is 24.

**SEASONLENGTH Keyword**

The SEASONLENGTH keyword specifies the length of the seasonal period (the number of observations in one period or season) for the data.

- To adjust the season length, specify a positive integer.
- If SEASONLENGTH is not specified, the periodicity that is specified by using the DATE command defines the season length.
- Note that SEASONLENGTH is used to override the date specification for the active dataset to make it consistent with the date specification from the model file. An error occurs if these two specifications are not consistent.

**MISSING Subcommand**

The MISSING subcommand controls the handling of user-missing values.

- By default, user-missing values are treated as missing (invalid) data.
- System-missing values are always treated as invalid.
- Cases with missing values of a dependent variable that occur within the estimation period are included in the model. The specific handling of the missing value depends on the estimation method.
- For ARIMA models, a warning is issued if a predictor has any missing values within the estimation period. Any models involving the predictor are not reestimated.
- If any predictor has missing values within the forecast period, the procedure issues a warning and forecasts as far as it can.

**Example**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSAPPLY
  /MODELDETAILS PRINT=FORECASTS
  /AUXILIARY REESTIMATE=YES
  /MISSING USERMISSING=INCLUDE
  /MODEL FILE=' /models/models.xml ' .
```

**USERMISSING Keyword**

The USERMISSING keyword controls the treatment of user-missing values and is required.

- |                |  |
|----------------|--|
| <b>EXCLUDE</b> | <i>Exclude user-missing values.</i> User-missing values are treated as missing. This setting is the default. |
| <b>INCLUDE</b> | <i>Include user-missing values.</i> User-missing values are treated as valid data.                           |

**MODEL Subcommand**

The MODEL subcommand is required and specifies an external file, referred to as a **model file**, containing models that are developed by using TSMODEL. A model file specifies lag structure, parameters, outliers, and variables for each model.

- TSAPPLY assumes that the variables that are specified in the model file exist in the active dataset.
- You can use multiple MODEL subcommands to apply two or more sets of models to your data. Each MODEL subcommand is used to load models from a different model file. For example, one model file might contain models for series that represent unit sales, and another model file might contain models for series that represent revenue.

**FILE Keyword**

The FILE keyword is used to specify an external file containing models, and the keyword is required.

- A warning is issued if a model variable that is specified in the file does not exist in the active dataset or is not numeric. Models that refer to the variable are ignored.



- A warning is issued if a model specifies a variable transformation (square root or natural log) and one or more model variables to be transformed contain values that are invalid for the specified transformation—for example, negative values. Any such models will not be applied.
- An error occurs if the date interval or cycle length that is specified in the file is inconsistent with the date specification for the active dataset. The `SEASONLENGTH` keyword on the `AUXILIARY` subcommand can be used to override the date specification for the active dataset.
- An error occurs if the file cannot be read or is invalid.
- An error occurs if two or more `MODEL` subcommands specify the same file.
- When multiple `MODEL` subcommands are used, a warning occurs if the same model ID occurs in more than one model file. The first model with that ID is loaded. The check for duplicate model IDs takes place after `DROP` and `KEEP` are applied to each model file.
- A warning is issued, and forecasts are not available for a model, if `REESTIMATE=NO` and the model's forecast period is inconsistent with the current `USE` or `PREDICT` interval. Specifically, forecasts are not available if model forecasts would start before the beginning of the `USE` period or after the end of the `PREDICT` interval.

### ***DROP Keyword***

The `DROP` keyword is used to exclude models.

- Specify one or more quoted model identifiers, enclosed in square brackets—for example, `['Model_1' 'Model_2']`. All other models in the file are loaded and applied to the data.
- Redundant items in the `DROP` list are ignored.
- An error occurs if all models are dropped.
- Model identifiers that do not exist in the file are ignored.

### ***KEEP Keyword***

The `KEEP` keyword specifies models to be applied.

- Specify one or more quoted model identifiers, enclosed in square brackets—for example, `['Model_1' 'Model_2']`. All other models in the file are excluded.
- Redundant items in the `KEEP` list are ignored.
- Model identifiers that do not exist in the file are ignored.
- `DROP` and `KEEP` are mutually exclusive. An error occurs if both are specified.

### ***OUTFILE Keyword***

The `OUTFILE` keyword specifies an external file to which updated models are written.

- You can use `OUTFILE` to save models whose parameters are reestimated. `OUTFILE` is ignored with a warning if `REESTIMATE=NO`.
- The filename must be specified in full. No extension is supplied.
- Models that are excluded by using `DROP` or `KEEP` are not saved.
- You can specify the same external file on multiple `MODEL` subcommands. Models from the different subcommands are then saved to that single file.

# TSET

```
TSET
  [PRINT={DEFAULT**}]  [/NEWVAR={CURRENT**}]  [/MXAUTO={16**}]
    {BRIEF   }           {NONE   }             {lags}
    {DETAILED}           {ALL    }
  [/MXCROSS={7**}]  [/MXNEWVARS={60**}]  [/MXPREDICT={60**}]
    {lags}             {n   }             {n   }
  [/MISSING={EXCLUDE**}]  [/CIN={95**}]  [/TOLER={0.0001**}]
    {INCLUDE  }         {value}             {value }
  [/CNVERGE={0.001**}]  [/ACFSE={IND**}]
    {value  }           {MA   }
  [/PERIOD=n]          [/ID=varname]
  [/{CONSTANT**}]
    {NOCONSTANT}
  [/DEFAULT]
```

\*\*Default if the subcommand is omitted.

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
TSET PERIOD 6 NEWVAR NONE MXAUTO 25.
```

## Overview

TSET sets global parameters to be used by procedures that analyze time series and sequence variables. To display the current settings of these parameters, use the TSHOW command.

### Basic Specification

The basic specification is at least one subcommand.

### Subcommand Order

- Subcommands can be specified in any order.

### Syntax Rules

- The slash between subcommands is optional.
- You can specify **DEFAULT** on any subcommand to restore the default setting for that subcommand.
- Subcommand **DEFAULT** restores all TSET subcommands to their defaults.

**Operations**

- TSET takes effect immediately.
- Only the settings specified are affected. All others remain at their previous settings or the default.
- Subcommands on other procedures that perform the same function as subcommands on TSET override the TSET specifications for those procedures.
- Procedures that are affected by TSET specifications are CURVEFIT, CASEPLOT, NPLOT, and TSPLOT.

**DEFAULT Subcommand**

DEFAULT resets all TSET settings back to their defaults. There are no additional specifications on DEFAULT.

**ID Subcommand**

ID specifies a variable whose values are used to label observations in plots.

- The only specification on ID is the name of a variable in the active dataset.
- If ID is not specified, the *DATE\_* variable is used to label observations.
- If ID is specified within the procedure, it overrides the TSET specification for that procedure.

**MISSING Subcommand**

MISSING controls the treatment of user-missing values.

- The specification on MISSING is keyword INCLUDE or EXCLUDE. The default is EXCLUDE.
- INCLUDE indicates that observations with user-missing values should be treated as valid values and included in analyses.
- EXCLUDE indicates that observations with user-missing values should be excluded from analyses.

**MXNEWVARS Subcommand**

MXNEWVARS indicates the maximum number of new variables that can be generated by a procedure.

- The specification on MXNEWVARS indicates the maximum and can be any positive integer.
- The default maximum number is 60 new variables per procedure.

**MXPREDICT Subcommand**

MXPREDICT indicates the maximum number of new cases that can be added to the active dataset per procedure when the PREDICT command is used.

- The specification on `MXPREDICT` can be any positive integer.
- The default maximum number of new cases is 60 per procedure.

## ***NEWVAR Subcommand***

`NEWVAR` controls the creation of new variables in the active dataset.

- The specification on `NEWVAR` can be `CURRENT`, `NONE`, or `ALL`. The default is `CURRENT`.
- `CURRENT` specifies to save new variables and replace any existing variables of the same name.
- `ALL` specifies to save new variables without replacing existing ones.
- `NONE` specifies that no new variables are saved.

## ***PERIOD Subcommand***

`PERIOD` indicates the size of the period to be used for seasonal differencing.

- The specification on `PERIOD` indicates how many observations are in one season or period and can be any positive integer.
- There is no default for the `PERIOD` subcommand.
- The specification on `TSET PERIOD` overrides the periodicity of `DATE` variables.
- If a period is specified within an individual procedure, it overrides the `TSET PERIOD` specification for that procedure.

## ***PRINT Subcommand***

`PRINT` controls how much output is produced.

- The specification on `PRINT` can be `BRIEF`, `DETAILED`, or `DEFAULT`. The amount of output produced by `DEFAULT` is generally between the amount produced by `BRIEF` and `DETAILED`.
- For procedures with multiple iterations, `BRIEF` generally means that final statistics are displayed with no iteration history. `DEFAULT` provides a one-line summary at each iteration in addition to the final statistics. `DETAILED` provides a complete summary of each iteration (where necessary) plus the final statistics.
- For some procedures, the `DEFAULT` and `DETAILED` output is the same. For many of the simpler procedures, `BRIEF`, `DEFAULT`, and `DETAILED` are all the same.

# TSHOW

TSHOW

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Overview

TSHOW displays a list of all the current specifications on the TSET, USE, PREDICT, and DATE commands.

### **Basic Specification**

The command keyword TSHOW is the only specification.

### **Operations**

- TSHOW is executed immediately.
- TSHOW lists every current specification for the TSET, USE, PREDICT, and DATE commands, as well as the default settings.

## Example

TSHOW.

- The TSHOW command produces a list of the current settings on the TSET, USE, PREDICT, and DATE commands.

# TSMODEL

TSMODEL is available in the Forecasting option.

*Note:* Square brackets that are shown in the TSMODEL syntax chart are not used to indicate optional elements. Where indicated, they are required parts of syntax when a list of values is specified, but they may be omitted when a single value is specified. Equals signs (=) that are used in the syntax chart are required elements. The MODEL subcommand is required. All other subcommands are optional.

TSMODEL

## *Global subcommands:*

```
/MODEL SUMMARY PRINT=[MODELFIT** RESIDACF RESIDPACF NONE]
                  PLOT=[SRSQUARE RSQUARE RMSE MAPE MAE MAXAPE
                        MAXAE NORMBIC RESIDACF RESIDPACF]

/MODEL STATISTICS DISPLAY={YES**}
                  {NO }
                  MODELFIT=[SRSQUARE** RSQUARE RMSE MAPE MAE
                        MAXAPE MAXAE NORMBIC]

/MODEL DETAILS PRINT=[PARAMETERS RESIDACF RESIDPACF FORECASTS]
                 PLOT=[RESIDACF RESIDPACF]

/SERIES PLOT OBSERVED FORECAST FIT FORECASTCI FITCI

/OUTPUT FILTER DISPLAY={ALLMODELS**
                      {[BESTFIT({N=integer }) WORSTFIT({N=integer })]}
                      {PCT=percent} {PCT=percent}}
MODELFIT={SRSQUARE**
          {RSQUARE }
          {RMSE }
          {MAPE }
          {MAE }
          {MAXAPE }
          {MAXAE }
          {NORMBIC }

/SAVE PREDICTED(rootname)
      LCL(rootname)
      UCL(rootname)
      NRESIDUAL(rootname)

/AUXILIARY CILEVEL={95** } MAXACFLAGS={24** } SEASONLENGTH=integer
            {number} {integer}

/MISSING USERMISSING={EXCLUDE**}
                    {INCLUDE }
```

## *Model block subcommands:*

```
/MODEL DEPENDENT=varlist
       INDEPENDENT=varspec list
       OUTFILE=model file
       PREFIX='prefix'

{ /EXPERTMODELER TYPE=[ARIMA** EXSMOOTH**] }
  TRYSEASONAL={YES**}
               {NO }
```

```

{ /EXSMOOTH TYPE={SIMPLE          } }
                  {SIMPLESEASONAL }
                  {HOLT           }
                  {BROWN          }
                  {DAMPEDTREND    }
                  {WINTERSADDITIVE }
                  {WINTERSMULTIPLICATIVE}
TRANSFORM={NONE**}
           {SQRT  }
           {LN   }

{ /ARIMA  AR={{integer integer...}} }
          ARSEASONAL={{0**}         }
                   {{integer integer...}}
MA={{integer integer...}}
MASEASONAL={{0**}                 }
           {{integer integer...}}
DIFF={0** }
     {integer}
DIFFSEASONAL={0** }
              {integer}
CONSTANT={YES**}
         {NO  }
TRANSFORM={NONE**}
           {SQRT  }
           {LN   }

/TRANSFERFUNCTION VARIABLES={ALL** }
                        {varlist}
NUM={{0**}             }
   {{integer integer...}}
NUMSEASONAL={{0**}    }
            {{integer integer...}}
DENOM={{0**}          }
      {{integer integer...}}
DENOMSEASONAL={{0**} }
              {{integer integer...}}
DIFF={0** }
     {integer}
DIFFSEASONAL={0** }
              {integer}
DELAY={0** }
      {integer}
TRANSFORM={NONE**}
           {SQRT  }
           {LN   }

/AUTOOUTLIER  DETECT={OFF**}
              {ON  }
              TYPE=[ADDITIVE** LEVELSHIFT** INNOVATIONAL TRANSIENT
                  SEASONALADDITIVE LOCALTREND ADDITIVEPATCH]

/OUTLIER LOCATION=[date specification] TYPE={ADDITIVE          }
                                                {LEVELSHIFT      }
                                                {INNOVATIONAL    }
                                                {TRANSIENT        }
                                                {SEASONALADDITIVE}
                                                {LOCALTREND       }

```

\*\*Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 14.0

- Command introduced.

**Example**

```
TSMODEL  
  /MODEL DEPENDENT=sku1 TO sku10.
```

**Overview**

The TSMODEL procedure estimates exponential smoothing, univariate Autoregressive Integrated Moving Average (ARIMA), and multivariate ARIMA (or transfer function models) models for time series, and the procedure produces forecasts. The procedure includes an Expert Modeler that identifies and estimates an appropriate ARIMA or exponential smoothing model for each dependent-variable series. Alternatively, you can specify a custom ARIMA or exponential smoothing model.

**Options**

**Automatic Model Identification.** The Expert Modeler can identify an appropriate seasonal or nonseasonal ARIMA or exponential smoothing model for each dependent-variable series. If predictor variables are specified, the Expert Modeler selects, for inclusion in ARIMA models, those variables that have a statistically significant relationship with the dependent series. Model variables are transformed where appropriate by using differencing and/or a square root or natural log transformation. You can limit the set of candidate models to ARIMA models only or exponential smoothing models only.

**Custom Models.** You can specify a custom exponential smoothing or ARIMA model for one or more series. Seven exponential smoothing methods are available: simple, simple seasonal, Holt's linear trend, Brown's linear trend, damped trend, Winters' additive, and Winters' multiplicative. For ARIMA models, you can specify seasonal and nonseasonal autoregressive and moving average orders, differencing, as well as transfer functions for predictor variables. For exponential smoothing and ARIMA models, you can request that model variables be transformed prior to model estimation.

**Outliers.** If you use the Expert Modeler or specify a custom ARIMA model, TSMODEL can detect and model outlier time points automatically. The following outlier types can be identified: additive, additive patch, innovational, level shift, transient, seasonal additive, and local trend. Alternatively, if you request a custom ARIMA model, you can specify that one or more time points be modeled as outliers.

**Output.** Available output includes plots and tables that summarize the distribution of model goodness of fit, residual autocorrelations, and residual partial autocorrelations across models. In addition, you can obtain a table of model goodness of fit, Ljung-Box Q, and number of outliers by model. You can also obtain details for each model, including parameter estimates, forecasts, as well as autocorrelation and partial autocorrelation functions. Output can be restricted to the best-fitting or worst-fitting models based on goodness-of fit-values.

**Saved Variables.** You can save fit and forecast values to the active dataset as well as confidence intervals and noise residuals.

**Missing Values.** You can control whether user-missing values are treated as valid or invalid values.



### **Basic Specification**

- The basic specification is a `MODEL` subcommand that specifies one or more dependent-variable series.
- By default, `TSMODEL` uses the Expert Modeler to identify and estimate the best ARIMA or exponential smoothing model for each series.
- Default output includes a summary of the distribution of goodness of fit across estimated models and a table of stationary  $R$ -square, Ljung-Box  $Q$ , and number of outliers by model.

### **Syntax Rules**

- The following subcommands are global and apply to all models specified in a single instance of the `TSMODEL` command: `MODELSUMMARY`, `MODELSTATISTICS`, `MODELDETAILS`, `SERIESPLOT`, `OUTPUTFILTER`, `SAVE`, `AUXILIARY`, and `MISSING`.
- Each global subcommand is optional, may be used only once, and must appear before any `MODEL` subcommand.
- Models are specified in blocks. The `MODEL` subcommand indicates the start of a block. The `MODEL` subcommand must be used at least once.
- The following subcommands apply to the preceding `MODEL` subcommand and constitute—along with a `MODEL` subcommand—a `MODEL` block: `EXPERTMODELER`, `EXSMOOTH`, `ARIMA`, `TRANSFERFUNCTION`, `AUTOOUTLIER`, and `OUTLIER`. An error occurs if any of these subcommands precedes the first `MODEL` subcommand.
- The `EXPERTMODELER`, `EXSMOOTH`, and `ARIMA` subcommands are used within a `MODEL` block to specify the estimation method. If none of these subcommands is specified, the implicit default is `EXPERTMODELER`. An error occurs if more than one of these subcommands is specified within a `MODEL` block.
- `AUTOOUTLIER` may be used only once within a `MODEL` block.
- `TRANSFERFUNCTION` and `OUTLIER` may be used more than once within a `MODEL` block.
- Each keyword may be specified only once within a subcommand.
- Empty subcommands are not allowed; all subcommands must be specified with options.
- All subcommands other than `MODEL` are optional.
- Subcommand names and keywords must be spelled in full.
- Subcommands can be specified in any order, with the exception that global subcommands must precede the first `MODEL` subcommand.
- Equals signs (=) that are shown in the syntax chart are required.

### **Operations**

- `TSMODEL` honors time intervals and periodicity specified with the `DATE` command.
- The procedure verifies that date variables that are implied by the current date specification exist and that their values coincide with the date specification within the estimation period. If `SPLIT FILE` is in effect, the check is done for each split. A split is skipped and a warning is issued if verification fails for that split. If the verification fails when `SPLIT FILE` is not in effect, an error occurs and processing terminates. If the data are filtered, values of date variables are verified within the filtered subset of data.

- The procedure honors the estimation (historical) period that is defined with the `USE` command. However, depending on available data, the actual use period may vary by dependent variable. For a given dependent variable, the estimation period is the period left after eliminating contiguous missing values of the dependent variable at the beginning and end of the `USE` period.
- If `SPLIT FILE` is in effect, and a split has fewer cases than implied by the `USE` period, a warning is issued and all available cases are used. If there are no cases in the `USE` period, the split is skipped altogether with a warning.
- The `PREDICT` command defines the end of the forecast period for tables, charts, and model variables that are produced by `TSMODEL`. The forecast period always starts after the end of the estimation (`USE`) period. Like the estimation period, the forecast period can vary by model, depending on available data. If a `PREDICT` specification ends on or before the effective end of the `USE` period, the forecast period is empty.
- The procedure does not honor the following commands: `MODEL NAME` or `SAVE MODEL`. Options for naming and saving models are provided with the `TSMODEL` procedure.
- The procedure does not honor `TSET`. The `TSMODEL` procedure provides options for handling missing values, setting the width of confidence intervals, setting the maximum number of lags displayed for autocorrelations, and setting season length.
- The `TDISPLAY` command does not display models that are saved by using `TSMODEL`.

### **Limitations**

- `WEIGHT` is ignored with a warning.

## **Examples**

This section provides simple examples that are designed to get you started with using the Expert Modeler, producing forecasts, saving results to the active dataset, and saving your model for later use. Further examples that are specific to each subcommand are provided in the subcommand topics.

### **Using the Expert Modeler**

```
TSMODEL
  /MODEL DEPENDENT=sku1 TO sku10.
```

- The Expert Modeler is used to find the best-fitting exponential smoothing or ARIMA model for each of the dependent series *sku1* thru *sku10*.
- The procedure invokes the Expert Modeler because the `MODEL` subcommand is not followed by one of the model type subcommands (i.e., `EXSMOOTH` or `ARIMA`). The absence of a model type subcommand is equivalent to specifying `/EXPERTMODELER TYPE=[ARIMA EXSMOOTH] TRYSEASONAL=YES`.

### **Obtaining Model Forecasts**

```
PREDICT THRU YEAR 2006 MONTH 6.
TSMODEL
  /SERIESPLOT FORECAST
```

```
/MODELDETAILS PRINT=FORECASTS
/MODEL DEPENDENT=revenue.
```

- The PREDICT command is used to specify the forecast period for the TSMODEL procedure.
- The SERIESPLOT subcommand specifies that the output contains a plot of the predicted values within the forecast period.
- The MODELDETAILS subcommand specifies that the output includes a table containing the predicted values within the forecast period.

### ***Saving Models to an External File***

```
TSMODEL
/MODEL DEPENDENT=sku1 TO sku50 OUTFILE='/models/models_sku1TOsku50.xml'.
```

- The OUTFILE keyword specifies that each of the resulting models is to be saved to the file */models/models\_sku1TOsku50.xml*.
- The saved models can be used to produce new forecasts with the TSAPPLY command when new data are available. [For more information, see the topic TSAPPLY on p. 2006.](#)

### ***Saving Model Predictions, Residuals, and Confidence Intervals as New Variables***

```
TSMODEL
/SAVE PREDICTED LCL UCL NRESIDUAL
/MODEL DEPENDENT=revenue.
```

- The SAVE subcommand specifies that new variables containing the model predictions, noise residuals, and confidence intervals are saved to the active dataset.

### ***Specifying Multiple Model Types***

```
TSMODEL
/MODEL DEPENDENT=sku1 TO sku10 INDEPENDENT=adspending
/EXPERTMODELER TYPE=[ARIMA]
/MODEL DEPENDENT=sku11 TO sku15
/EXSMOOTH TYPE=WINTERSMULTIPLICATIVE.
```

- The first MODEL block specifies that the Expert Modeler is used to find the best-fitting ARIMA model for each of the dependent series *sku1* thru *sku10*, using the predictor variable *adspending*.
- The second MODEL block specifies that the Winters' multiplicative method is used to model each of the dependent series *sku11* thru *sku15*.
- In this example, different model types were used for different dependent variables. You can also specify multiple model types for the same dependent variable, thereby obtaining multiple models for the variable.

## Goodness-of-Fit Measures

The following model goodness-of-fit measures are available in TSMODEL:

<b>SRSQUARE</b>	<b>Stationary R-squared.</b> A measure that compares the stationary part of the model to a simple mean model. This measure is preferable to ordinary R-squared when there is a trend or seasonal pattern. Stationary R-squared can be negative with a range of negative infinity to 1. Negative values mean that the model under consideration is worse than the baseline model. Positive values mean that the model under consideration is better than the baseline model.
<b>RSQUARE</b>	<b>R-squared.</b> An estimate of the proportion of the total variation in the series that is explained by the model. This measure is most useful when the series is stationary. R-squared can be negative with a range of negative infinity to 1. Negative values mean that the model under consideration is worse than the baseline model. Positive values mean that the model under consideration is better than the baseline model.
<b>RMSE</b>	<b>RMSE.</b> Root Mean Square Error. The square root of mean square error. A measure of how much a dependent series varies from its model-predicted level, expressed in the same units as the dependent series.
<b>MAPE</b>	<b>MAPE.</b> Mean Absolute Percentage Error. A measure of how much a dependent series varies from its model-predicted level. It is independent of the units used and can therefore be used to compare series with different units.
<b>MAE</b>	<b>MAE.</b> Mean absolute error. Measures how much the series varies from its model-predicted level. MAE is reported in the original series units.
<b>MAXAPE</b>	<b>MaxAPE.</b> Maximum Absolute Percentage Error. The largest forecasted error, expressed as a percentage. This measure is useful for imagining a worst-case scenario for your forecasts.
<b>MAXAE</b>	<b>MaxAE.</b> Maximum Absolute Error. The largest forecasted error, expressed in the same units as the dependent series. Like MaxAPE, it is useful for imagining the worst-case scenario for your forecasts. Maximum absolute error and maximum absolute percentage error may occur at different series points—for example, when the absolute error for a large series value is slightly larger than the absolute error for a small series value. In that case, the maximum absolute error will occur at the larger series value and the maximum absolute percentage error will occur at the smaller series value.
<b>NORMBIC</b>	<b>Normalized BIC.</b> Normalized Bayesian Information Criterion. A general measure of the overall fit of a model that attempts to account for model complexity. It is a score based upon the mean square error and includes a penalty for the number of parameters in the model and the length of the series. The penalty removes the advantage of models with more parameters, making the statistic easy to compare across different models for the same series.

## MODELSUMMARY Subcommand

The MODELSUMMARY subcommand controls the display of tables and charts that summarize goodness of fit, residual autocorrelations, and residual partial autocorrelations across estimated models. Each keyword is followed by an equals sign (=) and one or more of the available options enclosed in square brackets.

### Example

```
TSMODEL
  /MODELSUMMARY PLOT=[SRSQUARE MAXAPE]
```

```
/MODEL DEPENDENT=sku1 TO sku100.
```

- The output includes two histograms: one histogram for stationary R-squared and one histogram for the maximum absolute percentage error. Each histogram consists of results across all models (one model for each dependent variable).

### **PRINT Keyword**

The PRINT keyword controls the display of model summary tables.

<b>MODELFIT</b>	<i>Goodness of fit.</i> Table of summary statistics and percentiles for stationary R-square, R-square, root mean square error, mean absolute percentage error, mean absolute error, maximum absolute percentage error, maximum absolute error, and normalized Bayesian Information Criterion.
<b>RESIDACF</b>	<i>Residual autocorrelation function.</i> Table of summary statistics and percentiles for autocorrelations of the residuals.
<b>RESIDPACF</b>	<i>Residual partial autocorrelation function.</i> Table of summary statistics and percentiles for partial autocorrelations of the residuals.
<b>NONE</b>	<i>No tables are shown.</i> An error occurs if NONE is used in combination with any other PRINT option.

### **PLOT Keyword**

The PLOT keyword controls the display of model summary charts. By default, no charts are shown. (For detailed definitions of the following terms, see [Goodness-of-Fit Measures](#) on p. 2030.)

<b>SRSQUARE</b>	<i>Histogram of Stationary R-Square.</i>
<b>RSQUARE</b>	<i>Histogram of R-Square.</i>
<b>RMSE</b>	<i>Histogram of Root Mean Square Error.</i>
<b>MAPE</b>	<i>Histogram of Mean Absolute Percentage Error.</i>
<b>MAE</b>	<i>Histogram of Mean Absolute Error.</i>
<b>MAXAPE</b>	<i>Histogram of Maximum Absolute Percentage Error.</i>
<b>MAXAE</b>	<i>Histogram of Maximum Absolute Error.</i>
<b>NORMBIC</b>	<i>Histogram of Normalized Bayesian Information Criterion (BIC).</i>
<b>RESIDACF</b>	<i>Boxplot of Residual Autocorrelation Function by Lag.</i>
<b>RESIDPACF</b>	<i>Boxplot of Residual Partial Autocorrelation Function by Lag.</i>

## **MODELSTATISTICS Subcommand**

The MODELSTATISTICS subcommand controls display of a table that lists all models, along with chosen goodness-of-fit statistics.

### **Example**

```
TSMODEL
/MODELSTATISTICS DISPLAY=YES MODELFIT=[RSQUARE]
/MODEL DEPENDENT=sku1 TO sku25.
```

- The output includes a table displaying the value of R-squared for each model.

**DISPLAY Keyword**

The DISPLAY keyword controls whether the model statistics table is shown.

<b>YES</b>	<i>Model statistics table is shown.</i> Table of model goodness of fit, Ljung-Box Q statistic, and number of outliers detected by model. This setting is the default. The Q statistic measures the degree of pattern in the residuals. Large values of Q in relation to its degrees of freedom indicate that model residuals are not randomly distributed.
<b>NO</b>	<i>Model statistics table is not shown.</i>

**MODELFIT Keyword**

The MODELFIT keyword controls which fit statistics are shown in the model statistics table, and the keyword is ignored if DISPLAY=NO. The keyword is followed by an equals sign (=) and one or more of the following options enclosed in square brackets. (For detailed definitions of the following terms, see [Goodness-of-Fit Measures](#) on p. 2030.)

<b>SRSQUARE</b>	<i>Stationary R -Square.</i> This setting is the default.
<b>RSQUARE</b>	<i>R-Square.</i>
<b>RMSE</b>	<i>Root Mean Square Error.</i>
<b>MAPE</b>	<i>Mean Absolute Percentage Error.</i>
<b>MAE</b>	<i>Mean Absolute Error.</i>
<b>MAXAPE</b>	<i>Maximum Absolute Percentage Error.</i>
<b>MAXAE</b>	<i>Maximum Absolute Error.</i>
<b>NORMBIC</b>	<i>Normalized Bayesian Information Criterion (BIC).</i>

**MODELDETAILS Subcommand**

The MODELDETAILS subcommand controls the display of model parameters, model forecasts, and autocorrelations of noise residuals for individual models. Each keyword is followed by an equals sign (=) and one or more of the available options enclosed in square brackets.

**Example**

```
TSMODEL
  /MODELDETAILS PRINT=[PARAMETERS FORECASTS]
  /MODEL DEPENDENT=sku1 TO sku50.
```

- The output includes tables displaying the model parameters and forecasts for each model (one model for each dependent variable).

**PRINT Keyword**

The PRINT keyword controls display of tables for individual models. By default, no tables are shown.

<b>PARAMETERS</b>	<i>Model parameter estimates.</i> Shows separate tables for ARIMA and exponential smoothing models. If outliers exist, parameter estimates for outliers are also displayed.
<b>RESIDACF</b>	<i>Residual autocorrelation function.</i> Shows residual autocorrelations by lag.
<b>RESIDPACF</b>	<i>Residual partial autocorrelation function.</i> Shows residual partial autocorrelations by lag.
<b>FORECASTS</b>	<i>Forecasts and confidence intervals.</i> Shows model forecasts and confidence intervals.

**PLOT Keyword**

The PLOT keyword controls display of charts for individual models. By default, no charts are shown.

<b>RESIDACF</b>	<i>Residual autocorrelation function.</i> Shows residual autocorrelations by lag.
<b>RESIDPACF</b>	<i>Residual partial autocorrelation function.</i> Shows residual partial autocorrelations by lag.

**SERIESPLOT Subcommand**

The SERIESPLOT subcommand allows you to obtain plots of predicted values, observed values, and confidence intervals for each estimated model. By default, no plots are shown. The subcommand is followed by one or more of the following keywords:

<b>OBSERVED</b>	<i>Displays all observed values of the dependent series.</i>
<b>FORECAST</b>	<i>Displays model-predicted values within the forecast period.</i>
<b>FIT</b>	<i>Displays model-predicted values within the estimation period.</i>
<b>FORECASTCI</b>	<i>Displays upper and lower confidence limits within the forecast period.</i>
<b>FITCI</b>	<i>Displays upper and lower confidence limits within the estimation period.</i>

**Example**

```
TSMODEL
  /SERIESPLOT OBSERVED FORECAST
  /MODEL DEPENDENT=sku1 TO sku5
  /MODEL DEPENDENT=revenue INDEPENDENT=adspending.
```

- The plot includes the model-predicted values in the forecast period as well as all observed values for each of the dependent series.

## **OUTPUTFILTER Subcommand**

The OUTPUTFILTER subcommand controls which models are included in the model statistics table (MODELSTATISTICS subcommand), detailed model output (MODELDETAILS subcommand), and time series plots (SERIESPLOT subcommand).

- By default, all models are included in the output. OUTPUTFILTER can be used to display only the best-fitting or worst-fitting models (or both). If you request both, two sets of output are displayed.
- OUTPUTFILTER is ignored if no output is requested on the MODELSTATISTICS, MODELDETAILS, or SERIESPLOT subcommands.
- OUTPUTFILTER has no effect on output from the MODELSUMMARY subcommand.

### **Example**

```
TSMODEL
/MODELSTATISTICS DISPLAY=YES MODELFIT=[SRSQUARE MAXAPE NORMBIC]
/OUTPUTFILTER DISPLAY=[BESTFIT(N=5) WORSTFIT(PCT=10)]
/MODEL DEPENDENT=sku1 TO sku200.
```

- The output consists of two sets of results: those results for the 5 best-fitting models and those results for models with fit values in the bottom 10%. The stationary R-squared value is used as the goodness-of-fit measure.

### **DISPLAY Keyword**

The DISPLAY keyword controls which models are included in MODELSTATISTICS, MODELDETAILS, and SERIESPLOT output.

- By default, all models are displayed (DISPLAY=ALLMODELS).
- To display only those models with extreme goodness-of-fit values, specify BESTFIT, WORSTFIT, or both, in square brackets. One or both keywords must be specified.

<b>ALLMODELS</b>	<i>All models are displayed.</i> All models that could be computed are shown. This setting is the default.
<b>BESTFIT</b>	<i>Models having the highest fit values are shown.</i> To display the models with the N highest fit values, specify BESTFIT followed by N=count in parentheses—for example, BESTFIT(N=5). The count must be a positive integer. If the count exceeds the number of estimated models, all models are shown. To display the models with fit values in the top N%, specify PCT=percent in parentheses. For example, BESTFIT(PCT=5) displays models with fit values in the top 5% among estimated models. The percentage value must be greater than zero and less than 100.
<b>WORSTFIT</b>	<i>Models having the lowest fit values are shown.</i> To display the models with the N lowest fit values, specify WORSTFIT followed by N=count in parentheses—for example, WORSTFIT(N=5). The count must be a positive integer. If the count exceeds the number of estimated models, all models are shown. To display the models with fit values in the lowest N%, specify PCT=percent in parentheses. For example, WORSTFIT(PCT=5) displays models with fit values in the bottom 5% among estimated models. The percentage value must be greater than zero and less than 100.



**MODELFIT Keyword**

The `MODELFIT` keyword specifies the fit measure that is used to filter models. `MODELFIT` is ignored if `DISPLAY=ALLMODELS`.

Specify one of the following options. (For detailed definitions of the following terms, see [Goodness-of-Fit Measures](#) on p. 2030.)

<b>SRSQUARE</b>	<i>Stationary R-Square.</i> This setting is the default.
<b>RSQUARE</b>	<i>R-Square.</i>
<b>RMSE</b>	<i>Root Mean Square Error.</i>
<b>MAPE</b>	<i>Mean Absolute Percentage Error.</i>
<b>MAE</b>	<i>Mean Absolute Error.</i>
<b>MAXAPE</b>	<i>Maximum Absolute Percentage Error.</i>
<b>MAXAE</b>	<i>Maximum Absolute Error.</i>
<b>NORMBIC</b>	<i>Normalized Bayesian Information Criterion (BIC).</i>

**SAVE Subcommand**

The `SAVE` subcommand is used to save new variables representing predicted values, residuals, and confidence intervals to the active dataset. By default, no new variables are saved to the active dataset.

- Specify one or more keywords, each keyword followed by an optional rootname to be used as the prefix for new variable names. Enclose the rootname in parentheses. Each keyword gives rise to one new variable for each dependent variable.
- The rootname, if specified, must conform to the rules for valid variable names.
- If no rootname is specified, `TSMODEL` uses a default name.
- The full variable name is the concatenation of the rootname, the name of the associated dependent variable, and a model identifier. The variable name is extended if necessary to avoid variable naming conflicts.
- The `PREDICT` command controls whether new cases are added to the file when new variables are saved. New cases are added if `PREDICT` specifies a forecast period that extends beyond the length of the dependent-variable series.
- `SAVE` is ignored with a warning if temporary transformations are in effect.

**Example**

```
TSMODEL
  /SAVE PREDICTED(Pred) NRESIDUAL
  /MODEL DEPENDENT=skul TO sku10.
```

- Two new variables are created for each of the dependent variables `skul` thru `sku10`: one variable contains the model predictions, and the other variable contains the noise residuals.

- The rootname (prefix) for the variables containing model predictions is *Pred*.
- The rootname for the variables containing noise residuals is *NResidual* (the default).

<b>PREDICTED(rootname)</b>	<i>Model fit and forecast values.</i> The default rootname is <i>Predicted</i> .
<b>LCL(rootname)</b>	<i>Lower confidence limits.</i> The default rootname is <i>LCL</i> .
<b>UCL(rootname)</b>	<i>Upper confidence limits.</i> The default rootname is <i>UCL</i> .
<b>NRESIDUAL(rootname)</b>	<i>Noise residuals.</i> The default rootname is <i>NResidual</i> . When transformations of the dependent variable are performed (for example, natural log), these residuals are the residuals for the transformed series.

## **AUXILIARY Subcommand**

The **AUXILIARY** subcommand is used to set the confidence interval level for forecasts, set the maximum number of lags for autocorrelation and partial autocorrelation plots and tables, and set the season length.

### **Example**

```
TSMODEL
  /MODELDETAILS PLOT=[RESIDACF RESIDPACF]
  /AUXILIARY MAXACFLAGS=36 SEASONLENGTH=6
  /MODEL DEPENDENT=sku1 TO sku20.
```

### **CILEVEL Keyword**

The **CILEVEL** keyword sets the confidence level.

- Specify a positive number that is less than 100.
- The default is 95.

### **MAXACFLAGS Keyword**

The **MAXACFLAGS** keyword sets the maximum number of lags displayed in residual autocorrelation and partial autocorrelation tables and plots.

- Specify a positive integer.
- The default is 24.

### **SEASONLENGTH Keyword**

The **SEASONLENGTH** keyword is used to specify the length of the seasonal period (the number of observations in one period or season).

- The season length must be a positive integer.
- If **SEASONLENGTH** is not specified, the periodicity that is specified by using the **DATE** command defines the season length.

- If the Expert Modeler is used, a warning occurs and the season length is treated as 1 for any splits for which there are gaps in the data. If a custom model is specified, a warning is issued and the model is not estimated for any such splits.
- For custom ARIMA or exponential smoothing models, SEASONLENGTH is ignored if the specified model has no seasonal component.

## **MISSING Subcommand**

The `MISSING` subcommand controls the handling of user-missing values.

- By default, user-missing values are treated as missing (invalid) data.
- System-missing values are always treated as invalid.
- Cases with missing values of a dependent variable that occur within the estimation period are included in the model. The specific handling of the missing value depends on the estimation method.
- A warning is issued if a predictor has missing values within the estimation period. For the Expert Modeler, models involving the predictor are estimated without the predictor. For custom ARIMA, models involving the predictor are not estimated.
- If any predictor has missing values within the forecast period, the procedure issues a warning and forecasts as far as it can.

### **Example**

```
TSMODEL
  /MISSING USERMISSING=INCLUDE
  /MODEL DEPENDENT=revenue INDEPENDENT=adspending.
```

### **USERMISSING Keyword**

The `USERMISSING` keyword controls the treatment of user-missing values and is required.

- |                |  |
|----------------|--|
| <b>EXCLUDE</b> | <i>Exclude user-missing values.</i> User-missing values are treated as missing. This setting is the default. |
| <b>INCLUDE</b> | <i>Include user-missing values.</i> User-missing values are treated as valid data.                           |

## **MODEL Subcommand**

The `MODEL` subcommand is required and is used to signal the start of a model specification, as well as to specify model variable(s).

- All variables must be numeric. Any string variables are filtered out, with a warning.
- Variables that define split-file groups are filtered out of all variable lists, with a warning.
- `ALL` and `TO` can be used in dependent and independent variable lists. `ALL` represents all variables in the active dataset. `TO` refers to a range of variables in the active dataset.

*Note:* Specification of the model type is optional and is accomplished by including an `EXPERTMODELER`, `EXSMOOTH`, or `ARIMA` subcommand following the `MODEL` subcommand. If the model type is not specified, the best-fitting ARIMA or exponential smoothing model

will automatically be selected by the Expert Modeler, which is equivalent to specifying `/EXPERTMODELER TYPE=[ARIMA EXSMOOTH]`. If you are unsure of which type of model to choose, or you want Expert Modeler to choose for you, use the `MODEL` subcommand without a model type subcommand.

### **Example**

```
TSMODEL
  /MODEL DEPENDENT=store1 TO store100
        INDEPENDENT=adspending
        OUTFILE='/models/stores.xml'
        PREFIX='Expert'.
```

- `DEPENDENT` specifies that variables *store1* thru *store100* are to be modeled. The keyword `TO` refers to the order of the variables in the active dataset.
- The `INDEPENDENT` keyword specifies that the variable *adspending* is to be considered as a predictor variable.
- `OUTFILE` specifies that the resulting models are to be stored in the file */models/stores.xml*.
- `PREFIX` specifies that the models will be named *Expert\_1* (the model for *store1*), *Expert\_2* (the model for *store2*), etc.
- The absence of a model type subcommand (`EXPERTMODELER`, `ARIMA`, or `EXSMOOTH`) means that the Expert Modeler will be used to find the best-fitting exponential smoothing or ARIMA model.

### **DEPENDENT Keyword**

The `DEPENDENT` keyword is used to specify one or more dependent variables and is required.

- At least one dependent variable must be specified.
- Repeated instances of the same variable are filtered out of the list.
- A separate model is generated for each dependent variable in each `MODEL` block. Identical dependent variables in separate `MODEL` blocks generate separate models.

### **INDEPENDENT Keyword**

The `INDEPENDENT` keyword is used to specify one or more optional independent variables.

- Order of variables within the independent variable list matters when the Expert Modeler is used. The Expert Modeler drops nonsignificant independent variables one at a time, starting with the last variable in the list.
- In the case of custom ARIMA models, all independent variables are included in the model.
- Repeated instances of an independent variable are filtered out of the list. For example, `a b a c a` is equivalent to `a b c`. Assuming that the active dataset contains variables *a*, *b*, and *c*, `ALL` is also equivalent to `a b c`.
- An independent variable that also appears in the dependent variable list is treated solely as a dependent variable and is excluded from consideration as an independent variable for the model involving that particular dependent variable. For example, if you specify

DEPENDENT=*a b* and INDEPENDENT=*a c*, then *c* is the only independent variable when modeling *a*, but *a* and *c* are both used as independent variables when modeling *b*.

### **Events**

- Event variables are special independent variables that are used to model effects of occurrences such as a flood, strike, or introduction of a new product line. Any abrupt shift in the level of the dependent series can be modeled by using event variables.
- To designate an independent variable as an event variable, specify [E] following the name of the variable. For example, INDEPENDENT=*strike* [E] indicates that the variable *strike* is an event variable.
- When the Expert Modeler is used, event variables enter the model with linear terms only (as opposed to transfer functions). Thus, event designation can save processing time and guard against overfitting when using the Expert Modeler.
- Custom ARIMA models ignore event designation.
- For event variables, the presence of an effect is indicated by cases with a value of 1. Cases with values other than 1 are treated as indicating no effect.
- The effect of an event can last a single time period or several periods. If the effect lasts several periods, the corresponding event variable should contain a series of consecutive 1s.
- Event designation applies only to the variable that immediately precedes the designation in the independent variable list. For example, *x1 x2 x3* [E] and *x1 TO x3* [E] designate *x3* (only) as an event variable. If event designation follows ALL, it applies to all independent variables.
- If a variable appears more than once in the independent variable list, the event specification for the last instance of the variable is honored. For example, if ALL *x1* [E] is specified, *x1* is treated as an event variable; if ALL [E] *x1* is specified, *x1* is treated as an ordinary predictor.

### **OUTFILE Keyword**

The OUTFILE keyword is used to save models to an external file. Saved models can be used to obtain updated forecasts, based on more current data, using the TSAPPLY command. [For more information, see the topic TSAPPLY on p. 2006.](#)

- Models are written to an XML file. Each model is assigned a unique name and a description that includes the name of the dependent variable and the model type.
- The filename must be specified in full. No extension is supplied.
- If two or more MODEL subcommands (within a single invocation of TSMODEL) specify the same external file, all models that are created by those MODEL subcommands are saved to that file.
- OUTFILE is ignored with a warning if SPLIT FILE is in effect.

### **PREFIX Keyword**

Models are assigned unique names consisting of a customizable prefix, along with an integer suffix. The PREFIX keyword is used to specify a custom prefix.

- Custom prefixes must be specified in quotes—for example, `PREFIX='ExpertArima'`. The default prefix is 'Model'.
- Integer suffixes are unique across the set of models that have the same prefix; for example, `Model_1`, `Model_2`. Models with different prefixes, however, can have the same integer suffix—for example, `CustomArima_1`, `ExpertArima_1`.

## **EXPERTMODELER Subcommand**

The `EXPERTMODELER` subcommand controls options for automatic model identification that is performed by the Expert Modeler.

### **Example**

```
TSMODEL
  /MODEL DEPENDENT=store1 TO store100
        INDEPENDENT=adspending
  /EXPERTMODELER TYPE=[ARIMA].
```

- The keyword `ARIMA` specifies that the Expert Modeler is to limit the model search to the best-fitting ARIMA model.

### **TYPE Keyword**

The `TYPE` keyword is required and is used to specify the model types that are evaluated: ARIMA, exponential smoothing, or both.

- By default, both ARIMA and exponential smoothing models are considered.
- Either `ARIMA`, `EXSMOOTH`, or both must be specified in square brackets.
- An error occurs if you specify the `EXSMOOTH` keyword by itself and the model contains independent variables.

The keyword is followed by an equals sign (=) and one or both of the following options enclosed in square brackets:

<b>ARIMA</b>	<i>ARIMA models are considered.</i>
<b>EXSMOOTH</b>	<i>Exponential smoothing models are considered.</i>

### **TRYSEASONAL Keyword**

The `TRYSEASONAL` keyword specifies whether seasonal models are considered by the Expert Modeler.

- `TRYSEASONAL` is ignored if the data are nonperiodic (that is, season length is 1).

<b>YES</b>	<i>The Expert Modeler considers seasonal models. This setting is the default.</i>
<b>NO</b>	<i>The Expert Modeler does not consider seasonal models. Only nonseasonal models are considered.</i>

## EXSMOOTH Subcommand

The EXSMOOTH subcommand controls options for custom exponential smoothing models.

- Using EXSMOOTH generates an error if independent variables are defined.
- For seasonal models (SIMPLESEASONAL, WINTERSADDITIVE, and WINTERSMULTIPLICATIVE), an error is generated if the season length has not been defined via the DATE command or the SEASONLENGTH keyword.

### Example

```
TSMODEL
/AUXILIARY SEASONLENGTH=12
/MODEL DEPENDENT=store1 TO store100
/EXSMOOTH TYPE=WINTERSMULTIPLICATIVE.
```

### TYPE Keyword

The TYPE keyword is required and is used to specify the type of exponential smoothing model. There is no default method.

<b>SIMPLE</b>	<b>Simple exponential smoothing model.</b> This model is appropriate for series in which there is no trend or seasonality. Its only smoothing parameter is level. Simple exponential smoothing is most similar to an ARIMA model with zero orders of autoregression, one order of differencing, one order of moving average, and no constant.
<b>SIMPLESEASONAL</b>	<b>Simple seasonal exponential smoothing model.</b> This model is appropriate for series with no trend and a seasonal effect that is constant over time. Its smoothing parameters are level and season. Simple seasonal exponential smoothing is most similar to an ARIMA model with zero orders of autoregression, one order of differencing, one order of seasonal differencing, and orders 1, p, and p + 1 of moving average, where p is the number of periods in a seasonal interval (for monthly data, p = 12).
<b>HOLT</b>	<b>Holt's method.</b> This model is appropriate for series in which there is a linear trend and no seasonality. Its smoothing parameters are level and trend, which are not constrained by each other's values. Holt's model is more general than Brown's model but may take longer to compute for large series. Holt's exponential smoothing is most similar to an ARIMA model with zero orders of autoregression, two orders of differencing, and two orders of moving average.
<b>BROWN</b>	<b>Brown's method.</b> This model is appropriate for series in which there is a linear trend and no seasonality. Its smoothing parameters are level and trend, which are assumed to be equal. Brown's model is therefore a special case of Holt's model. Brown's exponential smoothing is most similar to an ARIMA model with zero orders of autoregression, two orders of differencing, and two orders of moving average, with the coefficient for the second order of moving average equal to the square of one-half of the coefficient for the first order.
<b>DAMPEDTREND</b>	<b>Damped trend method.</b> This model is appropriate for series with a linear trend that is dying out and with no seasonality. Its smoothing parameters are level, trend, and damping trend. Damped exponential smoothing is most similar to an ARIMA model with 1 order of autoregression, 1 order of differencing, and 2 orders of moving average.

**WINTERSADDITIVE**

**Winters' additive method.** This model is appropriate for series with a linear trend and a seasonal effect that does not depend on the level of the series. Its smoothing parameters are level, trend, and season. Winters' additive exponential smoothing is most similar to an ARIMA model with zero orders of autoregression, one order of differencing, one order of seasonal differencing, and  $p + 1$  orders of moving average, where  $p$  is the number of periods in a seasonal interval (for monthly data,  $p = 12$ ).

**WINTERSMULTIPLICATIVE**

**Winters' multiplicative method.** This model is appropriate for series with a linear trend and a seasonal effect that depends on the level of the series. Its smoothing parameters are level, trend, and season. Winters' multiplicative exponential smoothing is not similar to any ARIMA model.

**TRANSFORM Keyword**

The TRANSFORM keyword specifies a transformation that is performed on each dependent variable before it is modeled. By default, dependent variables are not transformed.

<b>NONE</b>	<i>Dependent variable series are not transformed.</i> This setting is the default.
<b>SQRT</b>	<i>A square root transformation is performed.</i>
<b>LN</b>	<i>A natural log transformation is performed.</i>

**ARIMA Subcommand**

The ARIMA subcommand controls options for custom ARIMA models. The model must contain at least one parameter. An error occurs if the model has no autoregressive component, moving average component, outlier, or constant term, unless independent variables are selected.

**Example: Basic Custom Seasonal ARIMA Model**

```
TSMODEL
  /AUXILIARY SEASONLENGTH=12
  /MODEL DEPENDENT=passengers
  /ARIMA AR=[0] ARSEASONAL=[0] MA=[1] MASEASONAL=[1]
        DIFF=1 DIFFSEASONAL=1 TRANSFORM=LN.
```

- A custom ARIMA(0,1,1)(0,1,1) model is specified, including a natural log transformation of the dependent variable. The seasonal length is 12.

**Example: Including Specific Lags in a Custom ARIMA Model**

```
TSMODEL
  /MODEL DEPENDENT=revenue
  /ARIMA AR=0 MA=[1 3].
```

- A moving average model is specified. Moving average lags of orders 1 and 3 are included.

**AR Keyword**

The AR keyword specifies the nonseasonal autoregressive order of the model and is required.



- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.

### ***ARSEASONAL Keyword***

The `ARSEASONAL` keyword specifies the seasonal autoregressive order of the model.

- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers. The default order is zero.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.
- An error occurs if `ARSEASONAL` specifies a nonzero value and the season length has not been defined via the `DATE` command or the `SEASONLENGTH` keyword.

### ***MA Keyword***

The `MA` keyword specifies the nonseasonal moving average order of the model and is required.

- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.

### ***MASEASONAL Keyword***

The `MASEASONAL` keyword specifies the seasonal moving average order of the model.

- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers. The default order is zero.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.
- An error occurs if `MASEASONAL` specifies a nonzero value and the season length has not been defined via the `DATE` command or the `SEASONLENGTH` keyword.

**CONSTANT Keyword**

The `CONSTANT` keyword controls whether the model for each dependent variable includes a constant term.

- YES**            *The model includes a constant term. This setting is the default.*  
**NO**             *No constant term is included in the model.*

**TRANSFORM Keyword**

The `TRANSFORM` keyword specifies a transformation that is performed on each dependent variable before it is modeled. By default, the variables are not transformed.

- NONE**           *Dependent-variable series are not transformed. This setting is the default.*  
**SQRT**           *A square root transformation is performed.*  
**LN**             *A natural log transformation is performed.*

**DIFF Keyword**

The `DIFF` keyword specifies the order of nonseasonal differencing.

- The order of nonseasonal differencing must be a nonnegative integer. The default order is zero.

**DIFFSEASONAL Keyword**

The `DIFFSEASONAL` keyword specifies the order of seasonal differencing.

- The order of seasonal differencing must be a nonnegative integer. The default order is zero.
- An error occurs if `DIFFSEASONAL` specifies a nonzero value and the season length has not been defined via the `DATE` command or the `SEASONLENGTH` keyword.

**TRANSFERFUNCTION Subcommand**

The `TRANSFERFUNCTION` subcommand specifies a transfer function and transformations for predictor variables in an ARIMA model.

- By default, for each predictor in an ARIMA model, the numerator and denominator orders of the transfer function (both seasonal and nonseasonal) are set to zero. The predictors are neither transformed nor differenced by default.
- The `TRANSFERFUNCTION` subcommand may be used more than once in a `MODEL` block to specify different transfer functions and/or transformations for different independent variables.
- If there are multiple `TRANSFERFUNCTION` subcommands that refer to the same predictor variable within a `MODEL` block, the last set of specifications is honored for that variable.
- The `TRANSFERFUNCTION` subcommand is ignored if the `EXPERTMODELER` or `EXSMOOTH` subcommand is used.

**Example**

TSMODEL

```

/MODEL DEPENDENT=revenue INDEPENDENT=adspending
/ARIMA AR=[1] MA=[0]
/TRANSFERFUNCTION VARIABLES=adspending NUM=[1] DENOM=[1].

```

- A custom ARIMA(1,0,0) model is specified and includes the independent variable *adspending*.
- The TRANSFERFUNCTION subcommand specifies that *adspending* is to be modeled with numerator and denominator lags of order 1.

### **VARIABLES Keyword**

The VARIABLES keyword is used to specify a list of predictors.

- Repeated instances of the same variable are filtered out of the list.
- Variables that have not been specified as independent (with the INDEPENDENT keyword) in the preceding MODEL subcommand are ignored.
- Use ALL to specify all independent variables. This setting is the default.

### **NUM Keyword**

The NUM keyword specifies the numerator order of the transfer function for nonseasonal models.

- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers. The default order is zero.
- The model always includes an implicit zero, even if not specified via the NUM keyword.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.

### **NUMSEASONAL Keyword**

The NUMSEASONAL keyword specifies the numerator order of the transfer function for seasonal models.

- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers. The default order is zero.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.
- An error occurs if NUMSEASONAL specifies a nonzero value and the season length has not been defined via the DATE command or the SEASONLENGTH keyword.

### **DENOM Keyword**

The DENOM keyword specifies the denominator order of the transfer function for nonseasonal models.

- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers. The default order is zero.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.

### ***DENOMSEASONAL Keyword***

The `DENOMSEASONAL` keyword specifies the denominator order of the transfer function for seasonal models.

- You can specify a single value or a list of values in square brackets. All values must be nonnegative integers. The default order is zero.
- If you specify a single value, square brackets are optional. The specified lag will be included in the model.
- If you specify a list of positive integer values—for example, [1 3]—the specified lags are included in the model. Duplicate lags are ignored.
- An error occurs if `DENOMSEASONAL` specifies a nonzero value and the season length has not been defined via the `DATE` command or the `SEASONLENGTH` keyword.

### ***TRANSFORM Keyword***

The `TRANSFORM` keyword specifies a transformation that is performed on all predictor variables. By default, predictor variables are not transformed.

<b>NONE</b>	<i>Predictor variables are not transformed. This setting is the default.</i>
<b>SQRT</b>	<i>A square root transformation is performed.</i>
<b>LN</b>	<i>A natural log transformation is performed.</i>

### ***DIFF Keyword***

The `DIFF` keyword specifies the order of nonseasonal differencing for predictor variables.

- The order of nonseasonal differencing must be a nonnegative integer. The default order is zero.

### ***DIFFSEASONAL Keyword***

The `DIFFSEASONAL` keyword specifies the order of seasonal differencing for predictor variables.

- The order of seasonal differencing must be a nonnegative integer. The default order is zero.
- `DIFFSEASONAL` generates an error if the season length has not been defined via the `DATE` command or the `SEASONLENGTH` keyword.

### ***DELAY Keyword***

The `DELAY` keyword specifies the delay order for predictor variables.

- The delay order must be a nonnegative integer. The default order is zero.

## AUTOOUTLIER Subcommand

The AUTOOUTLIER subcommand specifies whether outlier time points are identified automatically.

- By default, outliers are not identified.
- If identified, outliers are not discarded but are incorporated in the model.
- The AUTOOUTLIER subcommand is ignored with a warning if it is part of a MODEL block that contains an EXSMOOTH subcommand.

### Example

```
TSMODEL
/MODELDETAILS PRINT=PARAMETERS
/MODEL DEPENDENT=revenue
/EXPERTMODELER TYPE=[ARIMA]
/AUTOOUTLIER DETECT=ON.
```

### DETECT Keyword

The DETECT keyword controls whether automatic outlier detection is performed.

- OFF**        *Outliers are not identified.* This setting is the default.  
**ON**         *Outliers are identified.* ON is ignored with a warning if the model type is EXSMOOTH.

### TYPE Keyword

The optional TYPE keyword specifies the types of outliers that are identified when DETECT=ON. TYPE is ignored if DETECT=OFF.

- Additive and level shift outliers are identified by default.
- The specification of outlier types overrides the default. That is, if you specify a list of outlier types, additive and level shift outliers are not included unless explicitly requested.

The keyword is followed by an equals sign (=) and one or more of the following options enclosed in square brackets:

<b>ADDITIVE</b>	<b>Additive.</b> An outlier that affects a single observation. For example, a data coding error might be identified as an additive outlier. Default if DETECT=ON.
<b>LEVELSHIFT</b>	<b>Level shift.</b> An outlier that shifts all observations by a constant, starting at a particular series point. A level shift could result from a change in policy. Default if DETECT=ON.
<b>INNOVATIONAL</b>	<b>Innovational.</b> An outlier that acts as an addition to the noise term at a particular series point. For stationary series, an innovational outlier affects several observations. For nonstationary series, it may affect every observation starting at a particular series point.
<b>TRANSIENT</b>	<b>Transient.</b> An outlier whose impact decays exponentially to 0.
<b>SEASONALADDITIVE</b>	<b>Seasonal additive.</b> An outlier that affects a particular observation and all subsequent observations separated from it by one or more seasonal periods. All such observations are affected equally. A seasonal additive outlier might occur if, beginning in a certain year, sales are higher every January. The keyword generates an error if the season length has not been defined via the DATE command or the SEASONLENGTH keyword.

**LOCALTREND**  
**ADDITIVEPATCH**

**Local trend.** An outlier that starts a local trend at a particular series point.  
**Additive patch.** A group of two or more consecutive additive outliers. Selecting this outlier type results in the detection of individual additive outliers in addition to patches of them.

## ***OUTLIER Subcommand***

The **OUTLIER** subcommand specifies outlier time points to be included in an ARIMA model.

- By default, no time points are modeled as outliers.
- For each outlier, you must specify its time point (location) and type.
- Multiple **OUTLIER** subcommands can be used to define two or more outliers for an ARIMA model. Duplicate specifications (same location and type) are ignored and a warning is issued.
- A warning occurs and the **OUTLIER** subcommand is ignored if the subcommand is part of a **MODEL** block that contains an **EXPERTMODELER** or **EXSMOOTH** subcommand.
- If the **AUTOOUTLIER** and **OUTLIER** subcommands are both included in a **MODEL** block containing an **ARIMA** subcommand, a warning is issued and the **OUTLIER** specification takes precedence.

### ***Example***

```
TSMODEL
  /MODELDETAILS PRINT=PARAMETERS
  /MODEL DEPENDENT=revenue
  /ARIMA AR=[1] MA=[1]
  /OUTLIER LOCATION=[YEAR 2000 MONTH 8] TYPE=LOCALTREND.
```

- A custom ARIMA(1,0,1) model is specified with an outlier of type Local Trend for August, 2000.

### ***LOCATION Keyword***

The **LOCATION** keyword specifies the time point to be treated as an outlier and is required.

- Specify a date in square brackets. If the data are undated, you must specify a case number. Otherwise, specify the outlier location using date keywords and values. For example, **LOCATION=[YEAR 2000 MONTH 8]** specifies that August, 2000 be treated as an outlier. The following date keywords may be used: **CYCLE**, **YEAR**, **QUARTER**, **MONTH**, **WEEK**, **DAY**, **HOUR**, **MINUTE**, **SECOND**, and **OBS**. Every keyword in the IBM® SPSS® Statistics date specification must appear in **LOCATION**. Only keywords that correspond to system date variables may be used, and no date keyword may be used more than once. [For more information, see the topic DATE on p. 586.](#)
- A warning is issued if you specify an outlier location that is outside the estimation period or corresponds to a gap in the data. Any such models are ignored.

**TYPE Keyword**

The `TYPE` keyword specifies the type of outlier and is required. Specify one of the following types:

<b>ADDITIVE</b>	<b>Additive.</b> An outlier that affects a single observation. For example, a data coding error might be identified as an additive outlier.
<b>LEVELSHIFT</b>	<b>Level shift.</b> An outlier that shifts all observations by a constant, starting at a particular series point. A level shift could result from a change in policy.
<b>INNOVATIONAL</b>	<b>Innovational.</b> An outlier that acts as an addition to the noise term at a particular series point. For stationary series, an innovational outlier affects several observations. For nonstationary series, it may affect every observation starting at a particular series point.
<b>TRANSIENT</b>	<b>Transient.</b> An outlier whose impact decays exponentially to 0.
<b>SEASONALADDITIVE</b>	<b>Seasonal additive.</b> An outlier that affects a particular observation and all subsequent observations separated from it by one or more seasonal periods. All such observations are affected equally. A seasonal additive outlier might occur if, beginning in a certain year, sales are higher every January. The keyword generates an error if the season length has not been defined via the <code>DATE</code> command or the <code>SEASONLENGTH</code> keyword.
<b>LOCALTREND</b>	<b>Local trend.</b> An outlier that starts a local trend at a particular series point.

# TSPLIT

```
TSPLIT VARIABLES= variable names

[/DIFF={1}]
      {n}

[/SDIFF={1}]
      {n}

[/PERIOD=n]

[/{NOLOG**}]
      {LN      }

[/ID=varname]

[/MARK={varname} ]
      {date   }

[/SPLIT {UNIFORM**}]
      {SCALE  }

[/APPLY [= 'model name']]
```

*For plots with one variable:*

```
[/FORMAT=[{NOFILL**}] [{NOREFERENCE**}]]]
      {BOTTOM  }      {REFERENCE[(value)]}
```

*For plots with multiple variables:*

```
[/FORMAT={NOJOIN**}]
      {JOIN    }
      {HILO   }
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- For plots with one variable, REFERENCE keyword modified to allow specification of a value.

## **Example**

```
TSPLIT VARIABLES = TICKETS.
```

## **Overview**

TSPLIT produces a plot of one or more time series or sequence variables. You can request natural log and differencing transformations to produce plots of transformed variables. Several plot formats are available.



### **Options**

**Modifying the Variables.** You can request a natural log transformation of the variables using the `LN` subcommand and seasonal and nonseasonal differencing to any degree using the `SDIFF` and `DIFF` subcommands. With seasonal differencing, you can also specify the periodicity on the `PERIOD` subcommand.

**Plot Format.** With the `FORMAT` subcommand, you can fill in the space on one side of the plotted values on plots with one variable. You can also plot a reference line indicating the variable mean. For plots with two or more variables, you can specify whether you want to join the values for each observation with a vertical line. With the `ID` subcommand, you can label the horizontal axis with the values of a specified variable. You can mark the onset of an intervention variable on the plot with the `MARK` subcommand.

**Split-File Processing.** You can control how data that have been divided into subgroups by a `SPLIT FILE` command should be plotted using the `SPLIT` subcommand.

### **Subcommand Order**

- Subcommands can be specified in any order.

### **Syntax Rules**

- `VARIABLES` can be specified only once.
- Other subcommands can be specified more than once, but only the last specification of each one is executed.

### **Operations**

- Subcommand specifications apply to all variables named on the `TSPLLOT` command.
- If the `LN` subcommand is specified, any differencing requested on that `TSPLLOT` command is done on the log-transformed variables.
- Split-file information is displayed as part of the subtitle, and transformation information is displayed as part of the footnote.

### **Limitations**

- Maximum of one `VARIABLES` subcommand. There is no limit on the number of variables named on the list.

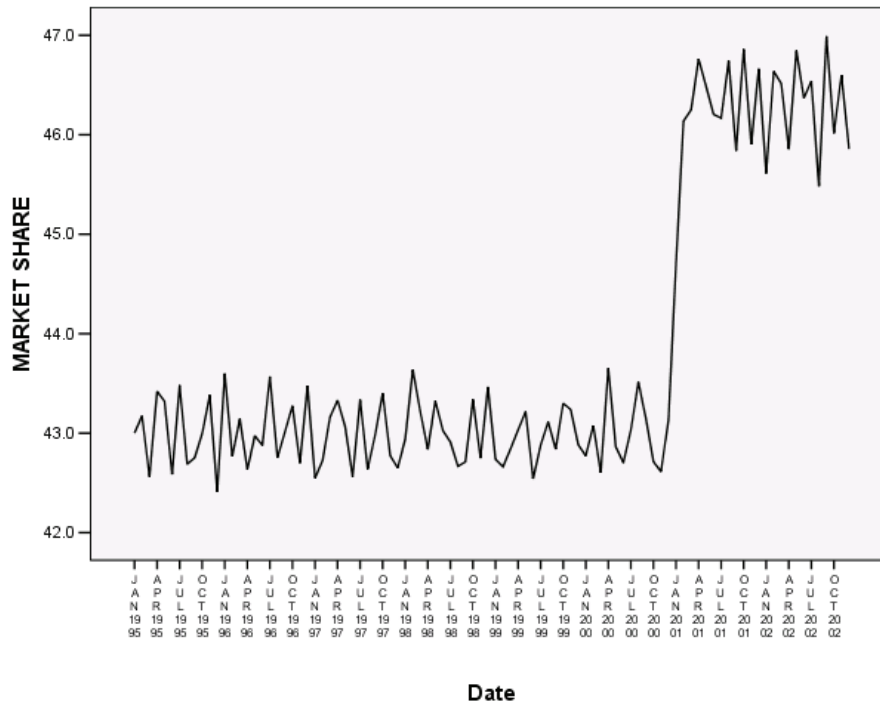
## **Basic Specification**

The basic specification is one or more variable names.

- If the `DATE` command has been specified, the horizontal axis is labeled with the `DATE_` variable at periodic intervals. Otherwise, sequence numbers are used. The vertical axis is labeled with the value scale of the plotted variable(s).

The following figure shows a default plot with `DATE YEAR 1995 MONTH`.

Figure 251-1  
 TSPLLOT VARIABLES=SHARE



## Example

```
TSPLLOT VARIABLES = TICKETS
/LN
/DIFF
/SDIFF
/PERIOD=12
/FORMAT=REFERENCE
/MARK=Y 55 M 6.
```

- This command produces a plot of *TICKETS* after a natural log transformation, differencing, and seasonal differencing have been applied.
- LN transforms the data using the natural logarithm (base  $e$ ) of *TICKETS*.
- DIFF differences the logged variable once.
- SDIFF and PERIOD apply one degree of seasonal differencing with a period of 12.
- FORMAT=REFERENCE adds a reference line representing the variable mean.
- MARK provides a marker on the plot at June 1955. The marker is displayed as a vertical reference line.

## VARIABLES Subcommand

VARIABLES specifies the names of the variables to be plotted and is the only required subcommand.

## ***DIFF Subcommand***

DIFF specifies the degree of differencing used to convert a nonstationary variable to a stationary one with a constant mean and variance before plotting.

- You can specify any positive integer on DIFF.
- If DIFF is specified without a value, the default is 1.
- The number of values plotted decreases by 1 for each degree of differencing.

### ***Example***

```
TSPLLOT VARIABLES = TICKETS  
/DIFF=2.
```

- In this example, *TICKETS* is differenced twice before plotting.

## ***SDIFF Subcommand***

If the variable exhibits a seasonal or periodic pattern, you can use the SDIFF subcommand to seasonally difference the variable before plotting.

- The specification on SDIFF indicates the degree of seasonal differencing and can be any positive integer.
- If SDIFF is specified without a value, the degree of seasonal differencing defaults to 1.
- The number of seasons plotted decreases by 1 for each degree of seasonal differencing.
- The length of the period used by SDIFF is specified on the PERIOD subcommand. If the PERIOD subcommand is not specified, the periodicity established on the TSET or DATE command is used (see the PERIOD subcommand).

## ***PERIOD Subcommand***

PERIOD indicates the length of the period to be used by the SDIFF subcommand.

- The specification on PERIOD indicates how many observations are in one period or season and can be any positive integer.
- If PERIOD is not specified, the periodicity established on TSET PERIOD is in effect. If TSET PERIOD is not specified, the periodicity established on the DATE command is used. If periodicity was not established anywhere, the SDIFF subcommand will not be executed.

### ***Example***

```
TSPLLOT VARIABLES = TICKETS  
/SDIFF=1  
/PERIOD=12.
```

- This command applies one degree of seasonal differencing with 12 observations per season to *TICKETS* before plotting.

## ***LN and NOLOG Subcommands***

LN transforms the data using the natural logarithm (base  $e$ ) of the variable and is used to remove varying amplitude over time. NOLOG indicates that the data should not be log transformed. NOLOG is the default.

- If you specify LN on TSPLLOT, any differencing requested on that command will be done on the log-transformed variables.
- There are no additional specifications on LN or NOLOG.
- Only the last LN or NOLOG subcommand on a TSPLLOT command is executed.
- If a natural log transformation is requested, any value less than or equal to zero is set to system-missing.
- NOLOG is generally used with an APPLY subcommand to turn off a previous LN specification.

### ***Example***

```
TSPLLOT VARIABLES = TICKETS  
/LN.
```

- In this example, *TICKETS* is transformed using the natural logarithm before plotting.

## ***ID Subcommand***

ID names a variable whose values will be used as labels for the horizontal axis.

- The only specification on ID is a variable name. If you have a variable named *ID* in your active dataset, the equals sign (=) after the subcommand is required.
- If the ID subcommand is not used and TSET ID has not been specified, the axis is labeled with the *DATE\_* variable created by the DATE command. If the DATE command has not been specified, the observation number is used as the label.

### ***Example***

```
TSPLLOT VARIABLES = VARA  
/ID=VARB.
```

- In this example, the values of *VARB* will be used to label the horizontal axis of *VARA* at periodic intervals.

## ***FORMAT Subcommand***

FORMAT controls the plot format.

- The specification on FORMAT is one of the keywords listed below.

- Keywords **NOFILL**, **BOTTOM**, **REFERENCE**, and **NOREFERENCE** apply to plots with one variable. **NOFILL** and **BOTTOM** are alternatives that indicate how the plot is filled. **NOREFERENCE** and **REFERENCE** are alternatives that specify whether a reference line is displayed.
- Keywords **JOIN**, **NOJOIN**, and **HILO** apply to plots with multiple variables and are alternatives. **NOJOIN** is the default. Only one keyword can be specified on a **FORMAT** subcommand for plots with multiple variables.

The following formats are available for plots with one variable:

<b>NOFILL</b>	<i>Plot only the values for the variable with no fill.</i> <b>NOFILL</b> produces a plot with no fill above or below the plotted values. This is the default format when one variable is specified.
<b>BOTTOM</b>	<i>Plot the values for the variable and fill in the area below the curve.</i> If the plotted variable has missing or negative values, <b>BOTTOM</b> is ignored and the default <b>NOFILL</b> is used instead.
<b>NOREFERENCE</b>	<i>Do not plot a reference line.</i> This is the default when one variable is specified.
<b>REFERENCE(value)</b>	<i>Plot a reference line at the specified value or at the variable mean if no value is specified.</i> A fill chart is displayed as an area chart with a reference line and a non-fill chart is displayed as a line chart with a reference line.

Figure 251-2  
FORMAT=BOTTOM

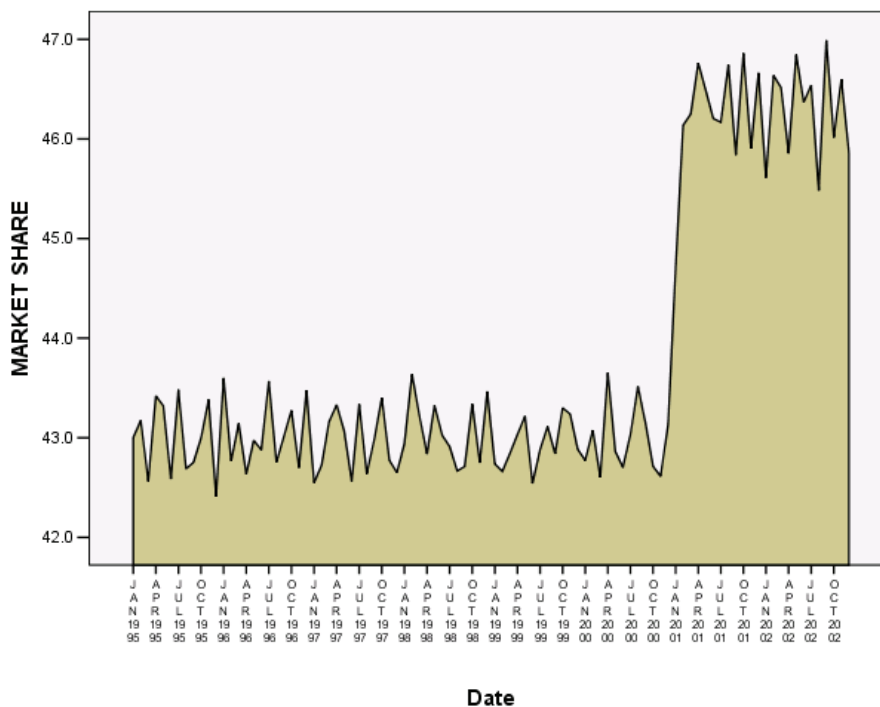
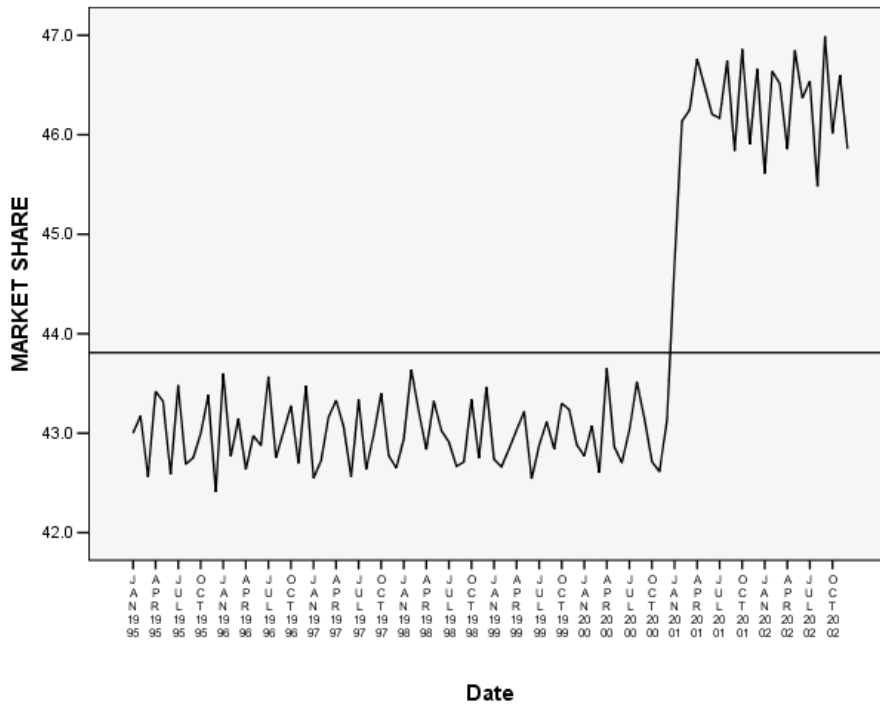


Figure 251-3  
 FORMAT=REFERENCE

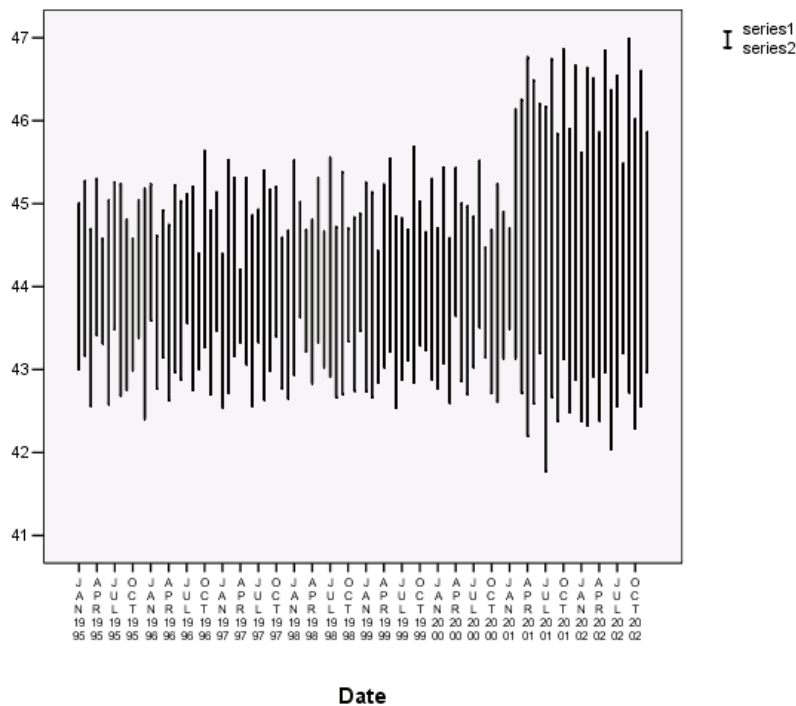


The following formats are available for plots with multiple variables:

- NOJOIN** *Plot the values of each variable named.* Different colors or line patterns are used for multiple variables. Multiple occurrences of the same value for a single observation are plotted using a dollar sign (\$). This is the default format for plots with multiple variables.
- JOIN** *Plot the values of each variable and join the values for each observation.* Values are plotted as described for NOJOIN, and the values for each observation are joined together by a line.
- HILO** *Plot the highest and lowest values across variables for each observation and join the two values together.* The high and low values are plotted as a horizontal bar and are

joined with a line. If more than three variables are specified, HILO is ignored and the default NOJOIN is used.

Figure 251-4  
FORMAT=HILO



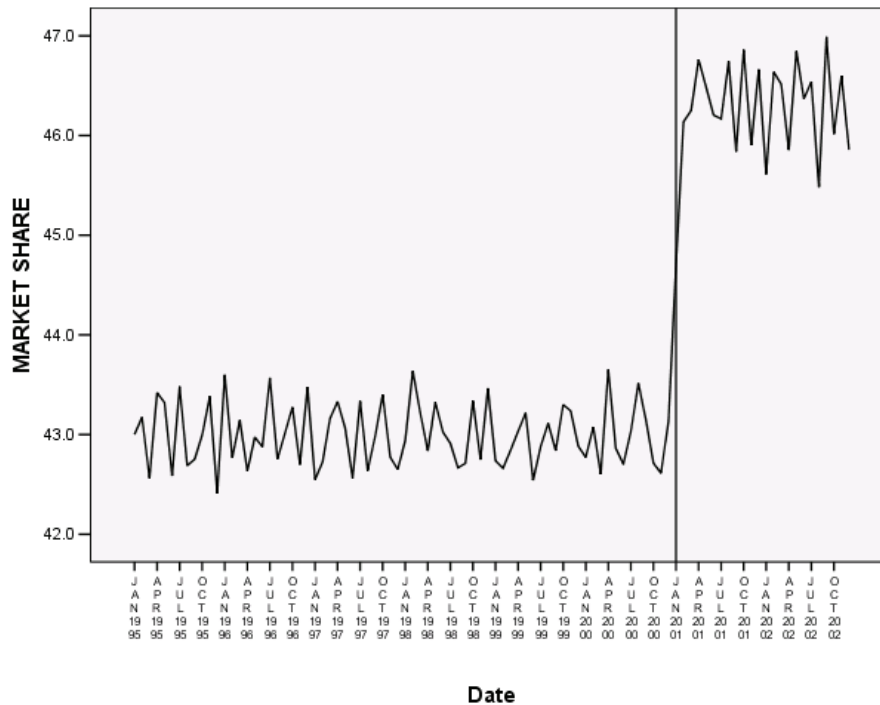
## **MARK Subcommand**

MARK indicates the onset of an intervention variable.

- The onset date is indicated by a vertical reference line.
- The specification on MARK can be either a variable name or an onset date if the *DATE\_* variable exists.
- If a variable is named, the plot indicates where the values of that variable change.
- A date specification follows the same format as the DATE command; that is, a keyword followed by a value. For example, the specification for June 1955 is Y 1955 M 6 (or Y 55 M 6 if only the last two digits of the year are used on DATE).

The following figure shows a plot with January 2001 marked as the onset date.

Figure 251-5  
MARK=Y 2001 M 1



## ***SPLIT Subcommand***

SPLIT specifies how to plot data that have been divided into subgroups by a SPLIT FILE command. The default is UNIFORM.

- UNIFORM**      *Scale uniformly.* The vertical axis is scaled according to the values of the entire dataset.
- SCALE**        *Scale individually.* The vertical axis is scaled according to the values of each individual subgroup.

- If FORMAT=REFERENCE is specified when SPLIT=SCALE, the reference line is placed at the mean of the subgroup. If FORMAT=REFERENCE is specified when SPLIT=UNIFORM, the reference line is placed at the overall mean.

### ***Example***

```
SPLIT FILE BY REGION.
TSPLLOT VARIABLES = TICKETS / SPLIT=SCALE.
```

- In this example, the data have been split into subgroups by *REGION*. The plots produced with the SCALE subcommand have vertical axes that are individually scaled according to the values of each particular region.



## ***APPLY Subcommand***

APPLY allows you to produce a plot using previously defined specifications without having to repeat the TSPLLOT subcommands.

- The only specification on APPLY is the name of a previous model enclosed in quotes. If a model name is not specified, the specifications from the previous TSPLLOT command are used.
- To change one or more specifications of the plot, specify the subcommands of only those portions you want to change after the subcommand APPLY.
- If no variables are specified, the variables that were specified for the original plot are used.
- To plot different variables, enter new variable names before or after the APPLY subcommand.

### ***Example***

```
TSPLLOT VARIABLES = TICKETS
  /LN
  /DIFF=1
  /SDIFF=1
  /PERIOD=12.
TSPLLOT VARIABLES = ROUNDTRP
  /APPLY.
TSPLLOT APPLY
  /NOLOG.
```

- The first command produces a plot of *TICKETS* after a natural log transformation, differencing, and seasonal differencing have been applied.
- The second command plots the values for *ROUNDTRP* using the same subcommands specified for *TICKETS*.
- The third command produces another plot of *ROUNDTRP* but this time without a log transformation. *ROUNDTRP* is still differenced once and seasonally differenced with a periodicity of 12.

# T-TEST

T-TEST is available in the Statistics Base option.

## *One-sample tests:*

```
T-TEST TESTVAL n /VARIABLE=varlist
```

## *Independent-samples tests:*

```
T-TEST GROUPS=varname ( {1,2** } ) /VARIABLES=varlist
                        {value }
                        {value,value}
```

## *Paired-samples tests:*

```
T-TEST PAIRS=varlist [WITH varlist [(PAIRED)]] [/varlist ...]
```

## *All types of tests:*

```
[/MISSING={ANALYSIS**} [INCLUDE]]
          {LISTWISE }
```

```
[/CRITERIA=CI({0.95**})
          {value }
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Examples**

```
T-TEST GROUPS=WORLD(1,3) /VARIABLES=NTCPRI NTCSAL NTCPUR.
```

```
T-TEST PAIRS=TEACHER CONSTRUC MANAGER.
```

## **Overview**

T-TEST compares sample means by calculating Student's *t* and displays the two-tailed probability of the difference between the means. Statistics are available for one-sample (tested against a specified value), independent samples (different groups of cases), or paired samples (different variables). Other procedures that compare group means are ANOVA, ONEWAY, UNIANOVA, GLM, and MANOVA (GLM and MANOVA are available in the Advanced Statistics option).

## **Options**

**Statistics.** There are no optional statistics. All statistics available are displayed by default.

### **Basic Specification**

The basic specification depends on whether you want a one-sample test, an independent-samples test or a paired-samples test. For all types of tests, T-TEST displays Student's  $t$ , degrees of freedom, and two-tailed probabilities, as well as the mean, standard deviation, standard error, and count for each group or variable.

- To request a one-sample test, use the TESTVAL and VARIABLES subcommands. The output includes a One-Sample Statistics table showing univariate statistics and a One-Sample Test table showing the test value, the difference between the sample mean and the test value, and the two-tailed probability level.
- To request an independent-samples test, use the GROUPS and VARIABLES subcommands. The output includes a Group Statistics table showing summary statistics by group for each dependent variable and an Independent-Samples Test table showing both pooled- and separate-variance estimates, along with the  $F$  value used to test homogeneity of variance and its probability. The two-tailed probability is displayed for the  $t$  value.
- To request a paired-samples test, use the PAIRS subcommand. The output includes a Paired Statistics table showing univariate statistics by pairs, a Paired Samples Correlations table showing correlation coefficients and two-tailed probability level for a test of the coefficient for each pair, and a Paired Samples Test table showing the paired differences between the means and two-tailed probability levels for a test of the differences.

### **Subcommand Order**

Subcommands can be named in any order.

### **Operations**

- If a variable specified on GROUPS is a long string, only the short-string portion is used to identify groups in the analysis.
- Probability levels are two-tailed. To obtain the one-tailed probability, divide the two-tailed probability by 2.

### **Limitations**

- Maximum of one TESTVAL and one VARIABLES subcommand per one-sample  $t$  test.
- Maximum of one GROUPS and one VARIABLES subcommand per independent-samples  $t$  test.

## **Examples**

### **One-Sample Test**

```
T-TEST
  /TESTVAL=322
  /VARIABLES=brake
  /CRITERIA=CIN (.90) .
```

- The procedure compares the mean of *brake* with the test value 322.
- CRITERIA requests a 90% confidence interval for the mean difference.

**Independent-Samples Test, using specific values**

```
T-TEST GROUPS=insert(0 1)
/MISSING=ANALYSIS
/VARIABLES=dollars
/CRITERIA=CIN(.95).
```

- The procedure compares the means of the two groups defined by values 0 and 1 of *insert* for variable *dollars*.
- All other options are set to their default values.

**Independent-Samples Test, using a cut point**

```
T-TEST GROUPS=score(50)
/MISSING=ANALYSIS
/VARIABLES=minutes
/CRITERIA=CIN(.95).
```

- The procedure compares the means of the two groups defined by values greater than or equal to 50 (versus all other values) of *score* for variable *minutes*.
- All other options are set to their default values.

**Paired-Samples Test**

```
T-TEST PAIRS=TEACHER CONSTRUC MANAGER.
```

- This paired-samples *t* test compares the means of *TEACHER* with *CONSTRUC*, *TEACHER* with *MANAGER*, and *CONSTRUC* with *MANAGER*.

**Paired-Samples Test, special pairing**

```
T-TEST
PAIRS= tg0 wgt0 WITH tg4 wgt4 (PAIRED)
/CRITERIA=CIN(.95)
/MISSING=ANALYSIS.
```

- The procedure compares the means of *tg0* with *tg4* and *wgt0* with *wgt4*.
- All other options are set to their default values.

**VARIABLES Subcommand**

VARIABLES specifies the dependent variables to be tested in a one-sample or an independent-samples *t* test.

- VARIABLES can specify multiple variables, all of which must be numeric.
- When specified along with TESTVAL, the mean of all cases for each variable is compared with the specified value.
- When specified along with GROUPS, the means of two groups of cases defined by the GROUPS subcommand are compared.
- If both TESTVAL and GROUPS are specified, a one-sample test and an independent-samples test are performed on each variable.

## **TESTVAL Subcommand**

TESTVAL specifies the value with which a sample mean is compared.

- Only one TESTVAL subcommand is allowed.
- Only one value can be specified on the TESTVAL subcommand.

## **GROUPS Subcommand**

GROUPS specifies a variable used to group cases for independent-samples *t* tests.

- GROUPS can specify only one variable, which can be numeric or string.

Any one of three methods can be used to define the two groups for the variable specified on GROUPS:

- Specify a single value in parentheses to group all cases with a value equal to or greater than the specified value into one group and the remaining cases into the other group.
- Specify two values in parentheses to include cases with the first value in one group and cases with the second value in the other group. Cases with other values are excluded.
- If no values are specified on GROUP, T-TEST uses 1 and 2 as default values for numeric variables. There is no default for string variables.

## **PAIRS Subcommand**

PAIRS requests paired-samples *t* tests.

- The minimum specification for a paired-samples test is PAIRS with an analysis list. Only numeric variables can be specified on the analysis list. The minimum analysis list is two variables.
- If keyword WITH is not specified, each variable in the list is compared with every other variable on the list.
- If keyword WITH is specified, every variable to the left of WITH is compared with every variable to the right of WITH. WITH can be used with PAIRED to obtain special pairing.
- To specify multiple analysis lists, use multiple PAIRS subcommands, each separated by a slash. Keyword PAIRS is required only for the first analysis list; a slash can be used to separate each additional analysis list.

**(PAIRED)**      *Special pairing for paired-samples test.* PAIRED must be enclosed in parentheses and must be used with keyword WITH. When PAIRED is specified, the first variable before WITH is compared with the first variable after WITH, the second variable before WITH is compared with the second variable after WITH, and so forth. The same number of variables should be specified before and after WITH; unmatched variables are ignored and a warning message is issued. PAIRED generates an error message if keyword WITH is not specified on PAIRS.

### **Example**

```
T-TEST PAIRS=TEACHER CONSTRUC MANAGER.
T-TEST PAIRS=TEACHER MANAGER WITH CONSTRUC ENGINEER.
```

T-TEST PAIRS=TEACHER MANAGER WITH CONSTRUC ENGINEER (PAIRED) .

- The first T-TEST compares *TEACHER* with *CONSTRUC*, *TEACHER* with *MANAGER*, and *CONSTRUC* with *MANAGER*.
- The second T-TEST compares *TEACHER* with *CONSTRUC*, *TEACHER* with *ENGINEER*, *MANAGER* with *CONSTRUC*, and *MANAGER* with *ENGINEER*. *TEACHER* is not compared with *MANAGER*, and *CONSTRUC* is not compared with *ENGINEER*.
- The third T-TEST compares *TEACHER* with *CONSTRUC* and *MANAGER* with *ENGINEER*.

## ***CRITERIA Subcommand***

CRITERIA resets the value of the confidence interval. Keyword CI is required. You can specify a value between 0 and 1 in the parentheses. The default is 0.95.

## ***MISSING Subcommand***

MISSING controls the treatment of missing values. The default is ANALYSIS.

- ANALYSIS and LISTWISE are alternatives; however, each can be specified with INCLUDE.

<b>ANALYSIS</b>	<i>Delete cases with missing values on an analysis-by-analysis or pair-by-pair basis. For independent-samples tests, cases with missing values for either the grouping variable or the dependent variable are excluded from the analysis of that dependent variable. For paired-samples tests, a case with a missing value for either of the variables in a given pair is excluded from the analysis of that pair. This is the default.</i>
<b>LISTWISE</b>	<i>Exclude cases with missing values listwise. A case with a missing value for any variable specified on either GROUPS or VARIABLES is excluded from any independent-samples test. A case with a missing value for any variable specified on PAIRS is excluded from any paired-samples test.</i>
<b>INCLUDE</b>	<i>Include user-missing values. User-missing values are treated as valid values.</i>

# TWOSTEP CLUSTER

TWOSTEP CLUSTER is available in the Statistics Base option.

TWOSTEP CLUSTER

```
[/CATEGORICAL VARIABLES = varlist]
[/CONTINUOUS VARIABLES = varlist]
[/CRITERIA [INITHRESHOLD({0** })] [MXBRANCH({8**})]
           {value}                    {n }
           [MXLEVEL({3**})] ]
           {n }

[/DISTANCE {EUCLIDEAN }
           {LIKELIHOOD**}]

[/HANDLENOISE {0**}
              {n }

[/INFILE FILE = filename]

[/MEMALLOCATE {64**}
              {n }

[/MISSING {EXCLUDE**}
          {INCLUDE }

[/NOSTANDARDIZE [VARIABLES = varlist]]

[/NUMCLUSTERS {AUTO** {15**} [{AIC }]}]
              {n } {BIC**}
              {FIXED = n }

[/OUTFILE [MODEL = 'file'] [STATE = 'file']]

[/PRINT [IC] [COUNT] [SUMMARY]]

[/VIEWMODEL [DISPLAY = {YES**}
               {NO }
             [EVALUATIONFIELDS = varlist]]

[/SAVE CLUSTER [VARIABLE = varname]]
```

\*\*Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
TWOSTEP CLUSTER
  /CONTINUOUS VARIABLES = INCOME
  /CATEGORICAL VARIABLES = GENDER RACE
  /PRINT SUMMARY.
```

## Overview

TWOSTEP CLUSTER groups observations into clusters based on a nearness criterion. The procedure uses a hierarchical agglomerative clustering procedure in which individual cases are successively combined to form clusters whose centers are far apart. This algorithm is designed to cluster large numbers of cases. It passes the data once to find the cluster centers and again to assign cluster memberships. In addition to the benefit of few data passes, the procedure allows the user to set the amount of memory used by the clustering algorithm.

### Basic Features

**Cluster Features (CF) Tree.** TWOSTEP CLUSTER clusters observations by building a data structure called the **CF tree**, which contains the cluster centers. The CF tree is grown during the first stage of clustering and values are added to its leaves if they are close to the cluster center of a particular leaf.

**Distance Measure.** Two types of distance measures are offered—the traditional Euclidean distance and the likelihood distance. The former is available when no categorical variables are specified. The latter is especially useful when categorical variables are used. The likelihood function is computed using the normal density for continuous variables and the multinomial probability mass function for categorical variables. All variables are treated as independent.

**Tuning the Algorithm.** You can control the values of algorithm-tuning parameters with the **CRITERIA** subcommand.

**Noise Handling.** The clustering algorithm can optionally retain any outliers that do not fit in the CF tree. If possible, these values will be placed in the CF tree after it is completed. Otherwise, TWOSTEP CLUSTER will discard them after preclustering.

**Missing Values.** TWOSTEP CLUSTER will delete listwise any records with missing fields.

**Numclusters.** This subcommand specifies the number of clusters into which the data will be partitioned. The user may tell TWOSTEP CLUSTER to automatically select the number of clusters.

**Optional Output.** You can specify output to an XML file with the **OUTFILE** subcommand. The cluster membership for each case used can be saved to the active dataset with the **SAVE** subcommand.

**Weights.** TWOSTEP CLUSTER ignores specification on the **WEIGHT** command.

### Basic Specification

- The minimum specification is a list of variables, either categorical or continuous, to be clustered and at least one of the following subcommands: **OUTFILE**, **PRINT**, or **SAVE**.
- The number of clusters may be specified with the **NUMCLUSTERS** subcommand.
- Unless the **NOSTANDARDIZE** subcommand is given, TWOSTEP CLUSTER will standardize all continuous variables.
- If **DISTANCE** is Euclidean, TWOSTEP CLUSTER will accept only continuous variables.



**Subcommand Order**

- The subcommands can be specified in any order.

**Syntax Rules**

- Minimum syntax: a variable must be specified.
- Empty subcommands are silently ignored.
- Variables listed in the CONTINUOUS subcommand must be numeric.
- If a subcommand is issued more than once, TWOSTEP CLUSTER will ignore all but the last issue.

**Variable List**

The variable lists specify those variables to be clustered. The first variable list specifies only continuous variables, and the second list specifies only categorical variables (that is, the two lists are disjoint).

**CATEGORICAL Subcommand**

The CATEGORICAL subcommand specifies a list of categorical variables.

**Example**

```
TWOSTEP CLUSTER
  /CATEGORICAL VARIABLES = RACE GENDER CITIZEN
  /PRINT SUMMARY COUNT.
```

This tells TWOSTEP CLUSTER to cluster the categorical variables *RACE*, *GENDER*, and *CITIZEN*. Summary statistics by cluster and cluster frequencies are output in tables.

**CONTINUOUS Subcommand**

The CONTINUOUS subcommand specifies a list of scale variables.

**Example**

```
TWOSTEP CLUSTER
  /CATEGORICAL VARIABLES = RACE GENDER CITIZEN
  /CONTINUOUS VARIABLES = INCOME
  /PRINT SUMMARY COUNT.
```

This tells TWOSTEP CLUSTER to cluster the categorical variables *RACE*, *GENDER*, and *CITIZEN*, and the numeric variable *INCOME*. Summary statistics by cluster and cluster frequencies are output in tables.

## **CRITERIA Subcommand**

The CRITERIA subcommand specifies the following settings for the clustering algorithm:

<b>INITTHRESHOLD</b>	<i>The initial threshold used to grow the CF tree. The default is 0. If inserting a specific case into a leaf of the CF tree would yield tightness less than the threshold, the leaf is not split. If the tightness exceeds the threshold, the leaf is split.</i>
<b>MXBRANCH</b>	<i>The maximum number of child nodes that a leaf node can have. The default is 8.</i>
<b>MXLEVEL</b>	<i>The maximum number of levels that the CF tree can have. The default is 3.</i>

## **DISTANCE Subcommand**

The distance subcommand determines how distances will be computed between clusters.

<b>EUCLIDEAN</b>	<i>Use the Euclidean distance to compute distances between clusters. You may select Euclidean distance if all variables are continuous. TWOSTEP CLUSTER will return a syntax error if you specify Euclidean distance with non-numeric variables.</i>
<b>LIKELIHOOD</b>	<i>Use the minus log-likelihood to compute distances between clusters. This is the default. The likelihood function is computed assuming all variables are independent. Continuous variables are assumed to be normally distributed, and categorical variables are assumed to be multinomially distributed.</i>

## **HANDLENOISE Subcommand**

The HANDLENOISE subcommand tells TWOSTEP CLUSTER to treat outliers specially during clustering. During growth of the CF tree, this subcommand is relevant only if the CF tree fills. The CF tree is full if it cannot accept any more cases in a leaf node and no leaf node can be split. The default value of HANDLENOISE is 0, equivalent to no noise handling.

- If the CF tree fills and HANDLENOISE is greater than 0, the CF tree will be regrown after placing any data in sparse leaves into their own noise leaf. A leaf is considered sparse if the ratio of the number of cases in the sparse leaf to the number of cases in the largest leaf is less than HANDLENOISE. After the tree is grown, the outliers will be placed in the CF tree, if possible. If not, the outliers are discarded for the second phase of clustering.
- If the CF tree fills and HANDLENOISE is equal to 0, the threshold will be increased and CF tree regrown with all cases. After final clustering, values that cannot be assigned to a cluster are labeled outliers. The outlier cluster is given an identification number of -1. The outlier cluster is not included in the count of the number of clusters; that is, if you specify  $n$  clusters and noise handling, TWOSTEP CLUSTER will output  $n$  clusters and one noise cluster.

### **Example**

```
TWOSTEP CLUSTER
  /CATEGORICAL VARIABLES = RACE GENDER CITIZEN
  /CONTINUOUS VARIABLES = INCOME
  /HANDLENOISE 25
  /PRINT SUMMARY COUNT.
```

This tells `TWOSTEP CLUSTER` to cluster the categorical variables *RACE*, *GENDER* and *CITIZEN*, and the numeric variable *INCOME*. If the CF tree fills, a noise leaf is constructed from cases whose leaves contain fewer than 25 percent of the cases contained by the largest leaf. The CF tree is then regrown, ignoring the noise leaf. After the tree is regrown, cases from the noise leaf are checked to see if they fit any of the leaves in the new tree. Any cases that still do not fit are discarded as outliers. Summary statistics by cluster and cluster frequencies are output in tables.

## ***INFILE Subcommand***

The `INFILE` subcommand causes `TWOSTEP CLUSTER` to update a cluster model whose CF Tree has been saved as an XML file with the `OUTFILE` subcommand and `STATE` keyword. The model will be updated with the data in the active file. The user must supply variable names in the active file in the order they are stored in the XML file. `TWOSTEP CLUSTER` will update the cluster model in memory only, leaving unaltered the XML file.

- If the `INFILE` subcommand is given, `TWOSTEP CLUSTER` will ignore the `CRITERIA`, `DISTANCE`, `HANDLENOISE` and `MEMALLOCATE` subcommands, if given.

## ***MEMALLOCATE Subcommand***

The `MEMALLOCATE` subcommand specifies the maximum amount of memory in megabytes (MB) that the cluster algorithm should use. If the procedure exceeds this maximum, it will use the disk to store information that will not fit in memory.

- The minimum value that you can specify is 4. If this subcommand is not specified, the default value is 64 MB.
- Consult your system administrator for the largest value you can specify on your system.

## ***MISSING Subcommand***

The `MISSING` subcommand specifies how to handle cases with user-missing values.

- If this subcommand is not specified, the default is `EXCLUDE`.
- `TWOSTEP CLUSTER` deletes any case with a system-missing value.
- Keywords `EXCLUDE` and `INCLUDE` are mutually exclusive. Only one of them can be specified.

**EXCLUDE**      *Exclude both user-missing and system-missing values. This is the default.*

**INCLUDE**      *User-missing values are treated as valid. System-missing values cannot be included in the analysis.*

## ***NOSTANDARDIZE Subcommand***

The `NOSTANDARDIZE` subcommand will prevent `TWOSTEP CLUSTER` from standardizing the continuous variables specified with the `VARIABLES` keyword. If this subcommand is not specified, or given without a variable list, `TWOSTEP CLUSTER` will standardize all continuous variables by subtracting the mean and dividing by the standard deviation.

## **NUMCLUSTERS Subcommand**

The NUMCLUSTERS subcommand specifies the number of clusters into which the data will be partitioned.

- AUTO**      *Automatic selection of the number of clusters.* Under AUTO, you may specify a maximum number of possible clusters. TWOSTEP CLUSTER will search for the best number of clusters between 1 and the maximum using the criterion that you specify. The criterion for deciding the number of clusters can be either the Bayesian Information Criterion (BIC) or Akaike Information Criterion (AIC). TWOSTEP CLUSTER will find at least one cluster if the AUTO keyword is given.
- FIXED**      *User-specified number of clusters.* Specify a positive integer.

### **Examples**

```
TWOSTEP CLUSTER
/CONTINUOUS VARIABLES = INCOME
/CATEGORICAL VARIABLES = GENDER RACE
/NUMCLUSTERS AUTO 10 AIC
/PRINT SUMMARY COUNT.
```

TWOSTEP CLUSTER uses the variables *RACE*, *GENDER* and *INCOME* for clustering. Specifications on the NUMCLUSTERS subcommand will instruct the procedure to automatically search for the number of clusters using the Akaike Information Criterion and require the answer to lie between 1 and 10.

```
TWOSTEP CLUSTER
/CONTINUOUS VARIABLES = INCOME
/CATEGORICAL VARIABLES = RACE GENDER
/NUMCLUSTERS FIXED 7
/PRINT SUMMARY COUNT.
```

Here the procedure will find exactly seven clusters.

## **OUTFILE Subcommand**

The OUTFILE subcommand directs TWOSTEP CLUSTER to write its output to the specified filename as XML.

- MODEL**      *Save the final model output.*
- STATE**      *Save the CF tree.* Use the STATE keyword if you want to update the model later.

You must supply a valid filename on your operating system. We recommend specifying the full path of the filename.

## **PRINT Subcommand**

The PRINT subcommand causes TWOSTEP CLUSTER to print tables related to each cluster.

<b>IC</b>	<i>Information criterion.</i> Prints the chosen information criterion (AIC or BIC) for different numbers of clusters. If this keyword is specified when the AUTO keyword is not used with the NUMCLUSTERS subcommand, TWOSTEP CLUSTER will skip this keyword and issue a warning. TWOSTEP CLUSTER will ignore this keyword if AUTO 1 is specified in the NUMCLUSTERS subcommand.
<b>SUMMARY</b>	<i>Descriptive statistics by cluster.</i> This option prints two tables describing the variables in each cluster. In one table, means and standard deviations are reported for continuous variables. The other table reports frequencies of categorical variables. All values are separated by cluster.
<b>COUNT</b>	<i>Cluster frequencies.</i> This option prints a table containing a list of clusters and how many observations are in each cluster.

## **VIEWMODEL Subcommand**

The VIEWMODEL subcommand controls options for display of model-related output, including tables and charts. Tables in the model view include a model summary and cluster-by-features grid. Graphical output in the model view includes a cluster quality chart, cluster sizes, variable importance, cluster comparison grid, and cell information.

### **DISPLAY Keyword**

The DISPLAY keyword specifies whether to display the model view in the output.

<b>YES</b>	<i>Display the model view.</i> This is the default.
<b>NO</b>	<i>Do not display the model view.</i>

### **EVALUATIONFIELDS Keyword**

The EVALUATIONFIELDS keyword calculates cluster data for variables that were not used in cluster creation. Evaluation fields can be displayed along with the input features in the model viewer by selecting them in the Display subdialog. The EVALUATIONFIELDS keyword is optional and its varlist can include fields (variables) in the dataset that are not used to create the cluster. Fields that are used to create clusters will be ignored. Fields with missing values are ignored.

```
TWOSTEP CLUSTER
  /CATEGORICAL VARIABLES=var1 var2 var3
  /CONTINUOUS VARIABLES=var4 var5
  /VIEWMODEL DISPLAY=YES EVALUATIONFIELDS=var1 var2 var8 var9
```

- The procedure creates a cluster model using *var1*, *var2*, *var3* as categorical variables and *var4* and *var5* as continuous variables, and produces a model viewer with *var8* and *var9* available as evaluation fields.

## ***SAVE Subcommand***

The `SAVE` subcommand allows you to save cluster output to the active dataset.

**CLUSTER**     *Save the cluster identification.* The cluster number for each case is saved; the user may specify a variable name using the `VARIABLE` keyword, otherwise, it is saved to `TSC_n`, where *n* is a positive integer indicating the ordinal of the `SAVE` operation completed by this procedure in a given session.

# UNIANOVA

UNIANOVA is available in the Statistics Base option.

```
UNIANOVA dependent var [BY factor list [WITH covariate list]]
[/RANDOM=factor factor...]
[/REGWGT=varname]
[/METHOD=SSTYPE({1  })]
                {2  }
                {3**}
                {4  }
[/INTERCEPT=[INCLUDE**] [EXCLUDE]]
[/MISSING=[INCLUDE] [EXCLUDE**]]
[/CRITERIA=[EPS({1E-8**})] [ALPHA({0.05**})]
           {a      }      {a      }
[/PRINT = [DESCRIPTIVE] [HOMOGENEITY] [PARAMETER][ETASQ]
          [GEF] [LOF] [OPOWER] [TEST(LMATRIX)]]
[/PLOT=[SPREADLEVEL] [RESIDUALS]
       [PROFILE (factor factor*factor factor*factor*factor ...)]
       [WITH(covariate={value} [...])]]
       {MEAN }
[/TEST=effect VS {linear combination [DF(df)]}]
               {value DF (df)          }
[/LMATRIX={{["label"] effect list effect list ...;...}}
           {{["label"] effect list effect list ...      }
           {"label"} ALL list; ALL...                    }
           {"label"} ALL list                            }
[/KMATRIX= {number      }
           {number;...}
[/CONTRAST (factor name)={DEVIATION[(refcat)]**      }
                       {SIMPLE [(refcat)]           }
                       {DIFFERENCE                  }
                       {HELMERT                     }
                       {REPEATED                    }
                       {POLYNOMIAL [({1,2,3...})]   }
                       {metric }                   }
                       {SPECIAL (matrix)            }
[/POSTHOC =effect [effect...]
           ({SNK} [TUKEY] [BTUKEY] [DUNCAN]
            [SCHEFFE] [DUNNETT(refcat)] [DUNNETTL(refcat)]
            [DUNNETTR(refcat)] [BONFERRONI] [LSD] [SIDAK]
            [GT2] [GABRIEL] [FREGW] [QREGW] [T2] [T3] [GH] [C]
            [WALLER ({100** })]])
           {kratio}
           [VS effect]
[/EMMEANS=TABLES({OVERALL      }) [WITH(covariate={value} [...])]]
               {factor          }      {MEAN }
               {factor*factor...}
           [COMPARE ADJ(LSD) (BONFERRONI) (SIDAK)]
[/SAVE=[tempvar [(name)]] [tempvar [(name)]]...]
[/OUTFILE={{COVB ('savfile'|'dataset')}}
          {CORB ('savfile'|'dataset')}}]
```

```

[EFFECT('savfile'|'dataset')] [DESIGN('savfile'|'dataset')]

[/DESIGN={{INTERCEPT...} } ]
      {{effect effect...}}

```

\*\* Default if the subcommand or keyword is omitted.

Temporary variables (tempvar) are:

PRED, WPRED, RESID, WRESID, DRESID, ZRESID, SRESID, SEPREP, COOK, LEVER

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 17.0

- POSTHOC subcommand: T2, T3, GH, and C keywords are not valid when multiple factors in the model.

### **Example**

```

UNIANOVA YIELD BY SEED FERT
  /DESIGN.

```

## **Overview**

This section describes the use of UNIANOVA for univariate analyses. The UNIANOVA procedure provides regression analysis and analysis of variance for one dependent variable by one or more factors and/or variables.

### **Options**

**Design Specification.** You can specify which terms to include in the design on the DESIGN subcommand. This allows you to estimate a model other than the default full factorial model, incorporate factor-by-covariate interactions or covariate-by-covariate interactions, and indicate nesting of effects.

**Contrast Types.** You can specify contrasts other than the default deviation contrasts on the CONTRAST subcommand.

**Optional Output.** You can choose from a wide variety of optional output on the PRINT subcommand. Output appropriate to univariate designs includes descriptive statistics for each cell, parameter estimates, Levene's test for equality of variance across cells, partial eta-squared for each effect and each parameter estimate, the general estimable function matrix, and a contrast coefficients table (**L'** matrix). The OUTFILE subcommand allows you to write out the covariance or correlation matrix, the design matrix, or the statistics from the between-subjects ANOVA table into a separate data file.



Using the `EMMEANS` subcommand, you can request tables of estimated marginal means of the dependent variable and their standard deviations. The `SAVE` subcommand allows you to save predicted values and residuals in weighted or unweighted and standardized or unstandardized forms. You can specify different means comparison tests for comparing all possible pairs of cell means using the `POSTHOC` subcommand. In addition, you can specify your own hypothesis tests by specifying an **L** matrix and a **K** matrix to test the univariate hypothesis  $\mathbf{LB} = \mathbf{K}$ .

### **Basic Specification**

- The basic specification is a variable list identifying the dependent variable, the factors (if any), and the covariates (if any).
- By default, UNIANOVA uses a model that includes the intercept term, the covariate (if any), and the full factorial model, which includes all main effects and all possible interactions among factors. The intercept term is excluded if it is excluded in the model by specifying the keyword `EXCLUDE` on the `INTERCEPT` subcommand. Sums of squares are calculated and hypothesis tests are performed using type-specific estimable functions. Parameters are estimated using the normal equation and a generalized inverse of the SSCP matrix.

### **Subcommand Order**

- The variable list must be specified first.
- Subcommands can be used in any order.

### **Syntax Rules**

- For many analyses, the UNIANOVA variable list and the `DESIGN` subcommand are the only specifications needed.
- If you do not enter a `DESIGN` subcommand, UNIANOVA will use a full factorial model, with main effects of covariates, if any.
- At least one dependent variable must be specified, and at least one of the following must be specified: `INTERCEPT`, a factor, or a covariate. The design contains the intercept by default.
- If more than one `DESIGN` subcommand is specified, only the last one is in effect.
- Dependent variables and covariates must be numeric, but factors can be numeric or string variables.
- If a string variable is specified as a factor, only the first eight characters of each value are used in distinguishing among values.
- If more than one `MISSING` subcommand is specified, only the last one is in effect.
- The following words are reserved as keywords or internal commands in the UNIANOVA procedure:  
`INTERCEPT`, `BY`, `WITH`, `ALL`, `OVERALL`, `WITHIN`  
Variable names that duplicate these words should be changed before you run UNIANOVA.

### **Limitations**

- Any number of factors can be specified, but if the number of factors plus the number of split variables exceeds 18, the Descriptive Statistics table is not printed even when you request it.

- Memory requirements depend primarily on the number of cells in the design. For the default full factorial model, this equals the product of the number of levels or categories in each factor.

## Example

```
UNIANOVA YIELD BY SEED FERT WITH RAINFALL
  /PRINT=DESCRIPTIVE PARAMETER
  /DESIGN.
```

- *YIELD* is the dependent variable; *SEED* and *FERT* are factors; *RAINFALL* is a covariate.
- The `PRINT` subcommand requests the descriptive statistics for the dependent variable for each cell and the parameter estimates, in addition to the default tables Between-Subjects Factors and Univariate Tests.
- The `DESIGN` subcommand requests the default design, a full factorial model with a covariate. This subcommand could have been omitted or could have been specified in full as

```
/DESIGN = INTERCEPT RAINFALL, SEED, FERT, SEED BY FERT.
```

### Two-Factor analysis of variance

```
UNIANOVA amtspent BY gender style
  /METHOD=SSTYPE(3)
  /INTERCEPT=INCLUDE
  /POSTHOC=style(TUKEY)
  /PLOT=PROFILE(style*gender)
  /EMMEANS = TABLES(gender*style)
  /PRINT=ETASQ HOMOGENEITY DESCRIPTIVE
  /PLOT=SPREADLEVEL
  /CRITERIA=ALPHA(.05)
  /DESIGN=gender style gender*style.
```

- The procedure specifies *amtspent* as the target variable and *gender* and *style* as factors.
- `POSTHOC` requests Tukey tests for *style*.
- The two `PLOT` subcommands request a profile plot for *style\*gender* and a spread vs. level plot.
- `EMMEANS` requests estimated marginal means for *gender\*style*
- `PRINT` requests descriptive statistics, estimates of effect size, and homogeneity tests.
- `DESIGN` specifies *gender*, *style*, and *gender\*style* as effects in the model.
- All other options are set to their default values.

### Analysis of covariance

```
UNIANOVA
  incaft BY prog WITH incbef
  /METHOD = SSTYPE(3)
  /INTERCEPT = INCLUDE
  /PRINT = DESCRIPTIVE ETASQ PARAMETER HOMOGENEITY
  /PLOT = SPREADLEVEL
  /CRITERIA = ALPHA(.05)
  /DESIGN = incbef prog .
```

- The procedure specifies *incaft* as the target variable, *prog* as a factor, and *incbef* as a covariate.

- PRINT requests descriptive statistics, estimates of effect size, parameter estimates, and homogeneity tests.
- PLOT requests a spread vs. level plot.
- DESIGN specifies *incbef* and *prog* as effects in the model.
- All other options are set to their default values.

### **Random effects**

```
UNIANOVA
  amtspent BY usecoup shopfor storeid
  /RANDOM = storeid
  /METHOD = SSTYPE(3)
  /INTERCEPT = INCLUDE
  /PRINT = ETASQ
  /CRITERIA = ALPHA(.05)
  /DESIGN = usecoup shopfor shopfor*usecoup storeid .
```

- The procedure specifies *amtspent* as the target variable and *usecoup*, *shopfor*, and *storeid* as factors.
- RANDOM specifies that *storeid* is a random effect.
- PRINT requests estimates of effect size.
- DESIGN specifies *usecoup*, *shopfor*, *shopfor\*usecoup*, and *storeid* as effects in the model.
- All other options are set to their default values.

## **UNIANOVA Variable List**

The variable list specifies the dependent variable, the factors, and the covariates in the model.

- The dependent variable must be the first specification on UNIANOVA.
- The names of the factors follow the dependent variable. Use the keyword BY to separate the factors from the dependent variable.
- Enter the covariates, if any, following the factors. Use the keyword WITH to separate covariates from factors (if any) and the dependent variable.

### **Example**

```
UNIANOVA DEPENDNT BY FACTOR1 FACTOR2, FACTOR3.
```

- In this example, three factors are specified.
- A default full factorial model is used for the analysis.

### **Example**

```
UNIANOVA Y BY A WITH X
  /DESIGN.
```

- In this example, the DESIGN subcommand requests the default design, which includes the intercept term, the covariate *X*, and the factor *A*.

## ***RANDOM Subcommand***

RANDOM allows you to specify which effects in your design are random. When the RANDOM subcommand is used, a table of expected mean squares for all effects in the design is displayed, and an appropriate error term for testing each effect is calculated and used automatically.

- Random always implies a univariate mixed-model analysis.
- If you specify an effect on RANDOM, higher-order effects containing the specified effect (excluding any effects containing covariates) are automatically treated as random effects.
- The keyword INTERCEPT and effects containing covariates are not allowed on this subcommand.
- When the RANDOM subcommand is used, the appropriate error terms for the hypothesis testing of all effects in the model are automatically computed and used.
- More than one RANDOM subcommand is allowed. The specifications are accumulated.

### ***Example***

```
UNIANOVA DEP BY A B
  /RANDOM = B
  /DESIGN = A,B, A*B.
```

- In the example, effects  $B$  and  $A*B$  are considered as random effects. Notice that if only effect  $B$  is specified in the RANDOM subcommand,  $A*B$  is automatically considered as a random effect.
- The hypothesis testing for each effect ( $A$ ,  $B$ , and  $A*B$ ) in the design will be carried out using the appropriate error term, which is calculated automatically.

## ***REGWGT Subcommand***

The only specification on REGWGT is the name of the variable containing the weights to be used in estimating a weighted least-squares model.

- Specify a numeric weight variable name following the REGWGT subcommand. Only observations with positive values in the weight variable will be used in the analysis.
- If more than one REGWGT subcommand is specified, only the last one is in effect.

### ***Example***

```
UNIANOVA OUTCOME BY TREATMNT
  /REGWGT WT.
```

- The procedure performs a weighted least-squares analysis. The variable  $WT$  is used as the weight variable.

## METHOD Subcommand

METHOD controls the computational aspects of the UNIANOVA analysis. You can specify one of four different methods for partitioning the sums of squares. If more than one METHOD subcommand is specified, only the last one is in effect.

- SSTYPE(1)**      *Type I sum-of-squares method.* The Type I sum-of-squares method is also known as the hierarchical decomposition of the sum-of-squares method. Each term is adjusted only for the terms that precede it on the DESIGN subcommand. Under a balanced design, it is an orthogonal decomposition, and the sums of squares in the model add up to the total sum of squares.
- SSTYPE(2)**      *Type II sum-of-squares method.* This method calculates the sum of squares of an effect in the model adjusted for all other “appropriate” effects. An appropriate effect is one that corresponds to all effects that do not contain the effect being examined. For any two effects  $F1$  and  $F2$  in the model,  $F1$  is said to be **contained** in  $F2$  under the following three conditions: 1) both effects  $F1$  and  $F2$  have the same covariate, if any, 2)  $F2$  consists of more factors than  $F1$ , and 3) all factors in  $F1$  also appear in  $F2$ . The intercept effect is treated as contained in all of the pure factor effects. However, it is not contained in any effect involving a covariate. No effect is contained in the intercept effect. Thus, for any one effect  $F$  of interest, all other effects in the model can be classified as in one of the following two groups: the effects that do not contain  $F$  or the effects that contain  $F$ . If the model is a main-effects design (that is, only main effects are in the model), the Type II sum-of-squares method is equivalent to the regression approach sums of squares. This means that each main effect is adjusted for every other term in the model.
- SSTYPE(3)**      *Type III sum-of-squares method.* This is the default. This method calculates the sum of squares of an effect  $F$  in the design as the sum of squares adjusted for any other effects that do not contain it, and *orthogonal* to any effects (if any) that contain it. The Type III sums of squares have one major advantage—they are invariant with respect to the cell frequencies as long as the general form of estimability remains constant. Hence, this type of sums of squares is often used for an unbalanced model with no missing cells. In a factorial design with no missing cells, this method is equivalent to the Yates’ weighted squares of means technique, and it also coincides with the overparameterized  $\Sigma$ -restricted model.
- SSTYPE(4)**      *Type IV sum-of-squares method.* This method is designed for a situation in which there are missing cells. For any effect  $F$  in the design, if  $F$  is not contained in any other effect, then Type IV = Type III = Type II. When  $F$  is contained in other effects, then Type IV distributes the contrasts being made among the parameters in  $F$  to all higher-level effects equitably.

### Example

```
UNIANOVA DEP BY A B C
  /METHOD=SSTYPE(3)
  /DESIGN=A, B, C.
```

- The design is a main-effects model.
- The METHOD subcommand requests that the model be fitted with Type III sums of squares.

## **INTERCEPT Subcommand**

INTERCEPT controls whether an intercept term is included in the model. If more than one INTERCEPT subcommand is specified, only the last one is in effect.

- INCLUDE**            *Include the intercept term.* The intercept (constant) term is included in the model. This is the default.
- EXCLUDE**           *Exclude the intercept term.* The intercept term is excluded from the model. Specification of the keyword INTERCEPT on the DESIGN subcommand overrides INTERCEPT = EXCLUDE.

## **MISSING Subcommand**

By default, cases with missing values for any of the variables on the UNIANOVA variable list are excluded from the analysis. The MISSING subcommand allows you to include cases with user-missing values.

- If MISSING is not specified, the default is EXCLUDE.
- Pairwise deletion of missing data is not available in UNIANOVA.
- Keywords INCLUDE and EXCLUDE are mutually exclusive.
- If more than one MISSING subcommand is specified, only the last one is in effect.

- EXCLUDE**           *Exclude both user-missing and system-missing values.* This is the default when MISSING is not specified.
- INCLUDE**           *User-missing values are treated as valid.* System-missing values cannot be included in the analysis.

## **CRITERIA Subcommand**

CRITERIA controls the statistical criteria used to build the models.

- More than one CRITERIA subcommand is allowed. The specifications are accumulated. Conflicts across CRITERIA subcommands are resolved using the conflicting specification given on the last CRITERIA subcommand.
- The keyword must be followed by a positive number in parentheses.

**EPS(n)**            *The tolerance level in redundancy detection.* This value is used for redundancy checking in the design matrix. The default value is 1E-8.

**ALPHA(n)**         *The alpha level.* This keyword has two functions. First, it gives the alpha level at which the power is calculated for the  $F$  test. Once the noncentrality parameter for the alternative hypothesis is estimated from the data, then the power is the probability that the test statistic is greater than the critical value under the alternative hypothesis. The second function of alpha is to specify the level of the confidence interval. If the alpha level specified is  $n$ , the value  $(1-n) \times 100$  indicates the level of confidence for all individual and simultaneous confidence intervals generated for the specified model. The value of  $n$  must be between 0 and 1, exclusive. The default value of alpha is 0.05. This means that the default power calculation is at the 0.05 level, and the default level of the confidence intervals is 95%, since  $(1-0.05) \times 100 = 95$ .

## **PRINT Subcommand**

PRINT controls the display of optional output.

- Some PRINT output applies to the entire UNIANOVA procedure and is displayed only once.
- Additional output can be obtained on the EMMEANS, PLOT, and SAVE subcommands.
- Some optional output may greatly increase the processing time. Request only the output you want to see.
- If no PRINT command is specified, default output for a univariate analysis includes a factor information table and a Univariate Tests table (ANOVA) for all effects in the model.
- If more than one PRINT subcommand is specified, only the last one is in effect.

The following keywords are available for UNIANOVA univariate analyses.

<b>DESCRIPTIVES</b>	<i>Basic information about each cell in the design.</i> Observed means, standard deviations, and counts for the dependent variable in all cells. The cells are constructed from the highest-order crossing of the factors. If the number of factors plus the number of split variables exceeds 18, the Descriptive Statistics table is not printed.
<b>HOMOGENEITY</b>	<i>Tests of homogeneity of variance.</i> Levene's test for equality of variances for the dependent variable across all level combinations of the factors. If there are no factors, this keyword is not valid.
<b>PARAMETER</b>	<i>Parameter estimates.</i> Parameter estimates, standard errors, <i>t</i> tests, and confidence intervals for each test.
<b>OPOWER</b>	<i>Observed power.</i> The observed power for each test.
<b>LOF</b>	<i>Lack of fit.</i> Lack of fit test that allows you to determine if the current model adequately accounts for the relationship between the response variable and the predictors.
<b>ETASQ</b>	Partial eta-squared ( $\eta^2$ ). This value is an overestimate of the actual effect size in an <i>F</i> test. It is defined as $\eta^2 = \frac{df_h \times F}{df_h \times F + df_e}$ where <i>F</i> is the test statistic and <i>dfh</i> and <i>dfe</i> are its degrees of freedom and degrees of freedom for error. The keyword EFSIZE can be used in place of ETASQ.
<b>GEF</b>	<i>General estimable function table.</i> This table shows the general form of the estimable functions.
<b>TEST(LMATRIX)</b>	<i>Set of contrast coefficients (L) matrices.</i> The transpose of the <b>L</b> matrix ( <b>L'</b> ) is displayed. This set always includes one matrix displaying the estimable function for each effect appearing or implied in the DESIGN subcommand. Also, any <b>L</b> matrices generated by the LMATRIX or CONTRAST subcommands are displayed. TEST(ESTIMABLE) can be used in place of TEST(LMATRIX).

### **Example**

```
UNIANOVA DEP BY A B WITH COV
  /PRINT=DESCRIPTIVE, TEST(LMATRIX), PARAMETER
  /DESIGN.
```

- Since the design in the `DESIGN` subcommand is not specified, the default design is used. In this case, the design includes the intercept term, the covariate `COV`, and the full factorial terms of `A` and `B`, which are `A`, `B`, and `A*B`.
- For each combination of levels of `A` and `B`, the descriptive statistics of `DEP` are displayed.
- The set of `L` matrices that generates the sums of squares for testing each effect in the design is displayed.
- The parameter estimates, their standard errors, *t* tests, confidence intervals, and the observed power for each test are displayed.

## **PLOT Subcommand**

`PLOT` provides a variety of plots useful in checking the assumptions needed in the analysis. The `PLOT` subcommand can be specified more than once. All of the plots requested on each `PLOT` subcommand are produced.

Use the following keywords on the `PLOT` subcommand to request plots:

<b>SPREADLEVEL</b>	<i>Spread-versus-level plots.</i> Plots of observed cell means versus standard deviations, and versus variances.
<b>RESIDUALS</b>	<i>Observed by predicted by standardized residuals plot.</i> A plot is produced for each dependent variable. In a univariate analysis, a plot is produced for the single dependent variable.
<b>PROFILE</b>	<p><i>Line plots of dependent variable means for one-way, two-way, or three-way crossed factors.</i> The <code>PROFILE</code> keyword must be followed by parentheses containing a list of one or more factor combinations. All factors specified (either individual or crossed) must be made up of only valid factors on the factor list. Factor combinations on the <code>PROFILE</code> keyword may use an asterisk (*) or the keyword <code>BY</code> to specify crossed factors. A factor cannot occur in a single factor combination more than once.</p> <p>The order of factors in a factor combination is important, and there is no restriction on the order of factors. If a single factor is specified after the <code>PROFILE</code> keyword, a line plot of estimated means at each level of the factor is produced. If a two-way crossed factor combination is specified, the output includes a multiple-line plot of estimated means at each level of the first specified factor, with a separate line drawn for each level of the second specified factor. If a three-way crossed factor combination is specified, the output includes multiple-line plots of estimated means at each level of the first specified factor, with separate lines for each level of the second factor, and separate plots for each level of the third factor.</p> <p>The <code>PROFILE</code> keyword may be followed by the <code>WITH</code> keyword and a parenthesized list of covariates and values. Plots are produced with the values of the covariates fixed at the specified values. The keyword <code>MEAN</code> causes the covariate to be fixed at its observed mean value.</p>

### **Example**

```
UNIANOVA DEP BY A B
  /PLOT = SPREADLEVEL PROFILE(A A*B A*B*C) WITH(X=1.0)
  /DESIGN.
```

Assume each of the factors `A`, `B`, and `C` has three levels.



- Spread-versus-level plots are produced showing observed cell means versus standard deviations and observed cell means versus variances.
- Five profile plots are produced. For factor  $A$ , a line plot of estimated means at each level of  $A$  is produced (one plot). For the two-way crossed factor combination  $A*B$ , a multiple-line plot of estimated means at each level of  $A$ , with a separate line for each level of  $B$ , is produced (one plot). For the three-way crossed factor combination  $A*B*C$ , a multiple-line plot of estimated means at each level of  $A$ , with a separate line for each level of  $B$ , is produced for each of the three levels of  $C$  (three plots). Plots are produced with the value of the covariate  $X$  fixed at 1.0.

## **TEST Subcommand**

The `TEST` subcommand allows you to test a hypothesis term against a specified error term.

- `TEST` is valid only for univariate analyses. Multiple `TEST` subcommands are allowed, each executed independently.
- You must specify both the hypothesis term and the error term. There is no default.
- The hypothesis term is specified before the keyword `VS`. It must be a valid effect specified or implied on the `DESIGN` subcommand.
- The error term is specified after the keyword `VS`. You can specify either a linear combination or a value. The linear combination of effects takes the general form: `coefficient*effect +/- coefficient*effect ...`
- All effects in the linear combination must be specified or implied on the `DESIGN` subcommand. Effects specified or implied on `DESIGN` but not listed after `VS` are assumed to have a coefficient of 0.
- Duplicate effects are allowed. `UNIANOVA` adds coefficients associated with the same effect before performing the test. For example, the linear combination `5*A-0.9*B-A` will be combined to `4*A-0.9B`.
- A coefficient can be specified as a fraction with a positive denominator—for example, `1/3` or `-1/3`, but `1/-3` is invalid.
- If you specify a value for the error term, you must specify the degrees of freedom after the keyword `DF`. The degrees of freedom must be a positive real number. `DF` and the degrees of freedom are optional for a linear combination.

### **Example**

```
UNIANOVA DEP BY A B
  /TEST = A VS B + A*B
  /DESIGN = A, B, A*B.
```

- $A$  is tested against the pooled effect of  $B + A*B$ .

## LMATRIX Subcommand

The `LMATRIX` subcommand allows you to customize your hypotheses tests by specifying the **L** matrix (contrast coefficients matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ , where  $\mathbf{K} = \mathbf{0}$  if it is not specified on the `KMATRIX` subcommand. The vector **B** is the parameter vector in the linear model.

- The basic format for the `LMATRIX` subcommand is an optional label in quotation marks, an effect name or the keyword `ALL`, and a list of real numbers. There can be multiple effect names (or the keyword `ALL`) and number lists.
- The optional label is a string with a maximum length of 255 characters. Only one label can be specified.
- Only valid effects appearing or implied on the `DESIGN` subcommand can be specified on the `LMATRIX` subcommand.
- The length of the list of real numbers must be equal to the number of parameters (including the redundant ones) corresponding to that effect. For example, if the effect  $A*B$  takes up six columns in the design matrix, then the list after  $A*B$  must contain exactly six numbers.
- A number can be specified as a fraction with a positive denominator—for example,  $1/3$  or  $-1/3$ , but  $1/-3$  is invalid.
- A semicolon (;) indicates the end of a row in the **L** matrix.
- When `ALL` is specified, the length of the list that follows `ALL` is equal to the total number of parameters (including the redundant ones) in the model.
- Effects appearing or implied on the `DESIGN` subcommand must be explicitly specified here.
- Multiple `LMATRIX` subcommands are allowed. Each is treated independently.

### Example

```
UNIANOVA DEP BY A B
  /LMATRIX = "B1 vs B2 at A1"
    B 1 -1 0 A*B 1 -1 0 0 0 0 0 0
  /LMATRIX = "Effect A"
    A 1 0 -1
    A*B 1/3 1/3 1/3
        0 0 0
        -1/3 -1/3 -1/3;
    A 0 1 -1
    A*B 0 0 0
        1/3 1/3 1/3
        -1/3 -1/3 -1/3
  /LMATRIX = "B1 vs B2 at A2"
    ALL 0
        0 0 0
        1 -1 0
        0 0 0 1 -1 0 0 0 0
  /DESIGN = A, B, A*B.
```

Assume that factors  $A$  and  $B$  each have three levels. There are three `LMATRIX` subcommands; each is treated independently.

- **B1 versus B2 at A1.** In the first `LMATRIX` subcommand, the difference is tested between levels 1 and 2 of effect  $B$  when effect  $A$  is fixed at level 1. Since there are three levels each in effects  $A$  and  $B$ , the interaction effect  $A*B$  takes up nine columns in the design matrix.

- **Effect A.** In the second `LMATRIX` subcommand, effect *A* is tested. Since there are three levels in effect *A*, at most two independent contrasts can be formed; thus, there are two rows in the **L** matrix, which are separated by a semicolon (;). The first row tests the difference between levels 1 and 3 of effect *A*, while the second row tests the difference between levels 2 and 3 of effect *A*.
- **B1 versus B2 at A2.** In the last `LMATRIX` subcommand, the keyword `ALL` is used. The first 0 corresponds to the intercept effect; the next three zeros correspond to effect *A*.

## ***KMATRIX* Subcommand**

The `KMATRIX` subcommand allows you to customize your hypothesis tests by specifying the **K** matrix (contrast results matrix) in the general form of the linear hypothesis  $\mathbf{LB} = \mathbf{K}$ . The vector **B** is the parameter vector in the linear model.

- The default **K** matrix is a zero matrix; that is,  $\mathbf{LB} = \mathbf{0}$  is assumed.
- For the `KMATRIX` subcommand to be valid, at least one of the following subcommands must be specified: the `LMATRIX` subcommand or the `INTERCEPT = INCLUDE` subcommand.
- If `KMATRIX` is specified but `LMATRIX` is not specified, the `LMATRIX` is assumed to take the row vector corresponding to the intercept in the estimable function, provided the subcommand `INTERCEPT = INCLUDE` is specified. In this case, the **K** matrix can be only a scalar matrix.
- If `KMATRIX` and `LMATRIX` are specified, then the number of rows in the requested **K** and **L** matrices must be equal. If there are multiple `LMATRIX` subcommands, then all requested **L** matrices must have the same number of rows, and **K** must have the same number of rows as these **L** matrices.
- A semicolon (;) can be used to indicate the end of a row in the **K** matrix.
- If more than one `KMATRIX` subcommand is specified, only the last one is in effect.

### ***Example***

```
UNIANOVA DEP BY A B
  /LMATRIX = "Effect A"
            A 1 0 -1; A 1 -1 0
  /LMATRIX = "Effect B"
            B 1 0 -1; B 1 -1 0
  /KMATRIX = 0; 0
  /DESIGN = A B.
```

In this example, assume that factors *A* and *B* each have three levels.

- There are two `LMATRIX` subcommands; both have two rows.
- The first `LMATRIX` subcommand tests whether the effect of *A* is 0, while the second `LMATRIX` subcommand tests whether the effect of *B* is 0.
- The `KMATRIX` subcommand specifies that the **K** matrix also has two rows, each with value 0.

## ***CONTRAST* Subcommand**

`CONTRAST` specifies the type of contrast desired among the levels of a factor. For a factor with *k* levels or values, the contrast type determines the meaning of its *k*–1 degrees of freedom.

- Specify the factor name in parentheses following the subcommand `CONTRAST`.
- You can specify only one factor per `CONTRAST` subcommand, but you can enter multiple `CONTRAST` subcommands.
- After closing the parentheses, enter an equals sign followed by one of the contrast keywords.
- This subcommand creates an **L** matrix such that the columns corresponding to the factor match the contrast given. The other columns are adjusted so that the **L** matrix is estimable.

The following contrast types are available:

- DEVIATION**      *Deviations from the grand mean.* This is the default for factors. Each level of the factor except one is compared to the grand mean. One category (by default, the last) must be omitted so that the effects will be independent of one another. To omit a category other than the last, specify the number of the omitted category (which is not necessarily the same as its value) in parentheses after the keyword `DEVIATION`. For example,  
`UNIANOVA Y BY B /CONTRAST(B)=DEVIATION(1)`.  
 Suppose factor *B* has three levels, with values 2, 4, and 6. The specified contrast omits the first category, in which *B* has the value 2. Deviation contrasts are not orthogonal.
- POLYNOMIAL**      *Polynomial contrasts.* This is the default for within-subjects factors. The first degree of freedom contains the linear effect across the levels of the factor, the second contains the quadratic effect, and so on. In a balanced design, polynomial contrasts are orthogonal. By default, the levels are assumed to be equally spaced; you can specify unequal spacing by entering a metric consisting of one integer for each level of the factor in parentheses after the keyword `POLYNOMIAL`. (All metrics specified cannot be equal; thus, (1, 1, . . . 1) is not valid.) For example,  
`UNIANOVA RESPONSE BY STIMULUS /CONTRAST(STIMULUS) =  
 POLYNOMIAL(1, 2, 4)`.  
 Suppose that factor *STIMULUS* has three levels. The specified contrast indicates that the three levels of *STIMULUS* are actually in the proportion 1:2:4. The default metric is always (1, 2, . . . *k*), where *k* levels are involved. Only the relative differences between the terms of the metric matter; (1, 2, 4) is the same metric as (2, 3, 5) or (20, 30, 50) because, in each instance, the difference between the second and third numbers is twice the difference between the first and second.
- DIFFERENCE**      *Difference or reverse Helmert contrasts.* Each level of the factor except the first is compared to the mean of the previous levels. In a balanced design, difference contrasts are orthogonal.
- HELMERT**      *Helmert contrasts.* Each level of the factor except the last is compared to the mean of subsequent levels. In a balanced design, Helmert contrasts are orthogonal.
- SIMPLE**      *Each level of the factor except the last is compared to the last level.* To use a category other than the last as the omitted reference category, specify its number (which is not necessarily the same as its value) in parentheses following the keyword `SIMPLE`. For example,  
`UNIANOVA Y BY B /CONTRAST(B)=SIMPLE(1)`.  
 Suppose that factor *B* has three levels with values 2, 4, and 6. The specified contrast compares the other levels to the first level of *B*, in which *B* has the value 2. Simple contrasts are not orthogonal.
- REPEATED**      *Comparison of adjacent levels.* Each level of the factor except the first is compared to the previous level. Repeated contrasts are not orthogonal.

**SPECIAL** *A user-defined contrast.* Values specified after this keyword are stored in a matrix in column major order. For example, if factor *A* has three levels, then `CONTRAST(A)=SPECIAL(1 1 1 1 -1 0 0 1 -1)` produces the following contrast matrix:

```
1 1 0 1 -1 1 1 0 -1
```

**Orthogonal contrasts** are particularly useful. In a balanced design, contrasts are orthogonal if the sum of the coefficients in each contrast row is 0 and if, for any pair of contrast rows, the products of corresponding coefficients sum to 0. `DIFFERENCE`, `HELMERT`, and `POLYNOMIAL` contrasts always meet these criteria in balanced designs.

### Example

```
UNIANOVA DEP BY FAC
  /CONTRAST (FAC) =DIFFERENCE
  /DESIGN.
```

- Suppose that the factor *FAC* has five categories and therefore four degrees of freedom.
- `CONTRAST` requests `DIFFERENCE` contrasts, which compare each level (except the first) with the mean of the previous levels.

## POSTHOC Subcommand

`POSTHOC` allows you to produce multiple comparisons between means of a factor. These comparisons are usually not planned at the beginning of the study but are suggested by the data in the course of study.

- Post hoc tests are computed for the dependent variable. The alpha value used in the tests can be specified by using the keyword `ALPHA` on the `CRITERIA` subcommand. The default alpha value is 0.05. The confidence level for any confidence interval constructed is  $(1-\alpha) \times 100$ . The default confidence level is 95.
- Only factors appearing in the factor list are valid in this subcommand. Individual factors can be specified.
- You can specify one or more effects to be tested. Only fixed main effects appearing or implied on the `DESIGN` subcommand are valid test effects.
- Optionally, you can specify an effect defining the error term following the keyword `VS` after the test specification. The error effect can be any single effect in the design that is not the intercept or a main effect named on a `POSTHOC` subcommand.
- A variety of multiple comparison tests are available. Some tests are designed for detecting homogeneity subsets among the groups of means, some are designed for pairwise comparisons among all means, and some can be used for both purposes.
- For tests that are used for detecting homogeneity subsets of means, non-empty group means are sorted in ascending order. Means that are not significantly different are included together to form a homogeneity subset. The significance for each homogeneity subset of means is displayed. In a case where the numbers of valid cases are not equal in all groups, for most post hoc tests, the harmonic mean of the group sizes is used as the sample size in the calculation. For `QREGW` or `FREGW`, individual sample sizes are used.

- For tests that are used for pairwise comparisons, the display includes the difference between each pair of compared means, the confidence interval for the difference, and the significance. The sample sizes of the two groups being compared are used in the calculation.
- Output for tests specified on the `POSTHOC` subcommand are available according to their statistical purposes. The following table illustrates the statistical purpose of the post hoc tests:

Post Hoc Tests	Statistical Purpose	
	Homogeneity Subsets Detection	Pairwise Comparison and Confidence Interval
LSD		Yes
SIDAK		Yes
BONFERRONI		Yes
GH		Yes
T2		Yes
T3		Yes
C		Yes
DUNNETT		Yes*
DUNNETTL		Yes*
DUNNETTR		Yes*
SNK	Yes	
BTUKEY	Yes	
DUNCAN	Yes	
QREGW	Yes	
FREGW	Yes	
WALLER	Yes <sup>†</sup>	
TUKEY	Yes	Yes
SCHEFFE	Yes	Yes
GT2	Yes	Yes
GABRIEL	Yes	Yes

\* Only CIs for differences between test group means and control group means are given.

† No significance for Waller test is given.

- Tests that are designed for homogeneity subset detection display the detected homogeneity subsets and their corresponding significances.
- Tests that are designed for both homogeneity subset detection and pairwise comparison display both kinds of output.
- For the `DUNNETT`, `DUNNETTL`, and `DUNNETTR` keywords, only individual factors can be specified.
- The default reference category for `DUNNETT`, `DUNNETTL`, and `DUNNETTR` is the last category. An integer greater than 0 within parentheses can be used to specify a different reference category. For example, `POSTHOC = A (DUNNETT(2))` requests a `DUNNETT` test for factor *A*, using the second level of *A* as the reference category.
- The keywords `DUNCAN`, `DUNNETT`, `DUNNETTL`, and `DUNNETTR` must be spelled out in full; using the first three characters alone is not sufficient.

- If the `REGWT` subcommand is specified, weighted means are used in performing post hoc tests.
- Multiple `POSTHOC` subcommands are allowed. Each specification is executed independently so that you can test different effects against different error terms.

<b>SNK</b>	<i>Student-Newman-Keuls procedure based on the Studentized range test.</i>
<b>TUKEY</b>	<i>Tukey's honestly significant difference.</i> This test uses the Studentized range statistic to make all pairwise comparisons between groups.
<b>BTUKEY</b>	<i>Tukey's b.</i> Multiple comparison procedure based on the average of Studentized range tests.
<b>DUNCAN</b>	<i>Duncan's multiple comparison procedure based on the Studentized range test.</i>
<b>SCHEFFE</b>	<i>Scheffé's multiple comparison t test.</i>
<b>DUNNETT(refcat)</b>	<i>Dunnett's two-tailed t test.</i> Each level of the factor is compared to a reference category. A reference category can be specified in parentheses. The default reference category is the last category. This keyword must be spelled out in full.
<b>DUNNETTL(refcat)</b>	<i>Dunnett's one-tailed t test.</i> This test indicates whether the mean at any level (except the reference category) of the factor is <i>smaller</i> than that of the reference category. A reference category can be specified in parentheses. The default reference category is the last category. This keyword must be spelled out in full.
<b>DUNNETTR(refcat)</b>	<i>Dunnett's one-tailed t test.</i> This test indicates whether the mean at any level (except the reference category) of the factor is <i>larger</i> than that of the reference category. A reference category can be specified in parentheses. The default reference category is the last category. This keyword must be spelled out in full.
<b>BONFERRONI</b>	<i>Bonferroni t test.</i> This test is based on Student's <i>t</i> statistic and adjusts the observed significance level for the fact that multiple comparisons are made.
<b>LSD</b>	<i>Least significant difference t test.</i> Equivalent to multiple <i>t</i> tests between all pairs of groups. This test does not control the overall probability of rejecting the hypotheses that some pairs of means are different, while in fact they are equal.
<b>SIDAK</b>	<i>Sidak t test.</i> This test provides tighter bounds than the Bonferroni test.
<b>GT2</b>	<i>Hochberg's GT2.</i> Pairwise comparisons test based on the Studentized maximum modulus test. Unless the cell sizes are extremely unbalanced, this test is fairly robust even for unequal variances.
<b>GABRIEL</b>	<i>Gabriel's pairwise comparisons test based on the Studentized maximum modulus test.</i>
<b>FREGW</b>	<i>Ryan-Einot-Gabriel-Welsch's multiple stepdown procedure based on an F test.</i>
<b>QREGW</b>	<i>Ryan-Einot-Gabriel-Welsch's multiple stepdown procedure based on the Studentized range test.</i>
<b>T2</b>	<i>Tamhane's T2.</i> Tamhane's pairwise comparisons test based on a <i>t</i> test. This test can be applied in situations where the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.
<b>T3</b>	<i>Dunnett's T3.</i> Pairwise comparisons test based on the Studentized maximum modulus. This test is appropriate when the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.

<b>GH</b>	<i>Games and Howell's pairwise comparisons test based on the Studentized range test.</i> This test can be applied in situations where the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.
<b>C</b>	<i>Dunnett's C.</i> Pairwise comparisons based on the weighted average of Studentized ranges. This test can be applied in situations where the variances are unequal. This test is invalid when there are multiple factors in the model, and the keyword is ignored with a warning in such cases.
<b>WALLER(kratio)</b>	<i>Waller-Duncan t test.</i> This test uses a Bayesian approach. It is restricted to cases with equal sample sizes. For cases with unequal sample sizes, the harmonic mean of the sample size is used. The <i>k</i> -ratio is the Type 1/Type 2 error seriousness ratio. The default value is 100. You can specify an integer greater than 1 within parentheses.

## **EMMEANS Subcommand**

EMMEANS displays estimated marginal means of the dependent variable in the cells (with covariates held at their overall mean value) and their standard errors for the specified factors. Note that these are predicted, not observed, means. The estimated marginal means are calculated using a modified definition by Searle, Speed, and Milliken (1980).

- TABLES, followed by an option in parentheses, is required. COMPARE is optional; if specified, it must follow TABLES.
- Multiple EMMEANS subcommands are allowed. Each is treated independently.
- If identical EMMEANS subcommands are specified, only the last identical subcommand is in effect. EMMEANS subcommands that are redundant but not identical (for example, crossed factor combinations such as  $A*B$  and  $B*A$ ) are all processed.

<b>TABLES(option)</b>	<p><i>Table specification.</i> Valid options are the keyword OVERALL, factors appearing on the factor list, and crossed factors constructed of factors on the factor list. Crossed factors can be specified using an asterisk (*) or the keyword BY. All factors in a crossed factor specification must be unique.</p> <p>If OVERALL is specified, the estimated marginal means of the dependent variable are displayed, collapsing over factors.</p> <p>If a factor, or a crossing of factors, is specified on the TABLES keyword, UNIANOVA collapses over any other factors before computing the estimated marginal means for the dependent variable.</p> <p>The TABLES keyword may be followed by the WITH keyword and a parenthesized list of covariates and values. Tables are produced with the values of the covariates fixed at the specified values. The keyword MEAN causes the covariate to be fixed at its observed mean value.</p>
<b>COMPARE(factor) ADJ(method)</b>	<p><i>Main- or simple-main-effects omnibus tests and pairwise comparisons of the dependent variable.</i> This option gives the mean difference, standard error, significance, and confidence interval for each pair of levels for the effect specified in the TABLES command, as well as an omnibus test for that effect. If only one factor is specified on TABLES, COMPARE can be specified by itself; otherwise, the factor specification is required. In this case, levels of the specified factor are compared with each other for each level of the other factors in the interaction.</p>



The optional ADJ keyword allows you to apply an adjustment to the confidence intervals and significance values to account for multiple comparisons. Methods available are LSD (no adjustment), BONFERRONI, or SIDAK.

If OVERALL is specified on TABLES, COMPARE is invalid.

### Example

```
UNIANOVA DEP BY A B
  /EMMEANS = TABLES(A*B) WITH(X=1.0) COMPARE(A) ADJ(LSD)
  /DESIGN.
```

- The output of this analysis includes a pairwise comparisons table for the dependent variable *DEP*.
- Assume that *A* has three levels and *B* has two levels. The first level of *A* is compared with the second and third levels, the second level with the first and third levels, and the third level with the first and second levels, with the value of the covariate *X* fixed at 1.0. The pairwise comparison is repeated for the two levels of *B*.

## SAVE Subcommand

Use *SAVE* to add one or more residual or fit values to the active dataset.

- Specify one or more temporary variables, each followed by an optional new name in parentheses.
- *WPRED* and *WRESID* can be saved only if *REGWGT* has been specified.
- Specifying a temporary variable on this subcommand results in a variable being added to the active data file for each dependent variable.
- You can specify variable names for the temporary variables. These names must be unique, valid variable names.
- If new names are not specified, *UNIANOVA* generates a rootname using a shortened form of the temporary variable name with a suffix.
- If more than one *SAVE* subcommand is specified, only the last one is in effect.

<b>PRED</b>	<i>Unstandardized predicted values.</i>
<b>WPRED</b>	<i>Weighted unstandardized predicted values.</i> Available only if <i>REGWGT</i> has been specified.
<b>RESID</b>	<i>Unstandardized residuals.</i>
<b>WRESID</b>	<i>Weighted unstandardized residuals.</i> Available only if <i>REGWGT</i> has been specified.
<b>DRESID</b>	<i>Deleted residuals.</i>
<b>ZRESID</b>	<i>Standardized residuals.</i>
<b>SRESID</b>	<i>Studentized residuals.</i>
<b>SEPREP</b>	<i>Standard errors of predicted value.</i>
<b>COOK</b>	<i>Cook's distances.</i>
<b>LEVER</b>	<i>Uncentered leverage values.</i>

## OUTFILE Subcommand

The OUTFILE subcommand writes IBM® SPSS® Statistics data files that can be used in other procedures.

- You must specify a keyword on OUTFILE. There is no default.
- You must specify a quoted file specification or a previously declared dataset name (DATASET DECLARE command), enclosed in quotes. The asterisk (\*) is not allowed.
- If you specify more than one keyword, a different file specification is required for each.
- If more than one OUTFILE subcommand is specified, only the last one is in effect.
- For COVB or CORB, the output will contain, in addition to the covariance or correlation matrix, three rows for each dependent variable: a row of parameter estimates, a row of residual degrees of freedom, and a row of significance values for the *t* statistics corresponding to the parameter estimates. All statistics are displayed separately by split.

**COVB** ('savfile'|'dataset')

*Writes the parameter covariance matrix.*

**CORB** ('savfile'|'dataset')

*Writes the parameter correlation matrix.*

**EFFECT** ('savfile'|'dataset')

*Writes the statistics from the between-subjects ANOVA table.*

**DESIGN** ('savfile'|'dataset')

*Writes the design matrix. The number of rows equals the number of cases, and the number of columns equals the number of parameters. The variable names are DES\_1, DES\_2, ..., DES\_p, where p is the number of the parameters.*

## DESIGN Subcommand

DESIGN specifies the effects included in a specific model. The cells in a design are defined by all of the possible combinations of levels of the factors in that design. The number of cells equals the product of the number of levels of all the factors. A design is *balanced* if each cell contains the same number of cases. UNIANOVA can analyze both balanced and unbalanced designs.

- Specify a list of terms to be included in the model, separated by spaces or commas.
- The default design, if the DESIGN subcommand is omitted or is specified by itself, is a design consisting of the following terms in order: the intercept term (if INTERCEPT=INCLUDE is specified), next the covariates given in the covariate list, and then the full factorial model defined by all factors on the factor list and excluding the intercept.
- To include a term for the main effect of a factor, enter the name of the factor on the DESIGN subcommand.
- To include the intercept term in the design, use the keyword INTERCEPT on the DESIGN subcommand. If INTERCEPT is specified on the DESIGN subcommand, the subcommand INTERCEPT=EXCLUDE is overridden.
- To include a term for an interaction between factors, use the keyword BY or the asterisk (\*) to join the factors involved in the interaction. For example, *A\*B* means a two-way interaction effect of *A* and *B*, where *A* and *B* are factors. *A\*A* is not allowed because factors inside an interaction effect must be distinct.

- To include a term for nesting one effect within another, use the keyword `WITHIN` or a pair of parentheses on the `DESIGN` subcommand. For example,  $A(B)$  means that  $A$  is nested within  $B$ . The expression  $A(B)$  is equivalent to the expression  $A$  *WITHIN*  $B$ . When more than one pair of parentheses is present, each pair of parentheses must be enclosed or nested within another pair of parentheses. Thus,  $A(B)(C)$  is not valid.
- Multiple nesting is allowed. For example,  $A(B(C))$  means that  $B$  is nested within  $C$ , and  $A$  is nested within  $B(C)$ .
- Interactions between nested effects are not valid. For example, neither  $A(C)*B(C)$  nor  $A(C)*B(D)$  is valid.
- To include a covariate term in the design, enter the name of the covariate on the `DESIGN` subcommand.
- Covariates can be connected, but not nested, through the `*` operator to form another covariate effect. Therefore, interactions among covariates such as  $X1*X1$  and  $X1*X2$  are valid, but not  $X1(X2)$ . Using covariate effects such as  $X1*X1$ ,  $X1*X1*X1$ ,  $X1*X2$ , and  $X1*X1*X2*X2$  makes fitting a polynomial regression model easy in UNIANOVA.
- Factor and covariate effects can be connected only by the `*` operator. Suppose  $A$  and  $B$  are factors, and  $X1$  and  $X2$  are covariates. Examples of valid factor-by-covariate interaction effects are  $A*X1$ ,  $A*B*X1$ ,  $X1*A(B)$ ,  $A*X1*X1$ , and  $B*X1*X2$ .
- If more than one `DESIGN` subcommand is specified, only the last one is in effect.

### **Example**

```
UNIANOVA Y BY A B C WITH X
  /DESIGN A B(A) X*A.
```

- In this example, the design consists of a main effect  $A$ , a nested effect  $B$  within  $A$ , and an interaction effect of a covariate  $X$  with a factor  $A$ .

# UPDATE

```
UPDATE FILE={master file}
           {*          }

[/RENAME=(old varnames=new varnames)...]

[/IN=varname]

/FILE={transaction file1}
      {*          }

[/FILE=transaction file2]

/BY key variables

[/MAP]

[/KEEP={ALL**  }] [/DROP=varlist]
      {varlist}
```

\*\*Default if the subcommand is omitted.

## Example

```
UPDATE FILE='/data/maillist.sav'
/FILE='/data/newlist.sav'
/BY=ID.
```

## Overview

UPDATE replaces values in a master file with updated values recorded in one or more files called transaction files. Cases in the master file and transaction file are matched according to a key variable.

The master file and the transaction files must be IBM® SPSS® Statistics data files or datasets available in the current session, including the active dataset. UPDATE replaces values and creates a new active dataset, which replaces the original active dataset.

UPDATE is designed to update values of existing variables for existing cases. Use MATCH FILES to add new variables to a data file and ADD FILES to add new cases.

## Options

**Variable Selection.** You can specify which variables from each input file are included in the new active dataset using the DROP and KEEP subcommands.

**Variable Names.** You can rename variables in each input file before combining the files using the RENAME subcommand. This permits you to combine variables that are the same but whose names differ in different input files, or to separate variables that are different but have the same name.

**Variable Flag.** You can create a variable that indicates whether a case came from a particular input file using IN. You can use the FIRST or LAST subcommand to create a variable that flags the first or last case of a group of cases with the same value for the key variable.

**Variable Map.** You can request a map showing all variables in the new active dataset, their order, and the input files from which they came using the `MAP` subcommand.

### ***Basic Specification***

The basic specification is two or more `FILE` subcommands and a `BY` subcommand.

- The first `FILE` subcommand must specify the master file. All other `FILE` subcommands identify the transaction files.
- `BY` specifies the key variables.
- All files must be sorted in ascending order by the key variables.
- By default, all variables from all input files are included in the new active dataset.

### ***Subcommand Order***

- The master file must be specified first.
- `RENAME` and `IN` must immediately follow the `FILE` subcommand to which they apply.
- `BY` must follow the `FILE` subcommands and any associated `RENAME` and `IN` subcommands.
- `MAP`, `DROP`, and `KEEP` must be specified after all `FILE` and `RENAME` subcommands.

### ***Syntax Rules***

- `BY` can be specified only once. However, multiple variables can be specified on `BY`. All files must be sorted in ascending order by the key variables named on `BY`.
- The master file cannot contain duplicate values for the key variables.
- `RENAME` can be repeated after each `FILE` subcommand and applies only to variables in the file named on the immediately preceding `FILE` subcommand.
- `MAP` can be repeated as often as needed.

### ***Operations***

- `UPDATE` reads all input files named on `FILE` and builds a new active dataset. The new active dataset is built when the data are read by one of the procedure commands or the `EXECUTE`, `SAVE`, or `SORT CASES` command.
- If the current active dataset is included and is specified with an asterisk (`FILE=*`), the new merged dataset replaces the active dataset. If that dataset is a named dataset, the merged dataset retains that name. If the current active dataset is not included or is specified by name (for example, `FILE=Dataset1`), a new unnamed, merged dataset is created, and it becomes the active dataset. For information on naming datasets, see [DATASET NAME](#).
- The new active dataset contains complete dictionary information from the input files, including variable names, labels, print and write formats, and missing-value indicators. The new active dataset also contains the documents from each input file, unless the `DROP DOCUMENTS` command is used.
- `UPDATE` copies all variables in order from the master file, then all variables in order from the first transaction file, then all variables in order from the second transaction file, and so on.

- Cases are updated when they are matched on the `BY` variable(s). If the master and transaction files contain common variables for matched cases, the values for those variables are taken from the transaction file, provided that the values are not missing or blanks. Missing or blank values in the transaction files are not used to update values in the master file.
- When `UPDATE` encounters duplicate keys within a transaction file, it applies each transaction sequentially to that case to produce one case per key value in the resulting file. If more than one transaction file is specified, the value for a variable comes from the last transaction file with a nonmissing value for that variable.
- Variables that are in the transaction files but not in the master file are added to the master file. Cases that do not contain those variables are assigned the system-missing value (for numerics) or blanks (for strings).
- Cases that are in the transaction files but not in the master file are added to the master file and are interleaved according to their values for the key variables.
- If the active dataset is named as an input file, any `N` and `SAMPLE` commands that have been specified are applied to the active dataset before files are combined.
- The `TEMPORARY` command cannot be in effect if the active dataset is used as an input file.

### **Limitations**

- A maximum of one `BY` subcommand. However, `BY` can specify multiple variables.

## **Examples**

### **Basic Example**

```
UPDATE FILE='/data/mailist.sav'
  /FILE='/data/newlist.sav'
  /BY=ID.
```

- *mailist.sav* is specified as the master file. *newlist.sav* is the transaction file. *ID* is the key variable.
- Both *mailist.sav* and *newlist.sav* must be sorted in ascending order of *ID*.
- If *newlist.sav* has cases or nonmissing variables that are not in *mailist.sav*, the new cases or variables are added to the resulting file.

### **Using Multiple Key Variables**

```
SORT CASES BY LOCATN DEPT.
UPDATE FILE='/data/master.sav' /FILE=* /BY LOCATN DEPT
  /KEEP AVGHOUR AVGRAISE LOCATN DEPT SEX HOURLY RAISE /MAP.
SAVE OUTFILE='/data/personnel.sav'.
```

- `SORT CASES` sorts the active dataset in ascending order of the variables to be named as key variables on `UPDATE`.
- `UPDATE` specifies *master.sav* as the master file and the sorted active dataset as the transaction file. The file *master.sav* must also be sorted by *LOCATN* and *DEPT*.
- `BY` specifies the key variables *LOCATN* and *DEPT*.

- **KEEP** specifies the subset and order of variables to be retained in the resulting file.
- **MAP** provides a list of the variables in the resulting file and the two input files.
- **SAVE** saves the resulting file as a data file in IBM® SPSS® Statistics format.

## ***FILE Subcommand***

**FILE** identifies each input file. At least two **FILE** subcommands are required on **UPDATE**: one specifies the master file and the other a transaction file. A separate **FILE** subcommand must be used to specify each transaction file.

- The first **FILE** subcommand must specify the master file.
- An asterisk on **FILE** refers to the active dataset.
- The files must be IBM® SPSS® Statistics data files or datasets available in the current session. File specifications should be enclosed in quotes.
- All files must be sorted in ascending order according to the variables specified on **BY**.
- The master file cannot contain duplicate values for the key variables. However, transaction files can and often do contain cases with duplicate keys (see “Operations”).

## ***Text Data Files***

You can update data from text data files — or from any data format that IBM® SPSS® Statistics can read — using defined datasets.

### ***Example***

```
DATA LIST FILE='/data/update.txt'
  ID 1-3 NAME 5-17 (A) ADDRESS 19-28 (A) ZIP 30-34.
SORT CASES BY ID.
DATASET NAME updatefile.
GET FILE='/maillist.sav'.
SORT CASES BY ID.
DATASET NAME=masterfile.
UPDATE FILE='masterfile'
  /RENAME=(STREET=ADDRESS)
  /FILE='updatefile'
  /BY=ID /MAP.
```

## ***BY Subcommand***

**BY** specifies one or more identification, or key, variables that are used to match cases between files.

- **BY** must follow the **FILE** subcommands and any associated **RENAME** and **IN** subcommands.
- **BY** specifies the names of one or more key variables. The key variables must exist in all input files and have the same names in all of the files. The key variables can be string variables (long strings are allowed).
- All input files must be sorted in ascending order of the key variables. If necessary, use **SORT CASES** before **UPDATE**.

- Missing values for key variables are handled like any other values.
- The key variables in the master file must identify unique cases. If duplicate cases are found, the program issues an error and UPDATE is not executed. The system-missing value is treated as one single value.

## **RENAME Subcommand**

RENAME renames variables on the input files *before* they are processed by UPDATE. RENAME must follow the FILE subcommand that contains the variables to be renamed.

- RENAME applies only to the immediately preceding FILE subcommand. To rename variables from more than one input file, specify a RENAME subcommand after each FILE subcommand.
- Specifications for RENAME consist of a left parenthesis, a list of old variable names, an equals sign, a list of new variable names, and a right parenthesis. The two variable lists must name or imply the same number of variables. If only one variable is renamed, the parentheses are optional.
- More than one rename specification can be specified on a single RENAME subcommand, each enclosed in parentheses.
- The TO keyword can be used to refer to consecutive variables in the file and to generate new variable names.
- RENAME takes effect immediately. Any KEEP and DROP subcommands entered prior to a RENAME must use the old names, while KEEP and DROP subcommands entered after a RENAME must use the new names.
- All specifications within a single set of parentheses take effect simultaneously. For example, the specification RENAME (A, B = B, A) swaps the names of the two variables.
- Variables cannot be renamed to scratch variables.

### **Example**

```
UPDATE FILE='/data/master.sav' /FILE=CLIENTS  
  /RENAME=(TEL_NO, ID_NO = PHONE, ID)  
  /BY ID.
```

- UPDATE updates the master phone list by using current information from the file *CLIENTS*.
- Two variables on *CLIENTS* are renamed prior to the match. *TEL\_NO* is renamed *PHONE* to match the name used for phone numbers in the master file. *ID\_NO* is renamed *ID* so that it will have the same name as the identification variable in the master file and can be used on the BY subcommand.
- The old variable names are listed before the equals sign, and the new variable names are listed in the same order after the equals sign. The parentheses are required.
- The BY subcommand matches cases according to client ID numbers.

## **DROP and KEEP Subcommands**

DROP and KEEP are used to include a subset of variables in the resulting file. DROP specifies a set of variables to exclude, and KEEP specifies a set of variables to retain.



- DROP and KEEP do not affect the input files on disk.
- DROP and KEEP must follow all FILE and RENAME subcommands.
- DROP and KEEP must specify one or more variables. If RENAME is used to rename variables, specify the new names on DROP and KEEP.
- DROP cannot be used with variables created by the IN subcommand.
- The keyword ALL can be specified on KEEP. ALL must be the last specification on KEEP, and it refers to all variables not previously named on KEEP.
- KEEP can be used to change the order of variables in the resulting file. With KEEP, variables are kept in the order in which they are listed on the subcommand. If a variable is named more than once on KEEP, only the first mention of the variable is in effect; all subsequent references to that variable name are ignored.
- Multiple DROP and KEEP subcommands are allowed. Specifying a variable that is not in the active dataset or that has been dropped because of a previous DROP or KEEP subcommand results in an error and the UPDATE command is not executed.

### Example

```
UPDATE FILE='/data/maillist.sav' /FILE='/data/newlist.sav' /RENAME=(STREET=ADDRESS)
  /BY ID /KEEP=NAME ADDRESS CITY STATE ZIP ID.
```

- KEEP specifies the variables to keep in the result file. The variables are stored in the order specified on KEEP.

## IN Subcommand

IN creates a new variable in the resulting file that indicates whether a case came from the input file named on the preceding FILE subcommand. IN applies only to the file specified on the immediately preceding FILE subcommand.

- IN has only one specification, the name of the flag variable.
- The variable created by IN has the value 1 for every case that came from the associated input file and the value 0 if the case came from a different input file.
- Variables created by IN are automatically attached to the end of the resulting file and cannot be dropped.

### Example

```
UPDATE FILE=WEEK10 /FILE=WEEK11 /IN=INWEEK11 /BY=EMPID.
```

- IN creates the variable *INWEEK11*, which has the value 1 for all cases in the resulting file that came from the input file *WEEK11* and the value 0 for those cases that were not in file *WEEK11*.

## MAP Subcommand

MAP produces a list of the variables in the new active dataset and the file or files from which they came. Variables are listed in the order in which they appear in the resulting file. MAP has no specifications and must be placed after all FILE, RENAME, and IN subcommands.

- Multiple `MAP` subcommands can be used. Each `MAP` shows the current status of the active dataset and reflects only the subcommands that precede the `MAP` subcommand.
- To obtain a map of the resulting file in its final state, specify `MAP` last.
- If a variable is renamed, its original and new names are listed. Variables created by `IN` are not included in the map, since they are automatically attached to the end of the file and cannot be dropped.

# USE

```
USE [{start date      }] [THRU [{end date      }]]  
    {start case number} {end case number}  
    {FIRST             } {LAST              }  
  
[ALL]  
  
/PERMANENT
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
USE Y 1960.
```

## **Overview**

USE designates a range of observations to be used with time series procedures. Cases outside the defined range are filtered but not deleted from the active dataset unless the /PERMANENT subcommand is specified.

### **Basic Specification**

The basic specification is either the start of the range, the end of the range, or both. You can also simply specify the keyword THRU or ALL.

- The default start is the first observation in the file, and the default end is the last observation.
- The keyword THRU is required if the end of the range is specified.
- The keyword ALL defines a USE range starting with the first observation and ending with the last observation in the series. It can be specified to restore processing to the entire series.
- The keyword THRU by itself is the same as specifying the keyword ALL.

### **Operations**

- USE is ignored by the utility procedures CREATE and RMV. These procedures process all of the available data.
- The DATE command turns off all existing USE and PREDICT specifications.
- FILTER and USE are mutually exclusive. USE automatically turns off any previous FILTER command, and FILTER automatically turns off any previous USE command.
- The SORT command automatically turns off any previous USE command specification.
- USE remains in effect in a session until it is changed by another USE command or until a new DATE, FILTER, or SORT command is issued.

**Limitations**

- A maximum of one range (one start and/or one end) can be specified.

**Syntax Rules**

- The start and end can be specified as either `DATE` specifications or case (observation) numbers.
- `DATE` specifications and case numbers cannot be mixed on a `USE` command.
- Any observation within the file can be used as the start or end as long as the starting observation comes before the end observation.

**DATE Specifications**

- A `DATE` specification consists of `DATE` keywords and values (see `DATE`). These specifications must correspond to existing `DATE` variables.
- If more than one `DATE` variable exists, the highest-order one must be used in the specification.
- Values on the keyword `YEAR` must have the same format (two or four digits) as the `YEAR` specifications on the `DATE` command.

**Case Specifications**

- The case number specification is the sequence number of the case (observation) as it is read by the program.

**Keywords FIRST and LAST**

- The start can also be specified with the keyword `FIRST`, and the end, with the keyword `LAST`. These keywords designate the first and last cases in the file, respectively.
- The keywords `FIRST` and `LAST` can be used along with either `DATE` or case specifications.

**PERMANENT Subcommand**

The `PERMANENT` subcommand deletes cases outside the defined range from the active dataset.

**Examples**

```
USE ALL.
```

- This command includes all observations in the file in the `USE` range.
- This specification is the same as `USE THRU` or `USE FIRST THRU LAST`.

```
USE Y 1960.
```

- This command selects observations starting with a `YEAR_` value of 1960 through the last observation in the file. It is equivalent to `USE Y 1960 THRU LAST`.

```
USE THRU D 5.
```

- This command selects all cases from the first case in the file to the last one with a *DAY\_* value of 5. It is equivalent to `USE FIRST THRU D 5.`

`USE THRU 5.`

- This command selects cases starting with the first case and ending with the fifth case.

`USE Y 1955 M 6 THRU Y 1960 M 6.`

- This selects cases from June, 1955, through June, 1960.

`USE W 16 D 3 THRU W 48 D 3.`

- This example selects cases from day 3 of week 16 through day 3 of week 48.

`USE CYCLE 2 OBS 4 THRU CYCLE 2 OBS 17.`

- This example selects observations 4 through 17 of the second cycle.

# VALIDATEDATA

VALIDATEDATA is available in the Data Preparation option.

```
VALIDATEDATA      [VARIABLES=varlist]
                  [ID=varlist]
                  [CROSSVARRULES=rulelist]
[/VARCHHECKS     [STATUS={ON**}]
                  {OFF }
                  [PCTMISSING={70** }
                  {number }
                  [PCTEQUAL={95** }
                  {number }
                  [PCTUNEQUAL={90** }
                  {number }
                  [CV={0.001**}
                  {number }
                  [STDDEV={0** }
                  {number }
[/IDCHECKS       [INCOMPLETE**] [DUPLICATE**] [NONE]]
[/CASECHECKS     [REPORTEMPTY={YES**}] [SCOPE={ANALYSISVARS}]
                  {NO } {ALLVARS }
[/RULESUMMARIES [BYVARIABLE**] [BYRULE] [NONE]]
[/CASEREPORT     DISPLAY={YES**}
                  {NO }
                  [MINVIOLATIONS={1** }
                  {integer}
                  [CASELIMIT={FIRSTN({500** }
                  {integer}
                  {NONE }
[/SAVE           [EMPTYCASE[ (varname)]]
                  [INCOMPLETEID[ (varname)]]
                  [DUPLICATEID[ (varname)]]
                  [RULEVIOLATIONS[ (varname)]]]
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Release History

Release 14.0

- Command introduced.

## Example

```
VALIDATEDATA VARIABLES=ALL.
```

## Overview

The VALIDATEDATA procedure identifies suspicious and invalid cases, variables, and data values in the active dataset.

The procedure can summarize two types of validation rules. Single-variable rules consist of a fixed set of checks that are applied to individual data values, such as range checks. Cross-variable rules are user-specified rules that typically examine combinations of data values for two or more variables.

### ***Options***

**Analysis Variables.** The procedure can identify variables that have a high proportion of missing values as well as variables that are constant (or nearly so). You can set the maximum acceptable percentage of missing values as well as thresholds for considering categorical and scale variables as constants.

**Case Identifiers.** The procedure can identify incomplete and duplicate case identifiers. A single variable or a combination of variables can be treated as the identifier for a case.

**Cases.** The procedure can identify empty cases. A case can be regarded as empty if all analysis variables are blank or missing or all non-ID variables are blank or missing.

**Single-Variable Rules.** Single-variable validation rules can be summarized by variable and rule. Single-variable rules are defined using the `DATAFILE ATTRIBUTE` command and are linked to analysis variables using the `VARIABLE ATTRIBUTE` command.

**Cross-Variable Rules.** The procedure can summarize violations of cross-variable validation rules. Cross-variable rules are defined using the `DATAFILE ATTRIBUTE` command.

**Saved Variables.** You can save variables that identify suspicious cases and values in the active dataset.

### ***Basic Specification***

- The minimum specification is a list of analysis variables, case identifier variables, or cross-variable rules.
- By default, if you specify analysis variables, the procedure reports all analysis variables that have a high proportion of missing values, as well as analysis variables that are constant or nearly so. If single-variable validation rules are defined for analysis variables, the procedure reports rule violations for each variable.
- By default, if case identifier variables are specified, the procedure reports all incomplete and duplicate case identifiers.
- If you specify cross-variable rules, the procedure reports the total number of cases that violated each rule.
- If you specify single-variable or cross-variable rules, a casewise report is shown that lists the first 500 cases that violated at least one validation rule.
- Empty cases are identified by default.

### ***Syntax Rules***

- Each subcommand is global and can be used only once.
- Subcommands can be used in any order.
- An error occurs if a keyword or attribute is specified more than once within a subcommand.

- Equals signs and slashes shown in the syntax chart are required.
- Subcommand names and keywords must be spelled out in full.
- Empty subcommands generate a procedure error.

### ***Operations***

- The procedure honors `SPLIT FILE` specifications. Split variables are filtered out of all variable lists.
- The procedure treats user- and system-missing values as missing values.
- An error occurs if no procedure output is requested.

Note: Since measurement level can affect the results, if any variables (fields) have an unknown measurement level, an initial data pass will be performed to determine default measurement level for any variables with an unknown measurement level. For information on the criteria used to determine default measurement level, see [SET SCALEMIN](#).

### ***Limitations***

- The weight variable, if specified, is ignored with a warning.

### ***Validation Rules***

- `VALIDATEDATA` ignores invalid rules and links with a warning.
- If an analysis variable is linked to multiple rules, `VALIDATEDATA` applies the rules independently; it does not check for conflicts or redundancies among the constituent rules.
- A rule outcome variable must be associated with each rule that indicates which cases violated the rule. Outcome variables are created outside the `VALIDATEDATA` procedure and should be coded such that 1 indicates an invalid value or combination of values and 0 indicates a valid value.
- An error occurs if the same outcome variable is referenced by two or more rules.
- Values of all outcome variables are assumed to be current with respect to the active dataset.

## ***Examples***

### ***Default Analysis Variable Validation***

```
VALIDATEDATA VARIABLES=ALL.
```

- The only specification is a list of analysis variables (all variables in the active dataset).
- The procedure reports all variables with a high proportion of missing values, as well as variables that are constant or nearly so.
- In addition, empty cases are identified.

### ***Default Case Identifier Variable Validation***

```
VALIDATEDATA ID=Firstname Lastname.
```



- Two case identifier variables are specified.
- By default, the procedure reports cases having incomplete or duplicate case identifiers. The combination of *Firstname* and *Lastname* values is treated as a case identifier.
- In addition, empty cases are identified.

### **Non-Default Analysis and Case Identifier Variable Validation**

```
VALIDATEDATA VARIABLES=Satis1 Satis2 Satis3
                ID=Firstname Lastname
/VARCHECKS STATUS=ON PCTMISSING=30
/IDCHECKS DUPLICATE
/SAVE EMPTYCASE.
```

- Three ordinal satisfaction variables (*Satis1*, *Satis2*, *Satis3*) are specified as analysis variables.
- ID specifies two case identifier variables, *Firstname* and *Lastname*.
- VARCHECKS reports analysis variables with more than 30% missing values. By default, the procedure also reports categorical analysis variables if more than 95% of their cases fall into a single category.
- IDCHECKS reports cases that have duplicate case identifiers and turns off the default check for incomplete case identifiers.
- Empty cases are identified by default.
- SAVE saves a variable to the active dataset that indicates which cases are empty.

### **Using Single-Variable Validation Rules**

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=$VD.SRule[1] ("Label='Likert 1 to 5',"+
                        "Description='Likert 5-point scale',"+
                        "Type='Numeric',"+
                        "Domain='Range',"+
                        "FlagUserMissing='No',"+
                        "FlagSystemMissing='No',"+
                        "FlagBlank='No',"+
                        "Minimum='1',"+
                        "Maximum='5',"+
                        "FlagNoninteger='Yes'").

COMPUTE Likert1to5_Satis_=NOT(Satis LE 5 AND Satis GE 1 AND
                             Satis=TRUNC(Satis)).

VARIABLE ATTRIBUTE
  VARIABLES= Likert1to5_Satis_
  ATTRIBUTE=$VD.RuleOutcomeVar("Yes").

VARIABLE ATTRIBUTE
  VARIABLES=Satis
  ATTRIBUTE=$VD.SRuleRef[1] ("Rule='$VD.SRule[1]',"+
                             "OutcomeVar='Likert1to5_Satis_').

VALIDATEDATA VARIABLES=Satis
/CASEREPORT DISPLAY=YES CASELIMIT=NONE
/SAVE RULEVIOLATIONS.
```

- DATAFILE ATTRIBUTE defines a single-variable validation rule named *\$VD.SRule[1]*. The rule flags any value that is not an integer within the range 1 to 5.

- The first `VARIABLE ATTRIBUTE` command links the rule to the variable *Satis*. It also references an outcome variable where 1 indicates an invalid value for the variable *Satis*.
- `COMPUTE` creates the rule outcome variable.
- The second `VARIABLE ATTRIBUTE` command marks *Likert1to5\_Satis\_* as an outcome variable in the data dictionary.
- `VALIDATEDATA` specifies *Satis* as an analysis variable. In addition to performing default checks for constants, variables with a large proportion of missing values, and empty cases, the procedure summarizes violations of single-variable validation rules defined for *Satis*.
- `CASEREPORT` reports all cases having a rule violation.
- `SAVE` saves a variable to the active dataset that indicates the total number of validation rule violations per case.

### **Using Cross-Variable Validation Rules**

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=$VD.CRule[1] ("Label='Pregnant Male',"+
    "Expression='Sex ='Male' AND Pregnant ''Yes''", "+
    "OutcomeVar='PregnantMale_'").

COMPUTE PregnantMale_= Sex ='Male' AND Pregnant = 'Yes'.
VARIABLE ATTRIBUTE
  VARIABLES=PregnantMale_
  ATTRIBUTE=$VD.RuleOutcomeVar("Yes").

VALIDATEDATA CROSSVARRULES=$VD.CRule[1]
  /CASECHECKS REPORTEEMPTY=NO.
```

- `DATAFILE ATTRIBUTE` defines the cross-variable rule *\$VD.CRule[1]*.
- `COMPUTE` creates the outcome variable *PregnantMale\_* referenced by the cross-variable rule. For *PregnantMale\_*, values of 1 identify cases containing males coded as being pregnant.
- `VARIABLE ATTRIBUTE` marks *PregnantMale\_* as an outcome variable in the data dictionary.
- `VALIDATEDATA` specifies *\$VD.CRule[1]* as a cross-variable rule to be summarized. The procedure reports the total number of pregnant males in the active dataset.
- `CASECHECKS` turns off the default check for empty cases.
- By default, the first 500 cases that violated at least one validation rule are listed.

## **Variable Lists**

You must specify `VARIABLES`, `ID`, or `CROSSVARRULES`.

### **VARIABLES Keyword**

The `VARIABLES` list specifies analysis variables to be validated.

- Specify one or more analysis variables. `ALL` and `TO` can be used to refer to all variables or a range of variables, respectively.
- Repeated instances of the same variable are filtered out of the list.
- Rule outcome variables are filtered out of the list.

**ID Keyword**

The ID list specifies case ID variables to be validated. ID variables are also used to label casewise output.

- Specify one or more variables. If two or more ID variables are specified, the combination of their values is treated as a case identifier.
- Repeated instances of the same variable are filtered out of the list.
- Rule outcome variables are filtered out of the list.

**CROSSVARRULES Keyword**

The CROSSVARRULES keyword specifies cross-variable rules to be summarized.

- Specify one or more cross-variable rules of the form  $\$VD.CRrule[n]$ , where  $n$  is a positive integer. Alternatively, you can provide a list of integers, where 1 means  $\$VD.CRrule[1]$ , 2 means  $\$VD.CRrule[2]$ , etc.
- The list accepts the keyword ALL, which signifies all cross-variable rules recorded in the data dictionary.
- Repeated instances of the same rule are filtered out of the list.
- A warning occurs if a specified rule does not exist. The rule is ignored.

**VARCHECKS Subcommand**

The VARCHHECKS subcommand specifies checks to be performed on analysis variables. The subcommand is ignored if no analysis variables are defined.

<b>STATUS</b>	<i>Perform variable checks.</i> By default, variable checks are performed. To turn off variable checks, specify STATUS=NO; any other VARCHHECKS keywords are then ignored.
<b>PCTMISSING</b>	<i>Maximum percentage of missing values.</i> Reports analysis variables with a percentage of missing values greater than the specified value. The specified value must be a positive number less than or equal to 100. The default value is 70.
<b>PCTEQUAL</b>	<i>Maximum percentage of cases representing a single category.</i> Reports categorical analysis variables with a percentage of cases representing a single nonmissing category greater than the specified value. The specified value must be a positive number less than or equal to 100. The default value is 95. The percentage is based on cases with nonmissing values of the variable. PCTEQUAL is ignored if no categorical analysis variables are specified.
<b>PCTUNEQUAL</b>	<i>Percentage of categories containing only one case in a categorical variable.</i> If the percentage of an analysis variable's categories containing only one case is greater than the specified value, the variable is reported. The specified value must be a positive number less than or equal to 100. The default value is 90. PCTUNEQUAL is ignored if no categorical analysis variables are specified.
<b>CV</b>	<i>Minimum absolute coefficient of variation.</i> A variable's coefficient of variation is defined as its standard deviation divided by its mean. If the absolute value of the coefficient of variation is less than the specified value, the variable is reported.

This setting applies only to continuous predictors and only if the mean is non-zero. The specified value must be a non-negative number. Specifying 0 turns off the coefficient of variation check. The default value is 0.001. CV is ignored if no scale analysis variables are specified.

**STDDEV** *Minimum standard deviation.* Reports variables whose standard deviation is less than the specified value. This setting applies only to scale analysis variables. The specified value must be a non-negative number. The default value is 0, which, in effect, turns off the standard deviation check. STDDEV is ignored if no scale analysis variables are specified.

## **IDCHECKS Subcommand**

The IDCHECKS subcommand specifies checks to be performed on case identifiers. The subcommand is ignored if no ID variables are specified.

**INCOMPLETE** *Incomplete case identifiers.* Reports cases with incomplete case identifiers. For a particular case, an identifier is considered incomplete if the value of any ID variable is blank or missing. Default.

**DUPLICATE** *Duplicate case identifiers.* Reports cases with duplicate case identifiers. Incomplete identifiers are excluded from the set of possible duplicates. Default.

**NONE** *Case identifiers are not validated.*

## **CASECHECKS Subcommand**

The CASECHECKS subcommand specifies checks to be performed on cases in the active dataset.

### **REPORTEMPTY Keyword**

Empty cases are identified by default. To turn off the check for empty cases, specify REPORTEMPTY=NO.

### **SCOPE Keyword**

The SCOPE keyword controls which variables are considered when the procedure determines whether a case is empty. SCOPE is ignored if REPORTEMPTY=NO.

**ANALYSISVARS** *Procedure considers analysis variables only.* A case is considered empty if values of all analysis variables are blank or missing. This is the default if analysis variables are specified. ANALYSISVARS is ignored if no analysis variables are specified.

**ALLVARS** *Procedure considers all variables.* A case is regarded as empty if all variables in the active dataset are blank or missing. This is the default if no analysis variables are specified. ID variables, split file variables, and rule outcome variables are always disregarded when determining whether a case is empty.

## **RULESUMMARIES Subcommand**

The RULESUMMARIES subcommand is used to summarize violations of single-variable validation rules.

- By default, single-variable rule violations are summarized by analysis variable.

- Cross-variable validation rules are always summarized by rule.
  - RULESUMMARIES generates a warning if summaries are requested but single-variable validation rules are not defined for at least one analysis variable.
- |                   |  |
|-------------------|--|
| <b>BYVARIABLE</b> | <i>Summarize by variable.</i> For each analysis variable, shows all single-variable validation rules that were violated and the number of values that violated each rule. Also reports the total number of single-variable rule violations for each variable. This is the default. |
| <b>BYRULE</b>     | <i>Summarize by rule.</i> For each single-variable validation rule, reports variables that violated the rule and the number of invalid values per variable. Also reports the total number of values that violated each rule across variables.                                      |
| <b>NONE</b>       | <i>No summaries are produced for single-variable rules.</i> An error occurs if NONE is specified along with any other RULESUMMARIES keyword.   |

## **CASEREPORT Subcommand**

The CASEREPORT subcommand is used to obtain a report that lists validation rule violations for individual cases.

- The case report includes violations of single-variable and cross-variable rules.
- CASEREPORT is ignored if no single-variable or cross-variable rules are applied.

### **DISPLAY Keyword**

- The case report is shown by default. To turn off the report, specify DISPLAY=NO.
- If you specify a case identifier, it is used to label cases in the report. Cases are also identified by case number.

### **MINVIOLATIONS Keyword**

- MINVIOLATIONS specifies the minimum number of rule violations required for a case to be included in the report. Specify a positive integer.
- The default is 1 violation.
- MINVIOLATIONS is ignored if DISPLAY=NO.

### **CASELIMIT Keyword**

- The CASELIMIT keyword specifies the maximum number of cases included in the case report. By default the maximum is 500.
- To change the maximum number of cases in the report, specify FIRSTN followed by a positive integer in parentheses.
- NONE turns off the case limit. All cases that meet the MINVIOLATIONS threshold are reported.
- CASELIMIT is ignored if DISPLAY=NO.

## **SAVE Subcommand**

The SAVE subcommand specifies optional variables to save. Specify one or more keywords, each followed by an optional variable name.

- The name, if specified, must be enclosed in parentheses—for example, `EMPTYCASE (Empty)`—and must be a valid variable name.
- If you do not specify a variable name, the procedure uses a default name.
- If the default name is used and it conflicts with that of an existing variable, a suffix is added to make the name unique.
- If you specify a name and it conflicts with that of an existing variable, an error occurs.

<b>EMPTYCASE</b>	<i>Empty case indicator.</i> Empty cases are assigned the value 1. All other cases are coded 0. The default rootname is <i>EmptyCase</i> . Values of <code>EMPTYCASE</code> reflect the scope specified via the <code>CASECHECKS</code> subcommand.
<b>INCOMPLETEID</b>	<i>Incomplete ID indicator.</i> Cases with empty or incomplete case identifiers are assigned a value 1. All other cases are coded 0. The default rootname is <i>IncompleteID</i> .
<b>DUPLICATEID</b>	<i>Duplicate ID group.</i> Cases that have the same case identifier (other than cases with incomplete identifiers) are assigned the same group number. Cases with unique or incomplete identifiers are coded 0. The default rootname is <i>DuplicateIDGroup</i> .
<b>RULEVIOLATIONS</b>	<i>Total count of single-variable and cross-variable validation rule violations.</i> The default rootname is <i>ValidationRuleViolations</i> . Ignored if no single-variable or cross-variable validation rules are applied.

## Single-Variable Validation Rules

To use a single-variable validation rule with the `VALIDATEDATA` procedure requires:

1. A rule definition (using the `DATAFILE ATTRIBUTE` command).

### Defining a Single-Variable Validation Rule

Single-variable validation rules are attributes of the data file and are generally defined according to the following syntax diagram:

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=$VD.SRule[n] ("Label='value',"
    "Description='value',"
    "Type='{Numeric}'"+
      {String }
      {Date  }
    "Format='dateformat'"
    "Domain='{Range}'"+
      {List }
    "FlagUserMissing='{Yes}'"+
      {No }
    "FlagSystemMissing='{Yes}'"+
      {No }
    "FlagBlank='{Yes}'"+
      {No }

  -- when Domain='Range' --
    "Minimum='value',"
    "Maximum='value',"
    "FlagUnlabeled='{Yes}'"+
      {No }
    "FlagNoninteger='{Yes}'"+
      {No }
```

```
-- when Domain='List' --
      "List='value1' 'value2' ... 'valuen'"+
      "CaseSensitive='{Yes}'"+
      "                {No }"+
      ").
```

- Single-variable rule attribute names must be of the form  $\$VD.SRule[n]$ , where  $n$  is a positive integer.
- Each rule specification includes a comma-separated list of properties that define the rule. Except where noted in the following table, each property is required. The order in which properties appear does not matter. Unrecognized rule properties are ignored. If the same property appears more than once in a rule specification, the value for the last one is honored.

Table 257-1

Properties for single-variable rule attributes

<i>Label</i>	The label is text that uniquely identifies the rule among single-variable and cross-variable rules. It cannot be the null string. Leading and trailing blanks are trimmed for the purpose of determining the uniqueness of the label.
<i>Description</i>	Any text (optional).
<i>Type</i>	This is the type of variable to which the rule can be applied. Select from Numeric, String, and Date. It is used by VALIDATEDATA to ensure that the rule is internally consistent and appropriate for the variable to which it is applied.
<i>Format</i>	This allows you to select the date format for rules that can be applied to date variables. Specify any IBM® SPSS® Statistics date or time format. <a href="#">For more information, see the topic Date and Time Formats on p. 63.</a> This property is not required if <i>Type</i> is not <i>Date</i> .
<i>FlagUserMissing</i>	Controls whether user-missing values are flagged as invalid. Specify <i>Yes</i> or <i>No</i> .
<i>FlagSystemMissing</i>	Controls whether system-missing values are flagged as invalid. This does not apply to string rule types. Specify <i>Yes</i> or <i>No</i> .
<i>FlagBlank</i>	Controls whether blank—that is, completely empty—string values are flagged as invalid. This does not apply to non-string rule types. Specify <i>Yes</i> or <i>No</i> .
<i>Domain</i>	You can specify the valid values either as a <i>Range</i> or a <i>List</i> of values. “Range” properties are ignored if <i>Domain</i> = 'List', and “List” properties are ignored if <i>Domain</i> = 'Range'.
<i>Minimum</i>	This property is optional**. Its value must be consistent with the rule type and format and be less than or equal to <i>Maximum</i> . The value is ignored if null. It cannot have a noninteger value if <i>FlagNoninteger</i> = 'Yes'.
<i>Maximum</i>	This property is optional**. Its value must be consistent with the rule type and format and be greater than or equal to <i>Minimum</i> . The value is ignored if null. It cannot have a non-integer value if <i>FlagNoninteger</i> = 'Yes'.
<i>FlagUnlabeled</i>	Controls whether unlabeled values are flagged as invalid. Specify <i>Yes</i> or <i>No</i> .
<i>FlagNoninteger</i>	Controls whether non-integer values are flagged as invalid. Specify <i>Yes</i> or <i>No</i> .
<i>List</i>	Specify one or more values consistent with the rule type and format. Null strings are ignored. Each item in the list must be quoted.
<i>CaseSensitive</i>	Controls whether case matters when string data values are checked against the list of acceptable values. Specify <i>Yes</i> or <i>No</i> .

\*\* Range properties are optional if *Domain*= 'Range'. However, if *Domain*= 'Range' and *FlagUserMissing*, *FlagSystemMissing*, and *FlagBlank* are all set to *No*, then *Minimum* or *Maximum* must be specified or *FlagUnlabeled* or *FlagNoninteger* must be *Yes*. Otherwise, no values would be validated when the rule is applied.

2. A rule outcome variable (using COMPUTE and VARIABLE ATTRIBUTE).

### **Creating a Rule Outcome Variable**

Rule outcome variables are generated to validate values of variables and are generally defined according to the following syntax diagram:

```
COMPUTE outcomevar1=expression for analysisvar1.
COMPUTE outcomevar2=...
```

```
VARIABLE ATTRIBUTE
  VARIABLES=outcomevarlist
  ATTRIBUTE=$VD.RuleOutcomeVar("Yes").
```

- The COMPUTE expression for each outcome variable should generate the value 1 for invalid cases of a variable, according to the associated single-variable rule definition.
  - The rule outcome variable is marked as such using the VARIABLE ATTRIBUTE command. Rule outcome variable attribute names must be *\$VD.RuleOutcomeVar* with the property *Yes*.
3. A link between the analysis and outcome variable (using VARIABLE ATTRIBUTE).

### **Linking Analysis and Outcome Variables**

Single-variable rule links are attributes of the analysis variables and are generally defined according to the following syntax diagram:

```
VARIABLE ATTRIBUTE
  VARIABLES=analysisvar1
  ATTRIBUTE=$VD.SRuleRef[n] ("Rule=' $VD.SRule[n] ', "+
                             "OutcomeVar='outcomevar1' ")
  /VARIABLES=analysisvar2...
```

- Single-variable rule link attribute names must be of the form *\$VD.SRuleRef[n]*, where *n* is a positive integer.
- The *Rule* and *OutcomeVar* properties are required. The order in which they appear does not matter. Unrecognized rule properties are ignored. If the same property appears more than once in a rule specification, the value for the last one is honored.

Table 257-2

*Properties for single-variable rule link attributes*

<i>Rule</i>	An existing single-variable validation rule whose <i>Type</i> property is consistent with that of the variable to which it is linked. String rules can be applied to string variables.
<i>Outcomevar</i>	A rule outcome variable associated with the analysis variable.

## **Cross-Variable Validation Rules**

To use a cross-variable validation rule with the VALIDATEDATA procedure requires:

1. A rule definition (using the DATAFILE ATTRIBUTE command).



### Creating a Rule Definition

Cross-variable validation rules are attributes of the data file and are generally defined according to the following syntax diagram:

```
DATAFILE ATTRIBUTE
  ATTRIBUTE=$VD.CRule[1] ("Label='value', "+
                          "Expression='value'."+
                          "OutcomeVar='outcomevar'").
```

Cross-variable rule attribute names must be of the form *\$VD.CRule[n]*, where *n* is a positive integer.

- Each rule specification includes a list of properties that define the rule. Each property is required. The order in which properties appear does not matter. Unrecognized rule properties are ignored. If the same property appears more than once in a rule specification, the value for the last one is honored.

Table 257-3

*Properties for cross-variable rule attributes*

<i>Label</i>	The label is text that uniquely identifies the rule among single-variable and cross-variable rules. It cannot be the null string. Leading and trailing blanks are trimmed for the purpose of determining the uniqueness of the label.
<i>Expression</i>	The expression must be valid command syntax for the right side of a COMPUTE statement. Any variables used in the expression must exist in the data file.
<i>OutcomeVar</i>	A rule outcome variable associated with the cross-variable rule.

2. A rule outcome variable (using COMPUTE and VARIABLE ATTRIBUTE).

### Creating a Rule Outcome Variable

```
COMPUTE outcomevar = expression.

VARIABLE ATTRIBUTE
  VARIABLES=outcomevar
  ATTRIBUTE=$VD.RuleOutcomeVar ("Yes").
```

- The COMPUTE expression for each outcome variable should generate the value 1 for invalid cases. It should be equivalent to the *Expression* in the cross-variable rule definition.
- The rule outcome variable is marked as such using the VARIABLE ATTRIBUTE command. Rule outcome variable attribute names must be *\$VD.RuleOutcomeVar* with the property *Yes*.

# VALUE LABELS

```
VALUE LABELS varlist value 'label' value 'label'... [/varlist...]  
[/datevarlist 'value' 'label'...]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- The maximum length of a value label is extended to 120 bytes (previous limit was 60 bytes).

Release 16.0

- Limitation preventing assignment of missing values to strings with a defined width greater than eight bytes removed.

## **Example**

```
VALUE LABELS JOBGRADE 'P' 'Parttime Employee' 'C' 'Customer Support'.
```

## **Overview**

VALUE LABELS deletes all existing value labels for the specified variable(s) and assigns new value labels. ADD VALUE LABELS can be used to add new labels or alter labels for specified values without deleting other existing labels.

## **Basic Specification**

The basic specification is a variable name and the individual values with their assigned labels.

## **Syntax Rules**

- Labels can be assigned to any previously defined variables.
- It is not necessary to enter value labels for all values for a variable.
- Each value label must be enclosed in quotes. For string variables, the values themselves must also be quotes.
- For date format variables (for example, DATE, ADATE), values expressed in date formats must be enclosed in quotes, and values must be expressed in the same date format as the defined date format for the variable.
- Value labels can contain any characters, including blanks. To enter an apostrophe as part of a label, enclose the label in double quotes or enter two apostrophes (two single quotes). [For more information, see the topic String Values in Command Specifications on p. 40.](#)
- Each value label can be up to 120 bytes long.

- The same labels can be assigned to the values of different variables by specifying a list of variable names. For string variables, the variables specified must be of equal length.
- Multiple sets of variable names and value labels can be specified on one `VALUE LABELS` command as long as the sets are separated by slashes.
- To continue a label from one command line to the next, specify a plus (+) sign before the continuation of the label. Each string segment of the label must be enclosed in quotes. To insert a blank between the strings, the blank must be included in the label specification.
- To control line wrapping of labels in pivot tables and charts, insert `\n` as part of the label wherever you want a line break. The `\n` is not displayed in output; it is interpreted as a line-break character. (*Note:* Labels will *always* wrap wherever `\n` appears in the defined label even if there is enough space to display the label without wrapping.)

### **Operations**

- Unlike most transformations, `VALUE LABELS` takes effect as soon as it is encountered in the command sequence. Thus, special attention should be paid to its position among commands. [For more information, see the topic Command Order on p. 41.](#)
- `VALUE LABELS` deletes all previously assigned value labels for the specified variables.
- The value labels assigned are stored in the dictionary of the working file and are automatically displayed on the output from many procedures.
- If a specified value is longer than the format of the variable, the program will be unable to read the full value and may not be able to assign the value label correctly.
- If the value specified for a string variable is shorter than the format of the variable, the value specification is right-padded without warning.

## **Examples**

### **Assigning Value Labels to Multiple Variables**

```
VALUE LABELS V1 TO V3 1 'Officials & Managers'
                6 'Service Workers'
/V4 'N' 'New Employee'.
```

- Labels are assigned to the values 1 and 6 for the variables between and including `V1` and `V3` in the active dataset.
- Following the required slash, a label for the value `N` of `V4` is specified. `N` is a string value and must be enclosed in quotes.
- If labels exist for the values 1 and 6 on `V1` to `V3` and the value `N` on `V4`, they are changed in the dictionary of the working file. If labels do not exist for these values, new labels are added to the dictionary.
- Existing labels for values other than 1 and 6 on `V1` to `V3` and the value `N` on `V4` are deleted.

### **Combining Strings to Construct Value Labels**

```
VALUE LABELS OFFICE88 1 "EMPLOYEE'S OFFICE ASSIGNMENT PRIOR"
+ " TO 1988".
```

- The label for *OFFICE88* is created by combining two strings with the plus sign. The blank between *PRIOR* and *TO* must be included in the first or second string to be included in the label.

### Value Labels for String Variables

```
VALUE LABELS=STATE REGION 'U' "UNKNOWN".
```

- The label *UNKNOWN* is assigned to the value *U* for both *STATE* and *REGION*.
- *STATE* and *REGION* must be string variables of equal length. If *STATE* and *REGION* have unequal lengths, a separate specification must be made for each, as in

```
VALUE LABELS STATE 'U' "UNKNOWN" / REGION 'U' "UNKNOWN".
```

### Using Value Labels with DATA LIST

```
DATA LIST / CITY 1-8(A) STATE 10-12(A).
VALUE LABELS STATE 'TEX' "TEXAS" 'TEN' "TENNESSEE"
              'MIN' "MINNESOTA".

BEGIN DATA
AUSTIN  TEX
MEMPHIS TEN
ST. PAUL MIN
END DATA.
FREQUENCIES VARIABLES=STATE.
```

- The `DATA LIST` command defines two variables. *CITY* is eight characters wide and *STATE* is three characters. The values are included between the `BEGIN DATA` and `END DATA` commands.
- The `VALUE LABELS` command assigns labels to three values of variable *STATE*. Each value and each label is specified in quotes.
- The format for the variable *STATE* must be at least three characters wide because the specified values, *TEX*, *TEN*, and *MIN*, are three characters. If the format for *STATE* were two characters, the program would issue a warning. This would occur even though the values named on `VALUE LABELS` and the values after `BEGIN DATA` agree.

### Forcing Value Labels to Wrap

```
VALUE LABELS myvar 1 "A long value label \n that always wraps".
FREQUENCIES myvar.
```

Figure 258-1  
Using `\n` to wrap value labels

A Fairly Long Label That Always Wraps

	Frequency	Percent	Valid Percent	Cumulative Percent
Valid A long value label that always wraps	1	100.0	100.0	100.0

# VARCOMP

VARCOMP is available in the Advanced Statistics option.

VARCOMP dependent variable BY factor list [WITH covariate list]

```
/RANDOM = factor [factor ...]

[/METHOD = {MINQUE({1})**}]
           {0}
           {ML      }
           {REML    }
           {SSTYPE({3}) }
           {1}

[/INTERCEPT = {INCLUDE**}]
               {EXCLUDE  }

[/MISSING = {EXCLUDE**}]
           {INCLUDE  }

[/REGWGT = varname]

[/CRITERIA = [CONVERGE({1.0E-8**})] [EPS({1.0E-8**})] [ITERATE({50**})]]
            {n      }      {n      }      {n      }

[/PRINT = [EMS] [HISTORY({1**})] [SS]]
         {n  }

[/OUTFILE = [VAREST] [{COVB} ('savfile'|'dataset') ]
            {CORB}

[/DESIGN = {[INTERCEPT] [effect effect ...]]]
```

\*\* Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
VARCOMP Y1 BY B C WITH X1 X2
/RANDOM = C.
```

## Overview

The VARCOMP procedure estimates variance components for mixed models. Following the general linear model approach, VARCOMP uses indicator variable coding to construct a design matrix and then uses one of the four available methods to estimate the contribution of each random effect to the variance of the dependent variable.

## Options

**Regression Weights.** You can use the REGWGT subcommand to specify regression weights for the model.

**Estimation Methods.** You can use the `METHOD` subcommand to use one of the four methods that are available for estimating variance components.

**Tuning the Algorithm.** You can control the values of algorithm-tuning parameters with the `CRITERIA` subcommand.

**Optional Output.** You can request additional output using the `PRINT` subcommand.

**Saving the Results.** You can save the variance component estimates and their asymptotic covariance matrix (if produced) to an external data file.

### **Basic Specification**

The basic specification is one dependent variable and one or more factor variables (that define the crosstabulation) and one or more factor variables on the `RANDOM` subcommand (to classify factors into either fixed or random factors). By default, `VARCOMP` uses the minimum norm quadratic unbiased estimator with unit prior weights to estimate variance components. Default output includes a factor-level information table and a variance component estimates table.

### **Subcommand Order**

- The variable specification must come first.
- Other subcommands can be specified in any order.

### **Syntax Rules**

- Only one dependent variable can be specified.
- At least one factor must be specified after `BY`.
- At least one factor must be specified on the `RANDOM` subcommand.

## **Example**

```
VARCOMP
  amtspent BY storeid  shopfor usecoup
  /RANDOM = storeid
  /METHOD = SSTYPE (3)
  /PRINT = SS
  /PRINT = EMS
  /DESIGN = shopfor usecoup shopfor*usecoup storeid
  /INTERCEPT = INCLUDE .
```

- The procedure fits a model to *amtspent* using *storeid*, *shopfor*, and *usecoup* as factors.
- `RANDOM` identifies *storeid* as a random effect.
- `METHOD` specifies ANOVA estimation using Type III sums of squares.
- `PRINT` requests tabular display of the sums of squares and expected mean squares.
- `DESIGN` indicates that the model fit is a factorial model for the fixed effects, plus a main-effects term for the random effect. The intercept is included.

## Variable List

The variable list specifies the dependent variable and the factors in the model.

- The dependent variable must be the first specification on VARCOMP.
- The factors follow the dependent variable and are separated from it by the keyword BY.
- The covariates, if any, follow the factors and are separated from the dependent variable and the factors by the keyword WITH.
- The dependent variable and the covariates must be numeric, but the factor variables can be either numeric or string.

## RANDOM Subcommand

The RANDOM subcommand allows you to specify random factors.

- You must specify at least one RANDOM subcommand with one random factor.
- You can specify multiple random factors on a RANDOM subcommand. You can also use multiple RANDOM subcommands. Specifications are accumulative.
- Only factors that are listed after the keyword BY in the variable list are allowed on the RANDOM subcommand.
- If you specify a factor on RANDOM, all effects containing the factor are automatically declared as random effects.

### Example

```
VARCOM Y BY DRUG SUBJECT
  /RANDOM = SUBJECT
  /DESIGN = DRUG DRUG*SUBJECT.
```

- This example specifies a mixed model where *DRUG* is the fixed factor and *SUBJECT* is a random factor.
- The default method MINQUE (1) is used to estimate the contribution of the random effect DRUG\*SUBJECT to the variance of the dependent variable.

## METHOD Subcommand

The METHOD subcommand offers four different methods for estimating the variances of the random effects. If more than one METHOD subcommand is specified, only the last subcommand is in effect. If the subcommand is not specified, the default method MINQUE (1) is used. METHOD cannot be specified without a keyword.

- |                  |  |
|------------------|--|
| <b>MINQUE(n)</b> | <i>Minimum norm quadratic unbiased estimator.</i> This method is the default method. When $n = 0$ , zero weight is assigned to the random effects and unit weight is assigned to the residual term. When $n = 1$ , unit weight is assigned to both the random effects and the residual term. By default, $n = 1$ . |
| <b>ML</b>        | <i>Maximum likelihood method.</i> Parameters of the fixed effects and variances of the random effects are estimated simultaneously. However, only the variances are reported.  |

<b>REML</b>	<i>Restricted maximum likelihood method.</i> Variances of the random effects are estimated based on residuals of the model after adjusting for the fixed effects.
<b>SSTYPE(n)</b>	<i>ANOVA method.</i> The ANOVA method equates the expected mean squares of the random effects to their observed mean squares. Their variances are then estimated by solving a system of linear equations. The expected mean squares are computed based on the chosen type of sum of squares. Two types are available in VARCOMP: Type I ( $n = 1$ ) and Type III ( $n = 3$ ). Type III is the default option for this method.

## ***INTERCEPT Subcommand***

The `INTERCEPT` subcommand controls whether an intercept term is included in the model. If more than one `INTERCEPT` subcommand is specified, only the last subcommand is in effect.

<b>INCLUDE</b>	<i>Include the intercept term.</i> The intercept (constant) term is included in the model. This setting is the default when <code>INTERCEPT</code> is not specified.
<b>EXCLUDE</b>	<i>Exclude the intercept term.</i> The intercept (constant) term is excluded from the model. <code>EXCLUDE</code> is ignored if you specify the keyword <code>INTERCEPT</code> on the <code>DESIGN</code> subcommand.

## ***MISSING Subcommand***

By default, cases with missing values for any of the variables on the `VARCOMP` variable list are excluded from the analyses. The `MISSING` subcommand allows you to include cases with user-missing values.

- Pairwise deletion of missing data is not available in `VARCOMP`.
- If more than one `MISSING` subcommand is specified, only the last subcommand is in effect.

<b>EXCLUDE</b>	<i>Exclude both user-missing and system-missing values.</i> This setting is the default when <code>MISSING</code> is not specified.
<b>INCLUDE</b>	<i>Treat user-missing values as valid.</i> System-missing values cannot be included in the analysis.

## ***REGWGT Subcommand***

`REGWGT` specifies the weight variable. Values of this variable are used as regression weights in a weighted least squares model.

- Specify one numeric variable name on the `REGWGT` subcommand.
- Cases with nonpositive values in the regression weight variable are excluded from the analyses.
- If more than one variable is specified on the same `REGWGT` subcommand, only the last variable is in effect.
- If more than one `REGWGT` subcommand is specified, only the last subcommand is in effect.

## ***CRITERIA Subcommand***

The `CRITERIA` subcommand specifies numerical tolerance for checking singularity and offers control of the iterative algorithm that is used for `ML` or `REML` estimation.



- Multiple CRITERIA subcommands are allowed.
- The last specified value for any keyword takes effect. If no value is specified, the default is used.

<b>EPS(n)</b>	<i>Epsilon value that is used as tolerance in checking singularity.</i> The value for <i>n</i> must be a positive value. The default is 1.0E-8.
<b>CONVERGE(n)</b>	<i>Convergence criterion.</i> Convergence is assumed if the relative change in the objective function is less than the specified value. The value for <i>n</i> must be a positive value. The default is 1.0E-8. This process is available only if you specify ML or REML on the METHOD subcommand.
<b>ITERATE(n)</b>	<i>Maximum number of iterations.</i> The value for <i>n</i> must be a positive integer. The default is 50. This process is available only if you specify ML or REML on the METHOD subcommand.

## **PRINT Subcommand**

The PRINT subcommand controls the display of optional output. If PRINT is not specified, the default output includes a factor information table and a variance component estimates table.

- For the maximum likelihood (ML) and restricted maximum likelihood (REML) methods, an asymptotic covariance matrix of the variance estimates table is also displayed.
- If more than one PRINT subcommand is specified, the specifications are accumulated. However, if you specify the keyword HISTORY more than once but with different values for *n*, the last specification is in effect.

<b>EMS</b>	<i>Expected mean squares.</i> This output shows expected mean squares of all of the effects. This process is available only if you specify SSTYPE (n) on the METHOD subcommand.
<b>HISTORY(n)</b>	<i>Iteration history.</i> The table contains the objective function value and variance component estimates at every <i>n</i> iteration. The value for <i>n</i> must be a positive integer. The default is 1. The last iteration is always printed if HISTORY is specified on PRINT. This process is available only if you specify ML or REML on the METHOD subcommand.
<b>SS</b>	<i>Sums of squares.</i> The table contains sums of squares, degrees of freedom, and mean squares for each source of variation. This process is available only if you specify SSTYPE (n) on the METHOD subcommand.

## **OUTFILE Subcommand**

The OUTFILE subcommand writes the variance component estimates to a data file or a previously declared dataset (DATASET DECLARE command) that can be used in other procedures. For the ML and REML methods, OUTFILE can also write the asymptotic covariance or correlation matrix. If more than one OUTFILE subcommand is specified, the last specification is in effect.

- OUTFILE writes to an external data file or previously declared dataset. You must specify a quoted file specification or dataset name in parentheses.

- COVB and CORB are available only if you specify ML or REML on the METHOD subcommand.
- COVB and CORB are mutually exclusive; only one of them can be specified on an OUTFILE subcommand.

<b>VAREST</b>	<i>Variance component estimates.</i> A variable will be created to contain the estimates, and another variable will be created to hold the labels of the variance components.
<b>COVB</b>	<i>Covariance matrix.</i> This matrix is the asymptotic covariance matrix of the variance component estimates. One variable is created for each variance component.
<b>CORB</b>	<i>Correlation matrix.</i> This matrix is the asymptotic correlation matrix of the variance component estimates. One variable is created for each variance component.
<b>(‘savfile’ ‘dataset’)</b>	<i>Output file specification or previously declared dataset name.</i> Specify a quoted file specification or dataset name, enclosed in parentheses. The variance component estimates and the asymptotic covariance or correlation matrix (if requested) are written to the same file.

## DESIGN Subcommand

The DESIGN subcommand specifies the effects in a model. DESIGN can be specified anywhere after the variable list. If more than one DESIGN subcommand is specified, only the last subcommand is in effect.

- Specify a list of effect terms to be included in the design. Each term must be separated from the next term by a comma or a space. Valid specifications include the keyword INTERCEPT, factors, covariates, and interaction or nested terms.
- The factors and covariates must have been specified on the variable list.
- If a factor is specified on the RANDOM subcommand, all effects that include that factor are random effects.
- If the DESIGN subcommand is omitted or specified without any term, the default design is generated. The default design includes the intercept term (if INTERCEPT=EXCLUDE is not specified), the covariates (if any) in the order in which they are specified on the variable list, the main factorial effects, and all orders of factor-by-factor interaction.

<b>INTERCEPT</b>	<i>Include the intercept term.</i> Specifying INTERCEPT on DESIGN explicitly includes the intercept term, regardless of the specification on the INTERCEPT subcommand.
<b>BY</b>	<i>Interaction.</i> You can also use the asterisk (*). Interaction terms can be formed among factors, among covariates, and between factors and covariates. Factors inside an interaction effect must be distinct. For factors <i>A</i> , <i>B</i> , and <i>C</i> , expressions like <i>A*C*A</i> or <i>A*A</i> are invalid. Covariates inside an interaction effect do not have to be distinct. For covariate <i>X</i> , <i>X*X</i> is the product of <i>X</i> and itself. This is equivalent to a covariate whose values are the square of the values of <i>X</i> .
<b>WITHIN</b>	<i>Nesting.</i> You can also use a pair of parentheses. Factors and covariates can be nested within factors, but no effects can be nested within covariates. Suppose that <i>A</i> and <i>B</i> are factors and <i>X</i> and <i>Y</i> are covariates. Both <i>A(B)</i> and <i>X(B)</i> are valid, but <i>X(Y)</i> is not valid. Factors inside a nested effect must be distinct. Expressions like <i>A(A)</i> are invalid.

Multiple-level nesting is supported. For example,  $A(B(C))$  or  $A$  WITHIN  $B$  WITHIN  $C$  means that factor  $B$  is nested within factor  $C$ , and factor  $A$  is nested within  $B(C)$ . The expression  $A(B)(C)$  is invalid.

Nesting within an interaction effect is valid. For example,  $A(B*C)$  means that factor  $A$  is nested within  $B*C$  while  $X(A*B)$  means that covariate  $X$  is nested within  $A*B$ .

Interactions among nested effects are allowed. For example,  $A*B(C)$  means interaction between  $A$  and  $B$  within levels of  $C$ .  $X*Y(A)$  means the product of  $X$  and  $Y$  nested within levels of  $C$ . The expression  $A(C)*B(C)$  is invalid.

### **Example**

```
VARCOM Y BY DRUG SUBJECT WITH X
  /RANDOM = SUBJECT
  /DESIGN = DRUG SUBJECT DRUG*SUBJECT X*SUBJECT.
```

- The `DESIGN` subcommand specifies two main effects and two interaction terms.
- All effects that involve the factor `SUBJECT` are assumed to be random.

# VARIABLE ALIGNMENT

```
VARIABLE ALIGNMENT varlist ({LEFT }) ... [/varlist...]  
                        {CENTER}  
                        {RIGHT }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
VARIABLE ALIGNMENT sales95 sales96 (LEFT)  
  /id gender (RIGHT).
```

## **Overview**

VARIABLE ALIGNMENT specifies the alignment of data values in the Data Editor. It has no effect on the format of the variables or the display of the variables or values in other windows or printed results.

### **Basic Specification**

The basic specification is a variable name and the keyword LEFT, RIGHT, or CENTER in parentheses.

# VARIABLE ATTRIBUTE

```
VARIABLE ATTRIBUTE
  VARIABLES=varlist
  ATTRIBUTE=name('value') name('value')...
              arrayname[1]('value') arrayname[2]('value')...
  DELETE=name name...
              arrayname[n] arrayname...
/VARIABLES...
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 14.0

- Command introduced.

## **Example**

```
VARIABLE ATTRIBUTE VARIABLES=Age Gender Region
  ATTRIBUTE=DemographicVars ('1').
```

## **Overview**

VARIABLE ATTRIBUTE provides you with the ability to define your own variable attributes and assign attribute values to variables in the active dataset. For example, you could create a variable attribute that identifies the type of response for survey questions (e.g., single selection, multiple selection, fill-in-the-blank) or the formulas used for computed variables.

- User-defined variable attributes are saved with the data file in the data dictionary.
- The VARIABLE ATTRIBUTE command takes effect immediately, updating the data dictionary without requiring a data pass.
- You can display a list of data file and variable attributes with DISPLAY ATTRIBUTES. [For more information, see the topic DISPLAY on p. 648.](#)

## **Basic Specification**

The basic specification is:

- The VARIABLES subcommand followed by an equals sign (=) and a list of valid variables.
- The ATTRIBUTE keyword followed by an equals sign (=) and one or more attribute names that follow variable naming rules, with each attribute name followed by a quoted attribute value enclosed in parentheses.

or

- The DELETE keyword followed by an equals sign (=) and a list of defined attribute names or attribute arrays.

**Syntax Rules**

- The VARIABLES subcommand is required.
- All subcommands and keywords (VARIABLES, ATTRIBUTE, DELETE) must be followed by an equals sign (=).
- Each ATTRIBUTE keyword must be followed by a name that follows variable naming rules and a single, quoted attribute value enclosed in parentheses. [For more information, see the topic Variable Names on p. 49.](#)
- Attribute names that begin with @ cannot be displayed in Variable View of the Data Editor and are not displayed by DISPLAY DICTIONARY or DISPLAY ATTRIBUTES. They can only be displayed with DISPLAY @ATTRIBUTES.
- Attribute names that begin with a dollar sign (\$) are reserved for internal IBM® SPSS® Statistics use.
- All attribute values must be quoted (single or double quotes), even if the values are numbers.
- Attribute values can be up to 32,767 bytes in length.

**Attribute Arrays**

If you append an integer enclosed in square brackets to the end of an attribute name, the attribute is interpreted as an array of attributes. For example,

```
VARIABLE ATTRIBUTE VARIABLES=Age
  ATTRIBUTE=MyAttribute[99]('not quite 100').
```

will create 99 attributes—*MyAttribute[01]* through *MyAttribute[99]*—and will assign the value “not quite 100” to the last one.

- Array subscripts (the value enclosed in square brackets) must be integers greater than 0. (Array subscript numbering starts with 1, not 0.)
- If the root name of an attribute array is the same as an existing attribute name for any variables specified on the VARIABLES subcommand, the attribute array replaces the existing attribute for those variables (and vice versa). If no value is assigned to the first element in the array (subscript [1]), the original attribute value is used for that element value.

With the DELETE keyword, the following rules apply to attribute arrays:

- If you specify DELETE followed by an array rootname and no value in square brackets, all attributes in the array are deleted.
- If you specify DELETE with an array name followed by an integer value in square brackets, the specified array element is deleted and the integer values for all subsequent attributes in the array (in numeric order) are changed to reflect the new order of array elements.

**Example**

```
VARIABLE ATTRIBUTE
  VARIABLES=Age Income1 Income2 Income3
  ATTRIBUTE=AnswerFormat('Fill-in')
  /VARIABLES=Gender OwnHome MortgageType
  ATTRIBUTE=AnswerFormat('Single select')
  /VARIABLES=AvgIncome
```

```

ATTRIBUTE=AnswerFormat('Computed value')
/VARIABLES=AvgIncome
ATTRIBUTE=Formula('mean(Income1, Income2, Income3)')
/VARIABLES=AvgIncome
ATTRIBUTE=DerivedFrom[1]('Income1')
DerivedFrom[2]('Income2')
DerivedFrom[3]('Income3').
DISPLAY ATTRIBUTES.

```

- The variables *Age*, *Income1*, *Income2*, and *Income3* are assigned the variable attribute *AnswerFormat* with a value of *Fill-in*.
- The variables *Gender*, *OwnHome*, and *MortgageType* are assigned the variable attribute *AnswerFormat* with a value of *Single select*.
- The variable *AvgIncome* is assigned the variable attribute *AnswerFormat* with a value of *Computed value*.
- The variable *AvgIncome* is assigned the variable attribute *Formula*, with a value the describes the formula used to calculate the value of this variable.
- The variable *AvgIncome* is assigned the variable attribute array *DerivedFrom*, and each array value identifies one of the variables used to calculate the value of *AvgIncome*.
- `DISPLAY ATTRIBUTES` lists the variable attributes for each variable in the active dataset.

Figure 261-1  
Variable attributes

Variable Attributes		
Variable	Attribute	Value
Age	AnswerFormat	Fill-in
Gender	AnswerFormat	Single select
Income1	AnswerFormat	Fill-in
Income2	AnswerFormat	Fill-in
Income3	AnswerFormat	Fill-in
OwnHome	AnswerFormat	Single select
MortgageType	AnswerFormat	Single select
AvgIncome	AnswerFormat	Computed value
	DerivedFrom[1]	Income1
	DerivedFrom[2]	Income2
	DerivedFrom[3]	Income3
	Formula	mean(Income1, Income2, Income3)

# VARIABLE LABELS

```
VARIABLE LABELS varname 'label' [/varname...]
```

## **Example**

```
VARIABLE LABELS YRHIRED 'YEAR OF FIRST HIRING'.
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Overview**

VARIABLE LABELS assigns descriptive labels to variables in the active dataset.

### **Basic Specification**

The basic specification is a variable name and the associated label in quotes.

### **Syntax Rules**

- Labels can be added to any previously defined variable. It is not necessary to enter labels for all variables in the active dataset.
- Each variable label must be enclosed in quotes.
- Variable labels can contain any characters, including blanks. To enter an apostrophe as part of a label, enclose the label in double quotes or enter two apostrophes (two single quotes). [For more information, see the topic String Values in Command Specifications on p. 40.](#)
- Each variable label can be up to 255 bytes long, although some procedures print fewer than the 255 bytes. All statistical procedures display at least 40 bytes.
- Multiple variables can be assigned labels on a single VARIABLE LABELS command. Only one label can be assigned to each variable, and each label can apply to only one variable.
- To continue a label from one command line to the next, specify a plus (+) sign before the continuation of the label. Each string segment of the label must be enclosed in quotes. To insert a blank between the strings, the blank must be included in the label specification.
- To control line wrapping of labels in pivot tables and charts, insert \n as part of the label wherever you want a line break. The \n is not displayed in output; it is interpreted as a line-break character. (*Note:* Labels will *always* wrap wherever \n appears in the defined label, even if there is enough space to display the label without wrapping.)

### **Operations**

- Unlike most transformations, VARIABLE LABELS takes effect as soon as it is encountered in the command sequence. Thus, special attention should be paid to its position among commands. [For more information, see the topic Command Order on p. 41.](#)



- Variable labels are automatically displayed in the output from many procedures and are stored in the dictionary of the active dataset.
- `VARIABLE LABELS` can be used for variables that have no previously assigned variable labels. If a variable has a previously assigned variable label, the new label replaces the old label.

## Examples

### Assigning Variable Labels to Multiple Variables

```
VARIABLE LABELS YRHIRED 'YEAR OF FIRST HIRING'
DEPT88 'DEPARTMENT OF EMPLOYMENT IN 1988'
SALARY88 'YEARLY SALARY IN 1988'
JOB CAT 'JOB CATEGORIES'.
```

- Variable labels are assigned to the variables `YRHIRED`, `DEPT88`, `SALARY88`, and `JOB CAT`.

### Combining Strings to Construct Variable Labels

```
VARIABLE LABELS OLDSAL "EMPLOYEE'S GROSS SALARY PRIOR"
+ " TO 1988".
```

- The label for `OLDSAL` is created by combining two strings with the plus sign. The blank between `PRIOR` and `TO` must be included in the first or second string to be included in the label.

### Forcing Variable Labels to Wrap

```
VARIABLE LABELS myvar "A Fairly Long Label \n That Always Wraps".
FREQUENCIES myvar.
```

Figure 262-1  
Using `\n` to wrap variable labels

**A Fairly Long Label  
That Always Wraps**

	Frequency	Percent	Valid Percent	Cumulative Percent
Valid	1.00	1	100.0	100.0

# VARIABLE LEVEL

```
VARIABLE LEVEL varlist ({SCALE }) ... [/varlist...]  
                        {ORDINAL}  
                        {NOMINAL}
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
VARIABLE LEVEL sales95 sales96 (SCALE)  
                /region division (NOMINAL)  
                /expense (ORDINAL).
```

## **Overview**

VARIABLE LEVEL specifies the level of measurement for variables.

### **Basic Specification**

The basic specification is a variable name, followed by a measurement level enclosed in parentheses. The measurement level can be:

- |                |  |
|----------------|--|
| <b>NOMINAL</b> | A variable can be treated as nominal when its values represent categories with no intrinsic ranking; for example, the department of the company in which an employee works. Examples of nominal variables include region, zip code, or religious affiliation.  |
| <b>ORDINAL</b> | A variable can be treated as ordinal when its values represent categories with some intrinsic ranking; for example, levels of service satisfaction from highly dissatisfied to highly satisfied. Examples of ordinal variables include attitude scores representing degree of satisfaction or confidence and preference rating scores. |
| <b>SCALE</b>   | A variable can be treated as scale when its values represent ordered categories with a meaningful metric, so that distance comparisons between values are appropriate. Examples of scale variables include age in years and income in thousands of dollars.  |

# VARIABLE ROLE

```
VARIABLE ROLE  
  [/INPUT varlist]  
  [/TARGET varlist]  
  [/BOTH varlist]  
  [/NONE varlist]  
  [/PARTITION varlist]  
  [/SPLIT varlist]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 18.0

- Command introduced.

## **Example**

```
VARIABLE ROLE  
  /INPUT Age Income Gender Education Debt  
  /TARGET Credit_rating  
  /NONE Customer_ID.
```

## **Overview**

Some dialogs support predefined roles that can be used to pre-select variables for analysis. When you open one of these dialogs, variables that meet the role requirements will be automatically displayed in the destination list(s). Available role subcommands are:

**Input.** The variable will be used as an input (e.g., predictor, independent variable).

**Target.** The variable will be used as an output or target (e.g., dependent variable).

**Both.** The variable will be used as both input and output.

**None.** The variable has no role assignment.

**Partition.** The variable will be used to partition the data into separate samples for training, testing, and validation.

**Split.** Included for round-trip compatibility with IBM® SPSS® Modeler. Variables with this role are *not* used as split-file variables in IBM® SPSS® Statistics.

- By default, all variables are assigned the *Input* role. This includes data from external file formats and data files from versions of SPSS Statistics prior to version 18.
- Role assignment only affects dialogs that support role assignment. It has no effect on command syntax.

**Basic Specification**

The basis specification is the command name, at least one subcommand, and at least one variable specified on that subcommand.

**Syntax Rules**

- At least one subcommand must be specified.
- Only a single instance of each subcommand is allowed.
- Each subcommand must be preceded by a forward slash.
- Each subcommand must specify at least one variable.
- The command name and subcommand names must be spelled in full.
- Keyword `ALL` can be used to specify all variables in the active dataset, and keyword `TO` can be used to specify a list of variables (in file order).

**Operations**

- If a variable is specified more than once, the last role assignment is used.
- This command takes effect immediately. It does not read the active dataset or execute pending transformations.

**Example**

```
VARIABLE ROLE  
  /INPUT Age Income Gender Education Debt  
  /TARGET Credit_rating  
  /NONE Customer_ID.
```

# VARIABLE WIDTH

```
VARIABLE WIDTH varlist (n) ... [/varlist...]
```

## **Example**

```
VARIABLE WIDTH sales95 sales96 (10)  
  /id gender (2).
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Overview**

VARIABLE WIDTH specifies the column width for the display of variables in the Data Editor. It has no effect on the format of the variable or the display of the variable or values in other windows or printed results.

### **Basic Specification**

The basic specification is a variable name and a positive integer in parentheses for the column width. Column width for proportional fonts is based on average character width. Depending on the characters used in the value, more or fewer characters may be displayed in the specified width.

# VARSTOCASES

```
VARSTOCASES

/MAKE new variable ["label"] [FROM] varlist [/MAKE ...]

[/INDEX = {new variable ["label"]                               }]
          {new variable ["label"] (make variable name)         }
          {new variable ["label"] (n) new variable ["label"] (n) ...}

[/ID = new variable ["label"]]

[/NULL = {DROP**}]
         {KEEP  }

[/COUNT=new variable ["label"]]

[/KEEP={ALL**  } ] [/DROP=varlist]
         {varlist}
```

**\*\***Default if the subcommand is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
VARSTOCASES /MAKE newvar FROM var1 TO var4.
```

## Overview

A **variable** contains information that you want to analyze, such as a measurement or a test score. A **case** is an observation, such as an individual or an institution. In a simple data structure, each variable is a single **column** in your data. So if you are measuring test scores, for example, all test score values would appear in only one column. In a simple data structure, each case is a single **row** in your data. So if you were measuring scores for all students in a class, there would be a row for each student.

VARSTOCASES restructures **complex** data structures (in which information about a variable is stored in more than one column) into a data file in which those measurements are organized into separate rows of a single column. It replaces the active dataset.

You can use VARSTOCASES to restructure data files in which repeated measurements of a single case were recorded in one row into a new data file in which each measurement for a case appears in a new row.

## Options

**Creating New Variables.** You can create an **identification** variable that identifies the row in the original data file that was used to create a group of new rows, a **count** variable that contains the number of new rows generated by a row in the original data, and one or more **index** variables that identify the original variable from which the new row was created.

**Variable Selection.** You can use the `DROP` and `KEEP` subcommands to specify which variables from the original data file are included in the new data file.

### **Basic Specification**

The basic specification is one or more `MAKE` subcommands, each of which specifies a list of variables to be combined into a single variable in which each value is displayed on a separate row.

### **Subcommand Order**

Subcommands can be specified in any order.

### **Syntax Rules**

- The `MAKE` subcommand is required and can be specified as many times as needed.
- The rest of the subcommands can be specified only once.

### **Operations**

- **Row order.** New rows are created in the order in which the variables are specified on the `FROM` list.
- **Propagated variables.** Variables that are not named on the `MAKE` or `DROP` subcommands are kept in the new data file. Their values are propagated for each new row.
- **Split file processing.** The `SPLIT FILE` command does not affect the results of `VARSTOCASES`. If split file processing is in effect, it will remain in effect in the new data file unless a variable that is used to split the file is named on the `MAKE` or `DROP` subcommands.
- **Weighted files.** The `WEIGHT` command does not affect the results of `VARSTOCASES`. If original data are weighted, the new data will be weighted unless the variable that is used to split the file is named on the `MAKE` or `DROP` subcommands.
- **Selected cases.** The `FILTER` and `USE` commands do not affect the results of `VARSTOCASES`. It processes all cases.

### **Limitations**

The `TEMPORARY` command cannot be in effect when `VARSTOCASES` is executed.

## **Example**

The following is the `LIST` output for a data file where repeated measurements for the same case are stored in variables on a single row:

caseid	var1	var2	var3	var4
001	.00	.05	5.00	3.00
002	7.00	1.00	5.00	4.00
003	6.00	3.00	6.00	2.00

VARSTOCASES

---

The command:

```
VARSTOCASES
  /MAKE newvar FROM var1 TO var4.
```

creates a new variable, *newvar*, using the values of *var1* through *var4*. The LIST output for the new active dataset is as follows:

caseid	newvar
001	.00
001	.05
001	5.00
001	3.00
002	7.00
002	1.00
002	5.00
002	4.00
003	6.00
003	3.00
003	6.00
003	2.00

The values for the new variable *newvar* are the values from *var1* through *var4* from the original data. There are now four rows for each case—one row for each variable that was named on the FROM list.

## **MAKE Subcommand**

The MAKE subcommand names, and optionally labels, the new variable to be created from the variables on the FROM list.

- One new variable is required on each MAKE subcommand. It must have a unique name.
- The label for the new variable is optional and, if specified, must be enclosed in quotes.
- The new variable will have the values of the variables listed on the FROM list. For each case in the original data, one new row will be created for each variable on the FROM list.
- All of the variables on the FROM list are required to be of the same type. For example, they must all be numeric or they must all be string.
- The dictionary information for the new variable (for example, value labels and format) is taken from the first variable in the FROM list. If string variables of different lengths are specified, the longest length is used.
- Rows are created in the order in which variables appear on the FROM list.
- Variables that appear on the FROM list will not appear in the new data file.
- Variables that are kept in the new data file and not named on the FROM list will have their values propagated for each new row.
- When multiple MAKE subcommands are used, a variable may not appear on more than one FROM list.
- A variable may be listed more than once on a FROM list. Its values are repeated.
- When multiple MAKE subcommands are used, the FROM lists must all contain the same number of variables (variables that are listed more than once must be included in the count).



## ***ID Subcommand***

The `ID` subcommand creates a new variable that identifies the permanent case sequence number (*\$casenum*) of the original row that was used to create the new rows. Use the `ID` subcommand when the original data file does not contain a variable that identifies cases.

- One new variable is named on the `ID` subcommand. It must have a unique name.
- The label for the new variable is optional and, if specified, must be enclosed in quotes.
- The format of the new variable is F8.0.

## ***INDEX Subcommand***

In the original data file, a case appears on a single row. In the new data file, that case will appear on multiple rows. The `INDEX` subcommand creates a new variable that sequentially identifies a group of new rows based on the original variables from which it was created.

- You can choose among three types of indices—a simple numeric index, an index that lists the variables on a `FROM` list, or multiple numeric indices.
- New variable(s) are named on the `INDEX` subcommand. A new variable must have a unique name.
- The label for the new variable is optional and, if specified, must be enclosed in quotes.

### ***Simple Numeric Index***

A simple numeric index numbers the rows sequentially within a new group.

- The basic specification is `/INDEX=ivar`, where *ivar* is a name for the new index variable.
- The new index variable starts with 1 and increments each time a `FROM` variable is encountered in a row in the original file. After the last `FROM` variable is encountered, the index restarts at 1.
- Gaps in the index sequence can occur if null data are dropped.

#### ***Example***

```
VARSTOCASES
  /MAKE newvar FROM var1 TO var4
  /INDEX=ivar.
```

caseid	ivar	newvar
001	1	.00
001	2	.05
001	3	5.00
001	4	3.00
002	1	7.00
002	2	1.00
002	3	5.00
002	4	4.00
003	1	6.00
003	2	3.00
003	3	6.00
003	4	2.00

## Variable Name Index

A variable name index works like the simple numeric index, except that it lists the name of the original FROM variable instead of a sequential number.

- The basic specification is `/INDEX=ivar (make variable name)`, where *ivar* is the name for the new index variable and *make variable name* is the name of a variable on the MAKE subcommand from which the index is to be constructed.
- The new index variable is a string that lists the name of the FROM variable from which the new row was created.

### Example

```
VARSTOCASES
  /MAKE newvar FROM var1 TO var4
  /INDEX=ivar (newvar).
```

caseid	ivar	newvar
001	VAR1	.00
001	VAR2	.05
001	VAR3	5.00
001	VAR4	3.00
002	VAR1	7.00
002	VAR2	1.00
002	VAR3	5.00
002	VAR4	4.00
003	VAR1	6.00
003	VAR2	3.00
003	VAR3	6.00
003	VAR4	2.00

## Multiple Numeric Indices

Multiple numeric indices are used to identify groups of new rows that share a particular combination of factors. You can create multiple numeric indices if the original variables are ordered so that levels of a given factor are grouped together.

- The basic specification is `/INDEX=ivar(n) ivar(n) ...`, where *ivar* is the name of the new index for a factor and *n* is the number of factor levels represented in the variable group for which the index is being constructed.
- The last index specified varies the fastest.

### Example

	B1	B2
A1	.00	.05
A2	5.00	3.00

- Data were collected for a designed experiment with two levels of factor A and two levels of factor B. The table shows the data for the first case.

caseid	v_a1b1	v_a1b2	v_a2b1	v_a2b2
001	.00	.05	5.00	3.00

- The original data file is structured so that each case has one variable for each combination of factors. Note that factor B varies fastest.

```
VARSTOCASES
  /MAKE newvar FROM v_a1b1 TO v_a2b2
  /INDEX=a(2) b(2).
```

caseid	a	b	newvar
001	1	1	.00
001	1	2	.05
001	2	1	5.00
001	2	2	3.00

- The command restructures the data file and creates two indices, *A* and *B*.

## NULL Subcommand

The `NULL` subcommand checks each potential new row for **null** values. A null value is a system-missing or blank value. By default, `VARSTOCASES` does not add a new row that contains null values for all variables created by `MAKE` subcommands. You can change the default null-value treatment with the `NULL` subcommand.

- DROP** *Do not include a new row when all MAKE variables are null.* A potential new row with null values for all of the variables created by `MAKE` subcommands is excluded from the new data file. This is the default. With this option, you may want to create a count variable to keep track of new rows because cases in the original data file are not guaranteed to appear in the new data file.
- KEEP** *Include a new row when all MAKE variables are null.* A potential new row with null values for all of the variables created by the `MAKE` subcommand is included in the new data. With this option, you may not need a count variable to keep track of cases because each row in the original data will result in a consistent number of rows in the new data file.

## COUNT Subcommand

When there are no null data, `VARSTOCASES` generates *n* new rows for each row in the original data file, where *n* is the number of variables on the `FROM` list(s). When the original data file contains null values and you drop them, it is possible to generate a different number of rows for a given subject in the original data file. The `COUNT` subcommand creates a new variable that contains the number of new rows generated by the original subject.

- One new variable is named on the `COUNT` subcommand. It must have a unique name.
- The label for the new variable is optional and, if specified, must be enclosed in quotes.
- The format of the new variable is F4.0.

## DROP and KEEP Subcommands

The `DROP` and `KEEP` subcommands are used to include only a subset of variables in the new active dataset. The `DROP` subcommand specifies a set of variables to exclude and the `KEEP` subcommand specifies a set of variables to retain. Variables not specified on the `KEEP` subcommand are dropped.

- `DROP` and `KEEP` cannot be used with variables that appear on a `FROM` list.

---

*VARSTOCASES*

- `DROP` and `KEEP` are mutually exclusive. Only one `DROP` or one `KEEP` subcommand can be used on the `VARSTOCASES` command.
- `KEEP` affects the order of variables in the new data file. The order of the variables kept in the new data file is the order in which they are named on the `KEEP` subcommand.

***Example***

```
VARSTOCASES  
  /MAKE newvar FROM var1 to var4  
  /DROP caseid.
```

- *Caseid* is dropped from the new data file. The new data file contains one variable, *newvar*.

# VECTOR

```
VECTOR {vector name=varname TO varname} [/vector name...]  
      {vector name(n [format])      }
```

## Example

```
VECTOR V=V1 TO V6.
```

## Overview

VECTOR associates a vector name with a set of existing variables or defines a vector of new variables. A vector is a set of variables that can be referred to using an index. The vector can refer to either string or numeric variables, and the variables can be permanent or temporary.

For each variable in the reference list, VECTOR generates an element. Element names are formed by adding a subscript in parentheses to the end of the vector name. For example, if the vector *AGES* has three elements, the element names are *AGES(1)*, *AGES(2)*, and *AGES(3)*. Although the VECTOR command has other uses within the transformation language, it is most often used with LOOP structures because the indexing variable on LOOP can be used to refer to successive vector elements.

## Options

**File Structures.** VECTOR can be used with the END CASE command to restructure data files. You can build a single case from several cases or, conversely, you can build several cases from a single case (see the examples for the END CASE command on p. 677).

**Short-Form Vectors.** VECTOR can be used to create a list of new variables and the vector that refers to them simultaneously. VECTOR in the short form can be used to establish the dictionary order of a group of variables before they are defined on a DATA LIST command (see “VECTOR: Short Form” on p. 2145).

## Basic Specification

- The basic specification is VECTOR, a vector name, a required equals sign, and the list of variables that the vector refers to. The TO keyword must be used to specify the variable list, and it defines a variable list based on file order.
- For the short form of VECTOR, the basic specification is VECTOR, an alphabetical prefix, and, in parentheses, the number of variables to be created.

## Syntax Rules

- Multiple vectors can be created on the same command by using a slash to separate each set of specifications.
- Variables specified on VECTOR must already be defined unless the short form of VECTOR is used to create variables (see “VECTOR: Short Form” on p. 2145).

- The `TO` convention must be used to specify the variable list. Thus, variables specified must be consecutive and must be from the same dictionary, permanent or scratch.
- A single vector must comprise all numeric variables or all string variables. The string variables must have the same length.
- A scalar (a variable named on `NUMERIC`), a function, and a vector can all have the same name, for example `MINI`. The scalar can be identified by the lack of a left parenthesis following the name. Where a vector has the same name as a function (or the abbreviation of a function), the vector name takes precedence (see the example on name conflicts in “VECTOR: Short Form” on p. 2145).
- Vector element names must always be specified with a subscript in parentheses.

### **Operations**

- `VECTOR` takes effect as soon as it is encountered in the command sequence, unlike most transformations, which do not take effect until the data are read. Thus, special attention should be paid to its position among commands. [For more information, see the topic Command Order on p. 41.](#)
- `VECTOR` is in effect only until the first procedure that follows it. The vector must be redeclared to be reused.

### **Limitations**

- A vector element is an *expression* and can be used wherever an expression can be used in transformation commands. It can also be the *target* of an assignment in a `COMPUTE` or `IF` command. It cannot be used in other transformation commands where a variable name is required. Specifically, vector elements cannot be used instead of variable names in `RECODE`, `COUNT`, `XSAVE`, `DO REPEAT`, or other transformation commands that require variable names but not expressions.
- Vectors cannot be used in procedure commands (for example, statistical or charting procedures).
- Vector names (for the whole vector rather than elements) cannot be used anywhere but in the `VECTOR` command itself.

## **Examples**

### **Example**

```
DATA LIST FREE /height age weight.  
BEGIN DATA  
1 2 3  
END DATA.  
VECTOR x=height TO weight.  
VECTOR y=age TO weight.  
VECTOR z=age TO height.
```

- The list of variables specified with `TO` defines a list of variables in file order.
- Vector `x` contains the variables *height*, *age*, and *weight*.
- Vector `y` contains the variables *age* and *weight*.

- The last `VECTOR` command generates an error because *height* comes before *age* in file order.

### Example

\* Replace a case's missing values with the mean of all nonmissing values for that case.

```
DATA LIST FREE /V1 V2 V3 V4 V5 V6 V7 V8.
MISSING VALUES V1 TO V8 (99).
COMPUTE MEANSUB=MEAN(V1 TO V8).
```

```
VECTOR V=V1 TO V8.
LOOP #I=1 TO 8.
+ DO IF MISSING (V(#I)).
+ COMPUTE V(#I)=MEANSUB.
+ END IF.
END LOOP.
```

```
BEGIN DATA
1 99 2 3 5 6 7 8
2 3 4 5 6 7 8 9
2 3 5 5 6 7 8 99
END DATA.
LIST.
```

- The first `COMPUTE` command calculates the variable *MEANSUB* as the mean of all nonmissing values for each case.
- `VECTOR` defines the vector *V* with the original variables as its elements.
- For each case, the loop is executed once for each variable. The `COMPUTE` command within the loop is executed only when the variable has a missing value for that case. `COMPUTE` replaces the missing value with the value of *MEANSUB*.
- For the first case, the missing value for the variable *V2* is changed to the value of *MEANSUB* for that case. The missing value for the variable *V8* for the third case is changed to the value of *MEANSUB* for that case.

For additional examples of `VECTOR`, see the examples for the `END CASE` command on p. 677 and the `IF` command on p. 962.

## VECTOR: Short Form

`VECTOR` can be used to create a list of new variables and the vector that refers to them simultaneously. The short form of `VECTOR` specifies a prefix of alphanumeric characters followed, in parentheses, by the length of the vector (the number of variables to be created).

- The new variable names must not conflict with existing variables. If the prefix starts with the `#` character, the new variables are created according to the rules for scratch variables.
- More than one vector of the same length can be created by naming two or more prefixes before the length specification.
- By default, variables created with `VECTOR` receive `F8.2` formats. Alternative formats for the variables can be specified by including a format specification with the length specification within the parentheses. The format and length can be specified in either order and must be separated by at least one space or comma. If multiple vectors are created, the assigned format applies to all of them unless you specify otherwise.

**Creating a Vector from a Set of New Scratch Variables**

```
VECTOR #WORK(10) .
```

- The program creates the vector *#WORK*, which refers to 10 scratch variables: *#WORK1*, *#WORK2*, and so on, through *#WORK10*. Thus, the element *#WORK(5)* of the vector is the variable *#WORK5*.

**Creating Multiple Vectors of the Same Length**

```
VECTOR X, Y(5) .
```

- *VECTOR* creates the vectors *X* and *Y*, which refer to the new variables *X1* through *X5* and *Y1* through *Y5*, respectively.

**Specifying the Format of Vector Variables**

```
VECTOR X(6, A5) .
```

- *VECTOR* assigns an A5 format to the variables *X1* through *X6*.

**Creating Multiple Vectors of Different Lengths and Formats**

```
VECTOR X, Y(A5, 6) Z(3, F2) .
```

- *VECTOR* assigns A5 formats to the variables *X1* to *X6* and *Y1* to *Y6*, and F2 formats to the variables *Z1* to *Z3*. It doesn't matter whether the format or the length is specified first within the parentheses.

**Predetermining Variable Order Using the Short Form of VECTOR**

```
INPUT PROGRAM.
VECTOR X Y (4, F8.2) .
DATA LIST / X4 Y4 X3 Y3 X2 Y2 X1 Y1 1-8.
END INPUT PROGRAM.
```

```
PRINT /X1 TO X4 Y1 TO Y4.
BEGIN DATA
49382716
49382716
49382716
END DATA.
```

- The short form of *VECTOR* is used to establish the dictionary order of a group of variables before they are defined on a *DATA LIST* command. To predetermine variable order, both *VECTOR* and *DATA LIST* must be enclosed within the *INPUT PROGRAM* and *END INPUT PROGRAM* commands.
- The order of the variables in the active dataset will be *X1*, *X2*, *X3*, and *X4*, and *Y1*, *Y2*, *Y3*, and *Y4*, even though they are defined in a different order on *DATA LIST*.
- The program reads the variables with the F1 format specified on *DATA LIST*. It writes the variables with the output format assigned on *VECTOR* (F8.2).



- Another method for predetermining variable order is to use `NUMERIC` (or `STRING` if the variables are string variables) before the `DATA LIST` command (see the example on variable order for the `NUMERIC` command on p. 1396). The advantage of using `NUMERIC` or `STRING` is that you can assign mnemonic names to the variables.

### ***Name Conflicts in Vector Assignments***

```
INPUT PROGRAM.
NUMERIC MIN MINI_A MINI_B MINIM(F2).
COMPUTE MINI_A = MINI(2). /*MINI is function MINIMUM.
VECTOR MINI(3,F2).
DO REPEAT I = 1 TO 3.
+ COMPUTE MINI(I) = -I.
END REPEAT.
COMPUTE MIN = MIN(1). /*The second MIN is function MINIMUM.
COMPUTE MINI_B = MINI(2). /*MINI now references vector MINI
COMPUTE MINIM = MINIM(3). /*The second MINIM is function MINIMUM.
END CASE.
END FILE.
END INPUT PROGRAM.
```

- In this example, there are potential name conflicts between the scalars (the variables named on `NUMERIC`), the vectors (named on `VECTOR`), and the statistical function `MINIMUM`.
- A name that is not followed by a left parenthesis is treated as a scalar.
- When a name followed by a left parenthesis may refer to a vector element or a function, precedence is given to the vector.

## ***VECTOR outside a Loop Structure***

`VECTOR` is most commonly associated with the loop structure, since the index variable for `LOOP` can be used as the subscript. However, the subscript can come from elsewhere, including from the data.

### ***Example***

```
* Create a single case for each of students 1, 2, and 3.

DATA LIST /STUDENT 1 SCORE 3-4 TESTNUM 6.
BEGIN DATA
1 10 1
1 20 2
1 30 3
1 40 4
2 15 2
2 25 3
3 40 1
3 55 3
3 60 4
END DATA.

VECTOR RESULT(4).
COMPUTE RESULT(TESTNUM)=SCORE.

AGGREGATE OUTFILE=* /BREAK=STUDENT
/RESULT1 TO RESULT4=MAX(RESULT1 TO RESULT4).

PRINT FORMATS RESULT1 TO RESULT4 (F2.0).
```

```
PRINT /STUDENT RESULT1 TO RESULT4.  
EXECUTE.
```

- Data are scores on tests recorded in separate cases along with a student identification number and a test number. In this example, there are four possible tests for three students. Not all students took every test.
- The vector *RESULT* creates the variables *RESULT1* through *RESULT4*.
- For each case, *COMPUTE* assigns the *SCORE* value to one of the four vector variables, depending on the value of *TESTNUM*. The other three vector variables for each case keep the system-missing value they were initialized to.
- Aggregating by the variable *STUDENT* creates new cases, as shown by the output from the *PRINT* command. The *MAX* function in *AGGREGATE* returns the maximum value across cases with the same value for *STUDENT*. If a student has taken a particular test, the one valid value is returned as the value for the variable *RESULT1*, *RESULT2*, *RESULT3*, or *RESULT4*.

**Figure 267-1**

*PRINT* output after aggregating

```
1 10 20 30 40  
2 . 15 25 .  
3 40 . 55 60
```

# VERIFY

```
VERIFY [VARIABLES=series name]
```

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
VERIFY.
```

## **Overview**

VERIFY produces a report on the status of the most current DATE, USE, and PREDICT specifications. The report lists the first and last observations in the active dataset, the current USE and PREDICT ranges, and any anomalies in the DATE variables. The number of missing values and the values of the first and last observations in the file and in the USE and PREDICT ranges can also be displayed for a specified series.

VERIFY should be used before a time series procedure whenever there is a possibility that DATE variables or USE and PREDICT ranges have been invalidated. In particular, the active dataset should be verified after you have modified the file structure with commands such as SELECT IF, SORT CASES, and AGGREGATE.

## **Options**

If a series is specified after VERIFY, the values of the first and last observations in the file and in the USE and PREDICT periods are reported for that series. In addition, the number of observations in the active dataset that have missing values for that series is displayed. This can be useful for determining the USE ranges that do not include any missing values.

## **Basic Specification**

The basic specification is the command keyword VERIFY.

- VERIFY displays the first and last observations in the active dataset and in the USE and PREDICT ranges. This information is presented by case number and by the values of the DATE variables.
- For DATE variables, VERIFY reports the number of missing or invalid values. In addition, DATE variables that are not properly nested within the next higher-level DATE variable, that have start or end values other than those expected at the beginning or end of a cycle, or that increment by more than the expected increment are flagged with an asterisk next to the problem. An explanation of the problem is given.

## **Operations**

- VERIFY reports on cases remaining after any preceding SELECT IF commands.

- The `USE` and `PREDICT` ranges are defined by the last `USE` and `PREDICT` commands specified before the `VERIFY` command. If `USE` and `PREDICT` have not been specified, the `USE` range is the entire series, and the `PREDICT` range does not exist.

**Limitations**

- A maximum of 1 `VARIABLES` subcommand. Only one series can be specified on `VARIABLES`.

**VARIABLES Subcommand**

`VARIABLES` names a series to include in the report and is optional. The actual keyword `VARIABLES` can be omitted.

- The series named on `VARIABLES` must be numeric. The `DATE_` series is non-numeric and cannot be specified.
- Only one `VARIABLES` subcommand can be specified, and it can name only one series.

**Examples**

```
VERIFY.
```

- This command produces a report on the status of the most recent `DATE`, `USE`, and `PREDICT` specifications, as well as the first and last valid cases in the file.

```
VERIFY VARIABLE=STOCK.
```

- In addition to the default `VERIFY` information, this command displays information on the series `STOCK`, including the values of the first and last cases and how many values in that series are missing.

# WEIGHT

```
WEIGHT {BY varname}
       {OFF }
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
WEIGHT BY V1.
```

## **Overview**

WEIGHT gives cases different weights (by simulated replication) for statistical analysis. WEIGHT can be used to weight a sample up to population size for reporting purposes or to replicate an example from a table or other aggregated data. With WEIGHT, you can arithmetically alter the sample size or its distribution.

To apply weights resulting from your sampling design, see the Complex Samples option.

## **Basic Specification**

The basic specification is keyword BY followed by the name of the weight variable. Cases are weighted according to the values of the specified variable.

## **Syntax Rules**

- Only one numeric variable can be specified. The variable can be a precoded weighting variable, or it can be computed with the transformation language.
- WEIGHT cannot be placed within a FILE TYPE-END FILE TYPE or INPUT PROGRAM-END INPUT PROGRAM structure. It can be placed nearly anywhere following these commands in a transformation program. [For more information, see the topic Commands and Program States on p. 2196.](#)

## **Operations**

- Unlike most transformations, WEIGHT takes effect as soon as it is encountered in the command sequence. Thus, special attention should be paid to its position among commands. [For more information, see the topic Command Order on p. 41.](#)
- Weighting is permanent during a session unless it is preceded by a TEMPORARY command, changed by another WEIGHT command, or turned off with the WEIGHT OFF specification.
- Each WEIGHT command overrides the previous one.
- WEIGHT uses the value of the specified variable to arithmetically replicate cases for subsequent procedures. Cases are not physically replicated.

- Weight values do not need to be integer, and some procedures, such as `FREQUENCIES`, `CROSSTABS`, and `CTABLES`, will use fractional values on the `WEIGHT` variable. However, most procedures treat the `WEIGHT` variable as a replication weight and will simply round fractional weights to the nearest integer. Some procedures ignore the `WEIGHT` variable completely, and this limitation is noted in the procedure-specific documentation.
- Cases with missing or nonpositive values for the weighting variable are treated as having a weight of 0 and are thus invisible to statistical procedures. They are not used in calculations even where unweighted counts are specified. These cases do remain in the file, however, and are included in case listings and saved when the file is saved.
- A file saved when weighting is in effect maintains the weighting.
- If the weighted number of cases exceeds the sample size, tests of significance are inflated; if it is smaller, they are deflated.

## Examples

```
WEIGHT BY V1.  
FREQ VAR=V2.
```

- The frequency counts for the values of variable *V2* will be weighted by the values of variable *V1*.

```
COMPUTE WVAR=1.  
IF (GROUP EQ 1) WVAR=.5.  
WEIGHT BY WVAR.
```

- Variable *WVAR* is initialized to 1 with the `COMPUTE` command. The `IF` command changes the value of *WVAR* to 0.5 for cases where *GROUP* equals 1.
- Subsequent procedures will use a case base in which cases from group 1 count only half as much as other cases.

# WLS

WLS is available in the Regression option.

WLS VARIABLES= dependent varname WITH independent varnames

```
[/SOURCE=varname]
[/DELTA={1.0**          }]
      {value list      }
      {value TO value BY value}
[/WEIGHT=varname]
[/{CONSTANT**}
 {NOCONSTANT}
[/PRINT={BEST}
 {ALL }
[/SAVE = WEIGHT]
[/APPLY [= 'model name' ]]
```

**\*\***Default if the subcommand or keyword is omitted.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
WLS VARIABLES = VARY WITH VARX VARZ
  /SOURCE=VARZ
  /DELTA=2.
```

## Overview

WLS (weighted least squares) estimates regression models with different weights for different cases. Weighted least squares should be used when errors from an ordinary regression are heteroscedastic—that is, when the size of the residual is a function of the magnitude of some variable, termed the **source**.

The WLS model is a simple regression model in which the residual variance is a function of the source variable, up to some power transform indicated by a delta value. For fuller regression results, save the weights produced by WLS and specify that weight variable on the REGWGT subcommand in REGRESSION.

## Options

**Calculated and Specified Weights.** WLS can calculate the weights based on a source variable and delta values (subcommands SOURCE and DELTA), or it can apply existing weights contained in a series (subcommand WEIGHT). If weights are calculated, each weight value is calculated as the source series value raised to the negative delta value.

**New Variables.** You can change `NEWVAR` settings on the `TSET` command prior to `WLS` to evaluate the regression coefficients and log-likelihood function without saving the weight variable, or save the new values to replace the values saved earlier, or save the new values without erasing values saved earlier (see the `TSET` command). You can also use the `SAVE` subcommand on `WLS` to override the `NONE` or the default `CURRENT` settings on `NEWVAR` for the current procedure.

**Statistical Output.** You can change the `PRINT` setting on the `TSET` command prior to `WLS` to display regression coefficients or the list of log-likelihood functions at each delta value, or to limit the output to only the regression statistics for the delta value at which the log-likelihood function is maximized (see the `TSET` command). You can also use the `PRINT` subcommand to override the `PRINT` setting on the `TSET` command for the current procedure and obtain regression coefficients at each value of delta in addition to the default output.

### **Basic Specification**

- The basic specification is the `VARIABLES` subcommand specifying one dependent variable, the keyword `WITH`, and one or more independent variables. Weights are calculated using the first independent variable as the source variable and a default delta value of 1.
- The default output for calculated weights displays the log-likelihood function for each value of delta. For the value of delta at which the log-likelihood function is maximized, the displayed summary regression statistics include  $R$ ,  $R^2$ , adjusted  $R^2$ , standard errors, analysis of variance, and  $t$  tests of the individual coefficients. A variable named `WGT#I` containing the calculated weights is automatically created, labeled, and added to the active dataset.

### **Syntax Rules**

- `VARIABLES` can be specified only once.
- `DELTA` can be specified more than once. Each specification will be executed.
- If other subcommands are specified more than once, only the last specification of each one is executed.
- You can specify either `SOURCE` and `DELTA`, or just the `WEIGHT` subcommand. You cannot specify all three, and you cannot specify `WEIGHT` with `SOURCE` or with `DELTA`.

### **Subcommand Order**

- Subcommands can be specified in any order.

### **Operations**

- If neither the `WEIGHT` subcommand nor the `SOURCE` and `DELTA` subcommands are specified, a warning is issued and weights are calculated using the default source and delta value.
- Only one `WGT#I` variable is created per procedure. If more than one delta value is specified, the weights used when the log-likelihood function is maximized are the ones saved as `WGT#I`.
- `WGT#I` is not created when the `WEIGHT` subcommand is used.
- The `WEIGHT` command specifies case replication weights, which are *not* the same as the weights used in weighted least squares. If the `WEIGHT` command and `WLS WEIGHT` subcommand are both specified, both types of weights are incorporated in `WLS`.



- WLS uses listwise deletion of missing values. Whenever one variable is missing a value for a particular observation, that observation will not be included in any computations.

### **Limitations**

- Maximum one VARIABLES subcommand.
- Maximum one dependent variable on the VARIABLES subcommand. There is no limit on the number of independent variables.
- Maximum 150 values specified on the DELTA subcommand.

## **Example**

```
WLS cost WITH sqft inorout yrexp
  /SOURCE sqft
  /POWER 0 TO 5 BY 0.5
  /CONSTANT
  /SAVE WEIGHT
  /PRINT BEST.
```

- This command specifies a weighted least-squares regression in which *cost* is the dependent variable and *sqft*, *inorout*, and *yrexp* are the independent variables.
- The SOURCE subcommand identifies *sqft* as the source of heteroscedasticity.
- The POWER subcommand specifies that several equations will be fit, one for each value from 0 to 5, incrementing by 0.5. Weights will be calculated using *sqft*.
- The SAVE subcommand requests that the weights that maximize the log-likelihood function be saved to the active dataset.
- All other options are set to their default values.

## **VARIABLES Subcommand**

VARIABLES specifies the variable list and is the only required subcommand.

## **SOURCE Subcommand**

SOURCE is used in conjunction with the DELTA subcommand to compute weights. SOURCE names the variable that is the source of heteroscedasticity.

- The only specification on SOURCE is the name of a variable to be used as the source of heteroscedasticity.
- Only one source variable can be specified.
- If neither SOURCE nor WEIGHT is specified, the first independent variable specified on the VARIABLES subcommand is assumed to be the source variable.

## ***DELTA Subcommand***

DELTA, alias POWER, is used in conjunction with the SOURCE subcommand to compute weights. DELTA specifies the values to use in computing weights. The weights are equal to  $1/(\text{SOURCE raised to the DELTA power})$ .

- The specification on DELTA is a list of possible delta values and/or value grids.
- Multiple values and grids can be specified on one DELTA subcommand.
- Delta values can be any value in the range of  $-6.5$  to  $+7.5$ . Values below this range are assigned the minimum ( $-6.5$ ), and values above are assigned the maximum ( $7.5$ ).
- A grid is specified by naming the starting value, the keyword TO, an ending value, the keyword BY, and an increment value. Alternatively, the keyword BY and the increment value can be specified after the starting value.
- More than one DELTA subcommand can be specified; each subcommand will be executed.
- If DELTA is not specified, the delta value defaults to 1.0.

### ***Example***

```
WLS VARIABLES = X1 WITH Y1 Z1
  /SOURCE=Z1
  /DELTA=0.5.
```

- In this example, weights are calculated using the source variable *Z1* and a delta value of 0.5. Thus, the weights are  $1 / (\text{SQRT} (Z1))$ .

### ***Example***

```
WLS VARIABLES = SHARES WITH PRICE
  /DELTA=0.5 TO 2.5 BY 0.5.
```

- In this example, several regression equations will be fit, one for each value of delta.
- Weights are calculated using the source variable *PRICE* (the default).
- The delta values start at 0.5 and go up to 2.5, incrementing by 0.5. This specification is equivalent to 0.5 BY 0.5 TO 2.5.
- The weights that maximize the log-likelihood function will be saved as variable *WGT#1*.

## ***WEIGHT Subcommand***

WEIGHT specifies the variable containing the weights to be used in weighting the cases. WEIGHT is an alternative to computing the weights using the SOURCE and DELTA subcommands. If a variable containing weights is specified, the output includes the regression coefficients, log-likelihood function, and summary regression statistics such as  $R$ ,  $R^2$ , adjusted  $R^2$ , standard errors, analysis of variance, and  $t$  tests of the coefficients. Since no new weights are computed, no new variable is created.

- The only specification on WEIGHT is the name of the variable containing the weights. Typically, *WGT* variables from previous WLS procedures are used.
- Only one variable can be specified.

**Example**

```
WLS VARIABLES = SHARES WITH PRICE
  /WEIGHT=WGT_1.
```

- This WLS command uses the weights contained in variable *WGT\_1* to weight cases.

**CONSTANT and NOCONSTANT Subcommands**

Specify `CONSTANT` or `NOCONSTANT` to indicate whether a constant term should be estimated in the regression equation. The specification of either subcommand overrides the `CONSTANT` setting on the `TSET` command for the current procedure.

- `CONSTANT` is the default and specifies that the constant term is used as an instrument.
- `NOCONSTANT` eliminates the constant term.

**SAVE Subcommand**

`SAVE` saves the weight variable generated during the current session to the end of the active dataset. The default name *WGT\_n* will be generated, where *n* increments to make the variable name unique. The only specification on `SAVE` is `WEIGHT`. The specification overrides the `NONE` or the default `CURRENT` setting on `NEWVAR` for the current procedure.

**PRINT Subcommand**

`PRINT` can be used to override the `PRINT` setting on the `TSET` command for the current procedure. Two keywords are available.

**BEST**                    *Display coefficients for the best weight only.* This is the default.  
**ALL**                     *Display coefficients for all weights.*

**APPLY Subcommand**

- The `APPLY` subcommand allows you to use a previously defined WLS model without having to repeat the specifications.
- The only specification on `APPLY` is the name of a previous model in quotes. If a model name is not specified, the model specified on the previous WLS command is used.
- To change one or more model specifications, specify the subcommands of only those portions you want to change after the `APPLY` subcommand.
- If no variables are specified on the command, the variables that were originally specified with the model being reapplied are used.

**Example**

```
WLS VARIABLES = X1 WITH Y1
  /SOURCE=Y1
  /DELTA=1.5.
WLS APPLY
```

```
/DELTA=2 .
```

- The first command produces a weighted least-squares regression of  $X1$ , with  $Y1$  as the source variable and delta equal to 1.5.
- The second command uses the same variable and source but changes the delta value to 2.

**Example**

```
WLS VARIABLES = X1 WITH Y1 Z1  
  /SOURCE=Z1  
  /DELTA=1 TO 3 BY 0.5  
WLS APPLY  
  /WEIGHT=WGT#1 .
```

- The first command regresses  $X1$  on  $Y1$  and  $Z1$ , using  $Z1$  as the source variable. The delta values range from 1 to 3, incrementing by 0.5.
- The second command again regresses  $X1$  on  $Y1$  and  $Z1$ , but this time applies the values of  $WGT\#1$  as the weights.

# WRITE

```
WRITE [OUTFILE='file'] [ENCODING='encoding specification']  
  [RECORDS={1}] [{NOTABLE}]  
      {n}    {TABLE }  
  
/{1    } varlist [{col location [(format)]}] [varlist...]  
{rec #}          {(format list)}           }  
                  {*                       }  
  
[/{2    }...]  
{rec #}
```

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Release History**

Release 16.0

- ENCODING subcommand added for Unicode support.

## **Example**

```
WRITE OUTFILE="/data/personnel.txt" / MOHIRED YRHIRED DEPT SALARY NAME.  
EXECUTE.
```

## **Overview**

WRITE writes files in a machine-readable format that can be used by other software applications. When used for this purpose, the OUTFILE subcommand is required. If OUTFILE is not specified, the output from WRITE that can be displayed is included with the output from your session in a format similar to that used by the PRINT command.

### **Options**

**Formats.** You can specify formats for the variables.

**Strings.** You can include strings within the variable specifications. The strings can be used to label values or to add extra space between values.

**Multiple Lines per Case.** You can write variables on more than one line for each case. See the RECORDS subcommand.

**Output File.** You can direct the output to a specified file using the OUTFILE subcommand.

**Summary Table.** You can display a table that summarizes the variable formats with the TABLE subcommand.

### ***Subcommand Order***

Subcommands can be specified in any order. However, all subcommands must be used before the slash that precedes the first variable list.

### ***Basic Specification***

The basic specification is a slash followed by a variable list. The values for all of the variables specified on the list are included with the rest of the output from your session.

### ***Syntax Rules***

- A slash must precede the variable specifications. The first slash begins the definition of the first (and possibly only) line per case of the WRITE output.
- Specified variables must already exist, but they can be numeric, string, scratch, temporary, or system variables. Subscripted variable names, such as  $X(I)$  for the first element in vector  $X$  cannot be used.
- Keyword ALL can be used to write the values of all user-defined variables in the active dataset.

### ***Operations***

- WRITE is executed once for each case constructed from the data file.
- Values are written to the file as the data are read.
- WRITE is a transformation and will not be executed unless it is followed by a procedure or the EXECUTE command.
- When writing to an external file with the OUTFILE subcommand, line/record width can be up to 2.1 billion bytes. When writing to the Viewer, however (when there is no OUTFILE subcommand), if a line width exceeds the line width defined by SET WIDTH, an error results and the WRITE command is not executed. The maximum line width you can specify with SET WIDTH is 255 bytes.
- There are no carriage control characters in the output file generated by WRITE.
- User-missing values are written just like valid values. System-missing values are represented by blanks.
- If you are writing a file to be used on another system, you should take into account that some data types cannot be read by all computers.
- If long records are less convenient than short records with multiple records per case, you can write out a case identifier and insert a string as a record identification number. The receiving system can then check for missing record numbers (see Strings on p. 2162 for an example).

## ***Examples***

### ***Writing a Specific Set of Variables***

```
WRITE OUTFILE="/data/personnel.txt" / MOHIRED YRHIRED DEPT SALARY NAME.  
FREQUENCIES VARIABLES=DEPT.
```

- `WRITE` writes values for each variable on the variable list to file *personnel.txt*. The `FREQUENCIES` procedure reads the data and causes `WRITE` to be executed.
- All variables are written with their dictionary formats.

### Writing All Variables

```
WRITE OUTFILE='/data/personnel.txt' /ALL.
EXECUTE.
```

- `WRITE` writes values for all user-defined variables in the active dataset to file *personnel.txt*. The `EXECUTE` command executes `WRITE`.

## Formats

By default, `WRITE` uses the dictionary write formats. You can specify formats for some or all variables specified on `WRITE`. For a string variable, the specified format must have the same width as that of the dictionary format.

- Format specifications can be either column-style or FORTRAN-like (see `DATA LIST`). The column location specified with column-style formats or implied with FORTRAN-like formats refers to the column in which the variable will be written.
- A format specification following a list of variables applies to all the variables in the list. Use an asterisk to prevent the specified format from applying to variables preceding the asterisk. The specification of column locations implies a default print format, and that format will apply to all previous variables if no asterisk is used.
- All available formats can be specified on `WRITE`. Note that hex and binary formats use different widths. For example, the `AHEX` format must have a width twice that of the corresponding `A` format. For more information on specifying formats and on the formats available, see `DATA LIST` and Variable Types and Formats on p. 54.
- Format specifications are in effect only for the `WRITE` command. They do not change the dictionary write formats.
- To specify a blank between variables in the output, use a string (see Strings on p. 2162), specify blank columns in the format, or use an `X` or `T` format element in the `WRITE` specifications (see `DATA LIST` for information on `X` and `T`).

### Example

```
WRITE OUTFILE='/data/personnel.txt' / TENURE (F2.0) ' ' MOHIRED YRHIRED DEPT *
      SALARY85 TO SALARY88 (4(DOLLAR8,1X)) NAME.
EXECUTE.
```

- Format `F2.0` is specified for `TENURE`. A blank between quotes is specified as a string after `TENURE` to separate values of `TENURE` from those of `MOHIRED`.
- `MOHIRED`, `YRHIRED`, and `DEPT` are written with default formats because the asterisk prevents them from receiving the `DOLLAR8` format specified for `SALARY85` to `SALARY88`. The `1X` format element is specified with `DOLLAR8` to add one blank after each value of `SALARY85` to `SALARY88`.
- `NAME` uses the default dictionary format.

## Strings

You can specify strings within the variable list. Strings must be enclosed in quotes.

- If a format is specified for a variable list, the application of the format is interrupted by a specified string. Thus, the string has the same effect within a variable list as an asterisk.

### Example

```
WRITE OUTFILE='/data/personnel.txt'
  /EMPLOYID '1' MOHIRED YRHIRED SEX AGE JOBCAT NAME
  /EMPLOYID '2' DEPT86 TO DEPT88 SALARY86 TO SALARY88.
EXECUTE.
```

- Strings are used to assign the constant 1 to record 1 of each case, and 2 to record 2 to provide record identifiers in addition to the case identifier *EMPLOYID*.

## RECORDS Subcommand

RECORDS indicates the total number of lines written per case. The number specified on RECORDS is informational only. The actual specification that causes variables to be written on a new line is a slash within the variable specifications. Each new line is requested by another slash.

- RECORDS must be specified before the slash that precedes the start of the variable specifications.
- The only specification on RECORDS is an integer to indicate the number of records for the output. If the number does not agree with the actual number of records indicated by slashes, the program issues a warning and ignores the specification on RECORDS.
- Specifications for each line of output must begin with a slash. An integer can follow the slash, indicating the line on which values are to be written. The integer is informational only. It cannot be used to rearrange the order of records in the output. If the integer does not agree with the actual record number indicated by the number of slashes in the variable specifications, the integer is ignored.
- A slash that is not followed by a variable list generates a blank line in the output.

### Examples

```
WRITE OUTFILE='/data/personnel.txt' RECORDS=2
  /EMPLOYID NAME DEPT
  /EMPLOYID TENURE SALARY.
EXECUTE.
```

- WRITE writes the values of an individual's name and department on one line, tenure and salary on the next line, and the employee identification number on both lines.

```
WRITE OUTFILE='/data/personnel.txt' RECORDS=2
  /1 EMPLOYID NAME DEPT
  /2 EMPLOYID TENURE SALARY.
EXECUTE.
```

- This command is equivalent to the command in the preceding example.

```
WRITE OUTFILE='/data/personnel.txt' / EMPLOYID NAME DEPT / EMPLOYID TENURE SALARY.
```



EXECUTE.

- This command is equivalent to the commands in both preceding examples.

## ***OUTFILE Subcommand***

OUTFILE specifies the target file for the output from the WRITE command. By default, the output is included with the rest of the output from the session.

- OUTFILE must be specified before the slash that precedes the start of the variable specifications.

### ***Example***

```
WRITE OUTFILE='/data/writeout.txt'  
  /1 EMPLOYID DEPT SALARY /2 NAME.  
EXECUTE.
```

- OUTFILE specifies *writeout.txt* as the file that receives the WRITE output.

## ***ENCODING Subcommand***

ENCODING specifies the encoding format of the file. The keyword is followed by an equals sign and a quoted encoding specification.

- In Unicode mode, the default is UTF8. For more information, see [SET command, UNICODE subcommand](#).
- In code page mode, the default is the current locale setting. For more information, see [SET command, LOCALE subcommand](#).
- The quoted encoding value can be: *Locale* (the current locale setting), UTF8, UTF16, UTF16BE (big endian), UTF16LE (little endian), a numeric Windows code page value (for example, '1252'), or an IANA code page value (for example, 'iso8859-1' or cp1252).
- If there is no OUTFILE subcommand, the ENCODING subcommand is ignored.

## ***TABLE Subcommand***

TABLE requests a table showing how the variable information is formatted. NOTABLE is the default.

- TABLE must be specified before the slash that precedes the start of the variable specifications.

### ***Example***

```
WRITE OUTFILE='/data/personnel.txt' TABLE  
  /1 EMPLOYID DEPT SALARY /2 NAME.  
EXECUTE.
```

- TABLE requests a summary table describing the WRITE specifications.

# WRITE FORMATS

```
WRITE FORMATS varlist (format) [varlist...]
```

This command takes effect immediately. It does not read the active dataset or execute pending transformations. [For more information, see the topic Command Order on p. 41.](#)

## Example

```
WRITE FORMATS SALARY (DOLLAR8)
/ HOURLY (DOLLAR7.2)
/ RAISE BONUS (PCT2).
```

## Overview

WRITE FORMATS changes variable write formats. Write formats are output formats and control the form in which values are written by the WRITE command.

WRITE FORMATS changes only write formats. To change print formats, use the PRINT FORMATS command. To change both the print and write formats with a single specification, use the FORMATS command. For information on assigning input formats during data definition, see DATA LIST. For detailed information on available formats and specifications, see [Variable Types and Formats](#).

## Basic Specification

The basic specification is a variable list followed by the new format specification in parentheses. All specified variables receive the new format.

## Syntax Rules

- You can specify more than one variable or variable list, followed by a format in parentheses. Only one format can be specified after each variable list. For clarity, each set of specifications can be separated by a slash.
- You can use keyword TO to refer to consecutive variables in the active dataset.
- The specified width of a format must include enough positions to accommodate any punctuation characters such as decimal points, commas, dollar signs, or date and time delimiters. (This differs from assigning an *input* format on DATA LIST, where the program automatically expands the input format to accommodate punctuation characters in output.)
- Custom currency formats (CCw, CCw.d) must first be defined on the SET command before they can be used on WRITE FORMATS.
- For string variables, you can only use WRITE FORMATS to switch between A and AHEx formats. WRITE FORMATS cannot be used to change the length of string variables. To change the length of a string variable, declare a new variable of the desired length with the STRING command and then use COMPUTE to copy values from the existing string into the new variable.

**Operations**

- Unlike most transformations, `WRITE FORMATS` takes effect as soon as it is encountered in the command sequence. Special attention should be paid to its position among commands. [For more information, see the topic Command Order on p. 41.](#)
- Variables not specified on `WRITE FORMATS` retain their current formats in the active dataset. To see the current formats, use the `DISPLAY` command.
- The new write formats are changed only in the active dataset and are in effect for the duration of the session or until changed again with a `WRITE FORMATS` or `FORMATS` command. Write formats in the original data file (if one exists) are not changed, unless the file is resaved with the `SAVE` or `XSAVE` command.
- New numeric variables created with transformation commands are assigned default print and write formats of `F8.2` (or the format specified on the `FORMAT` subcommand of `SET`). The `WRITE FORMATS` command can be used to change the new variable's write formats.
- New string variables created with transformation commands are assigned the format specified on the `STRING` command that declares the variable. `WRITE FORMATS` cannot be used to change the format of a new string variable.
- Date and time formats are effective only with the `LIST` and `TABLES` procedures and the `PRINT` and `WRITE` transformation commands. All other procedures use `F` format regardless of the date and time formats specified. [For more information, see the topic Date and Time Formats on p. 63.](#)
- If a numeric data value exceeds its width specification, the program attempts to write some value nevertheless. First the program rounds decimal values, then removes punctuation characters, then tries scientific notation, and finally, if there is still not enough space, produces asterisks indicating that a value is present but cannot be written in the assigned width.

**Examples****Specifying Write Formats for Multiple Variables**

```
WRITE FORMATS SALARY (DOLLAR8)
              / HOURLY (DOLLAR7.2)
              / RAISE BONUS (PCT2) .
```

- The write format for `SALARY` is changed to `DOLLAR` with eight positions, including the dollar sign and comma when appropriate. An eight-digit number would require a `DOLLAR11` format specification: eight characters for the digits, two characters for commas, and one character for the dollar sign.
- The write format for `HOURLY` is changed to `DOLLAR` with seven positions, including the dollar sign, decimal point, and two decimal places.
- The write format for both `RAISE` and `BONUS` is changed to `PCT` with two positions: one for the percentage and one for the percent sign.

**Changing the Default Format of a New Variable**

```
COMPUTE V3=V1 + V2 .
WRITE FORMATS V3 (F3.1) .
```

- COMPUTE creates the new numeric variable *V3*. By default, *V3* is assigned an F8.2 format.
- WRITE FORMATS changes the write format for *V3* to F3.1.

### **Working With Custom Currency Formats**

```
SET CCA='-/- .Df1 . .-' .  
WRITE FORMATS COST (CCA14.2) .
```

- SET defines a European currency format for the custom currency format type CCA.
- WRITE FORMATS assigns the write format CCA to variable *COST*. See the SET command for more information on custom currency formats.

# XGRAPH

XGRAPH is available in the Statistics Base option.

XGRAPH is available only on systems with high-resolution graphics capabilities.

*Note:* Square brackets used in the XGRAPH syntax chart are required parts of the syntax and are not used to indicate optional elements. Any equals signs (=) displayed in the syntax chart are required. All subcommands except the chart expression in the first line are optional.

```
XGRAPH CHART=yvars BY xvars BY zvars

  /BIN START={AUTO**} SIZE={AUTO** }
      {x      }           {WIDTH (x)}
                        {COUNT (n)}

  /DISPLAY DOT={ASYMMETRIC**}
      {SYMMETRIC  }
      {FLAT      }

  /DISTRIBUTION TYPE=NORMAL

  /COORDINATE SPLIT={NO**}
      {YES  }

  /ERRORBAR {CI      {(95)}}
      { (n) }
      {STDDEV {(2)}}
      { (n) }
      {SE      {(2)}}
      { (n) }

  /MISSING USE={LISTWISE** } REPORT={NO**}
      {VARIABLEWISE}           {YES  }

  /PANEL COLVAR=varlist COLOP={CROSS**} ROWVAR=varlist ROWOP={CROSS**}
      {NEST  }                               {NEST  }

  /TEMPLATE FILE='filespec'

  /TITLES TITLE='line' 'line2' SUBTITLE='line1'
      FOOTNOTE='line1' 'line2'
```

**\*\*** Default if the subcommand is omitted.

yvars defines the y dimension and has the general form:

```
(varname [function] [data_element_type] + ...) > {varname {[COLOR]  }}
                                                    {[PATTERN]}
                                                    {1      {[COLOR]  }}
                                                    {[PATTERN]}
```

xvars and zvars define the x and z dimensions and have the general form:

```
{varname {[c]}                } > varname {[COLOR]  }
      {[s]}                }           {[PATTERN]}
{$CASENUM [LABEL=varname]  }
{1      }                        }
```

The + operator blends the y-axis to accommodate all the y-axis variables. It is used when separate variables are displayed on the same axis. The number 1 is a placeholder for the variables on the axis. The > operator stacks data elements on the y-axis and clusters data elements on the x-axis

and z-axis. [c] or [s] indicates whether the variable should be treated as categorical or scale. \$CASENUM is used when individual cases are displayed on the same axis.

Summary functions available for *yvars*: VALUE, COUNT, PCT, CUPCT, CUFREQ, MINIMUM, MAXIMUM, VALIDN, SUM, CSUM, MEAN, STDDEV, VARIANCE, MEDIAN, GMEDIAN, MODE, PTILE(x), GPTILE(x), PLT(x), PGT(x), NLT(x), NGT(x), PIN(x1, x2), and NIN(x1, x2).

Data element types available for *yvars*: BAR, HISTOBAR, and POINT.

This command reads the active dataset and causes execution of any pending commands. [For more information, see the topic Command Order on p. 41.](#)

### **Release History**

Release 13.0

- Command introduced.

### **Example**

```
XGRAPH CHART=( [COUNT] [BAR] ) BY jobcat [c] BY gender [c] .
```

## **Overview**

XGRAPH generates a high-resolution chart by computing statistics from variables in the active dataset and constructing the chart according to your specification. The chart can be a 3-D bar chart, population pyramid, or dot plot. The chart is displayed where high-resolution display is available and can be edited with the Chart Editor and saved in an Output file.

*Note:* Although you can create other charts with XGRAPH, these are not currently supported.

### **Basic Specification**

The basic specification is the command name followed by a chart expression. The chart expression must identify one or more variables and a data element for the chart.

### **Syntax Rules**

- The chart expression is required and must appear first. It can optionally be preceded by CHART=. If CHART is used, the = must be present. All other subcommands are optional.
- After the chart expression, subcommands can appear in any order.
- Only a single instance of each subcommand is allowed.
- Subcommand names and keywords must be spelled in full.
- Required square brackets cannot be merged for different types of specifications. For example, when a summary function and a data element are adjacent, they cannot be merged. That is, [VALUE] [BAR] cannot be written as [VALUE BAR].

**Limitations**

- XGRAPH does not produce charts with more than 1,000,000 data values.
- When a weight is in effect and the `ERRORBAR` subcommand is used on a chart that displays medians, counts, or percentages, the case weights must be integers.

**CHART Expression**

The **chart expression** defines the data and the role of that data in the chart. It also specifies the data element type and any clustering and stacking. For clustered and stacked charts, the chart expression specifies whether color or patterns are used to distinguish the clustered and stacked groups. For charts of individual cases (that is, when `$CASENUM` defines a categorical axis), the chart expression can also specify a label variable.

- The `CHART` keyword is optional. If it is used, it must be followed by an equals sign (=).
- The order within the chart expression matters. The `yvars` (the *y* dimension) must be first, followed by `xvars` and then `zvars`.
- The final `BY` variable is used as the splitter variable in a population pyramid when the `COORDINATE` subcommand specifies `SPLIT`.
- There is a limit to the number of stack and cluster specifications. There can be two clusters, or one stack and one cluster.

**Functions**

Functions must be enclosed in square brackets.

*Value function:*

The `VALUE` function yields the value of the specified *y*-axis variable for each case. It always produces one data element for each case.

*Aggregation functions:*

Two groups of aggregation functions are available: count functions and summary functions.

*Count functions:*

<b>COUNT</b>	<i>Frequency of cases in each category.</i>
<b>PCT</b>	<i>Frequency of cases in each category expressed as a percentage of the whole.</i>
<b>CUPCT</b>	<i>Cumulative percentage sorted by category value.</i>
<b>CUFREQ</b>	<i>Cumulative frequency sorted by category value.</i>

- Count functions yield the count or percentage of valid cases within categories determined by the `xvars` and the `zvars`, as in:  

```
XGRAPH ([PCT] [BAR]) BY jobcat [c] BY gender [c].
```
- Count functions do not operate on a variable.
- Count functions do not have any arguments.

*Summary functions:*

<b>MINIMUM</b>	<i>Minimum value of the variable.</i>
<b>MAXIMUM</b>	<i>Maximum value of the variable.</i>
<b>VALIDN</b>	<i>Number of cases for which the variable has a nonmissing value.</i>
<b>SUM</b>	<i>Sum of the values of the variable.</i>
<b>CUSUM</b>	<i>Sum of the summary variable accumulated across values of the category variable.</i>
<b>MEAN</b>	<i>Mean.</i>
<b>STDDEV</b>	<i>Standard deviation.</i>
<b>VARIANCE</b>	<i>Variance.</i>
<b>MEDIAN</b>	<i>Median.</i>
<b>GMEDIAN</b>	<i>Group median.</i>
<b>MODE</b>	<i>Mode.</i>
<b>PTILE(x)</b>	<i>Xth percentile value of the variable. X must be greater than 0 and less than 100.</i>
<b>PLT(x)</b>	<i>Percentage of cases for which the value of the variable is less than x.</i>
<b>PGT(x)</b>	<i>Percentage of cases for which the value of the variable is greater than x.</i>
<b>NLT(x)</b>	<i>Number of cases for which the value of the variable is less than x.</i>
<b>NGT(x)</b>	<i>Number of cases for which the value of the variable is greater than x.</i>
<b>PIN(x1,x2)</b>	<i>Percentage of cases for which the value of the variable is greater than or equal to x1 and less than or equal to x2. x1 cannot exceed x2.</i>
<b>NIN(x1,x2)</b>	<i>Number of cases for which the value of the variable is greater than or equal to x1 and less than or equal to x2. x1 cannot exceed x2.</i>

- Summary functions operate on summary variables (variables that record continuous values, like age or expenses). To use a summary function, specify the name of one or more variables before the name of the function and the data element type, both of which are in square brackets, as in:

```
XGRAPH (salary [SUM] [BAR]) BY jobcat[c] BY gender[c].
```

- You can specify a list of variables on which the function operates, as in:

```
XGRAPH ((salary + salbegin) [SUM] [BAR]) BY jobcat[c] BY gender[c].
```

This syntax is equivalent to:

```
XGRAPH (salary [SUM] [BAR] + salbegin [SUM] [BAR]) BY jobcat[c] BY gender[c].
```

## **Data Element Types**

The data element type is a basic description of how the data are represented on the chart.

<b>BAR</b>	<i>Draws a bar data element. Bars show the results of a summary function. Used for 3-D bar charts and population pyramids that show the distribution of a categorical variable.</i>
<b>HISTOBAR</b>	<i>Draws a histogram data element. Histogram data elements show the counts in a range of cases. A range of cases is called a bin. Used for population pyramids.</i>
<b>POINT</b>	<i>Draws a point marker. A point marker shows a case with a specific x-axis value. Used for dot plots.</i>



## Measurement Level

The chart expression allows the user to specify a measurement level for BY variables. If measurement level is not specified, the following occurs:

- In a chart of BAR elements, `xvars` and `zvars` are treated as categorical.
- In a chart of HISTOBAR elements, `xvars` is treated as scale. `zvars` is treated as categorical, because the COORDINATE subcommand and SPLIT keyword are always used.
- In charts of POINT elements, the measurement level is obtained from the IBM® SPSS® Statistics data dictionary for each variable. If no measurement levels have been defined, numeric variables are treated as scale and string variables are treated as categorical.

[c] *Specifies that the variable is treated as a categorical variable.*

[s] *Specifies that the variable is treated as a scale variable.*

## Variable Placeholder

When summaries of separate variables are used to define a categorical axis, clustering, or stacking, 1 must be used as a variable placeholder.

- 1 *Uses the separate variables specified in `yvars` as categories on the axis, cluster categories, or stack categories. Must be defined as an integer (no decimal places).*

### Examples

```
XGRAPH CHART=((salary + salbegin) [MEAN] [BAR]) BY 1 BY gender[c].
```

- Separate variables appear on the y-axis.

```
XGRAPH CHART=((salary + salbegin) [MEAN] [BAR]) BY gender[c] BY 1.
```

- Separate variables appear on the z-axis.

## Case Numbers

Whenever the VALUE function is used, one categorical axis consists of cases, which are identified by a sequential number.

- \$CASENUM** *Uses the sequential numbering of selected cases as the categories on the axis. This is not the actual case number.*

### Example

```
XGRAPH CHART=(salary [VALUE]) BY gender[c] BY $CASENUM.
```

## Blending, Clustering, and Stacking

Blending on an axis is indicated by +. This is used when separate variables are displayed on the same axis.

Clustering and stacking are indicated by >. On the *x*- or *z*-axis, this indicates clustering. On the *y*-axis, this indicates stacking. Stacking is allowed with one clustering variable; or two clustering variables are allowed, one on the *x*-axis and one on the *z*-axis.

- + Scales the axis to accommodate all of the variables on that dimension.
- > Stacks groups on the *y*-axis, and clusters groups on the *x*- and *z*-axes.

### **Stacking Example**

```
XGRAPH CHART=(salary [MEAN] [BAR]) > educ BY jobcat[c] by gender[c].
```

### **Clustering Example**

```
XGRAPH CHART=(salary [MEAN] [BAR]) BY jobcat[c] > educ by gender[c].
```

When you use clustering or stacking, you can specify how the clusters and stacks are differentiated in the chart.

- [COLOR] Uses color to differentiate the clusters or stacks.
- [PATTERN] Uses pattern to differentiate the clusters or stacks.

[COLOR] and [PATTERN] follow these rules:

- The actual colors and patterns are specified in the Options. You can access these by choosing Options from the Edit menu. Then click the Charts tab.
- [COLOR] and [PATTERN] follow the stacking or clustering variable in the chart expression.
- If the specification is omitted, color is used to differentiate the clusters or stacks. The color cycle starts over for each dimension.
- If neither [COLOR] nor [PATTERN] is specified, XGRAPH uses the Style Cycle Preference specified by the Options.
- If [COLOR] or [PATTERN] is specified for both dimensions, these specifications are honored. The color or pattern cycle starts over for each dimension.
- If [COLOR] or [PATTERN] is specified for only one dimension, the specification is honored, and the other dimension uses the other specification.

### **Example**

```
XGRAPH CHART=(salary [MEAN] [BAR]) BY jobcat[c] > educ[PATTERN]
by gender[c].
```

## **Labels**

On categorical charts of individual cases, you can specify a labeling variable by using [LABEL=varname] in the chart expression. [LABEL=varname] follows \$CASENUM in the chart expression.

- [LABEL=varname] The variable whose values are used as the tick labels instead of the case number.

**Example**

```
XGRAPH CHART=(salary [VALUE] [BAR]) BY $CASENUM[LABEL=gender]
> jobcat BY educ.
```

**BIN Subcommand**

For population pyramids, the `BIN` subcommand controls the starting point and size of the bins. The bins are the bars in the chart and represent a group of data values. For example, if there are four bins in a chart and the values range from 0-100, the bin width is 25. Therefore, all points with values of 0-24 would be in the first bin, those with values of 25-49 would be in the second bin, and so on. Each keyword is followed by an equals sign (=) and the value for that keyword.

**Example**

```
XGRAPH CHART=( [HISTOBAR] ) BY age[s] BY gender[c]
/COORDINATE SPLIT=YES
/BIN START=30 SIZE=WIDTH(5).
```

- The first bin begins at age 30.
- Each bin includes cases with *age* values that span 5 years.
- So the first bin contains the cases with ages 30-34, the second bin contains the cases with ages 35-39, and so on.

**START Keyword**

The `START` keyword specifies the starting value of the first bin in the chart.

- AUTO** *Adjust the starting value of the first bin according to the data values.* The starting value is set so the first bin includes the lowest data value and the bin boundaries are at good values. This is the default.
- value** *User-specified starting value.* The value can be set lower than the lowest data value so the first bin includes values that are not in the dataset. For example, if the lowest value in the dataset is 6, the starting value can be set to 0 so the values of 0–5 are included in the first bin.
- Note:* If the variable being binned is a date, the starting value must be a data literal in quotes and in the date format specified for the variable in the IBM® SPSS® Statistics data dictionary (for example, 'January 1, 2001').

**SIZE Keyword**

The `SIZE` keyword specifies the width of the bins. The width can be determined by an exact value or by the number of bins (fewer bins result in wider bins).

- AUTO** *Adjust bin width according to the data values.* The size is set so the bins are neither too few nor too narrow. This is the default.
- WIDTH (value)** *User-specified value for the bin width.*

*Note:* If the variable being binned is a date, the width must be specified in quotes and in the duration format ddd hh:mm (for example, '30 12:00' for bin widths of 30 days and 12 hours). The hh:mm portion of the duration format is optional.

**COUNT (value)** *User-specified value for the number of bins.*

## **DISPLAY Subcommand**

DISPLAY controls how the data elements are displayed. Currently this subcommand applies to dot plots.

### **DOT Keyword**

The DOT keyword controls the display of points in a dot plot.

**ASYMMETRIC** *Stack the points on the x-axis. This is the default.*

**SYMMETRIC** *Arrange the points symmetrically across a central horizontal line in the chart. In other words, the point stacks are vertically centered.*

**FLAT** *Do not stack the points.*

## **DISTRIBUTION Subcommand**

The DISTRIBUTION subcommand adds distribution curves to the chart. This subcommand applies to population pyramids.

### **Example**

```
XGRAPH CHART=( [HISTOGRAM] ) BY age[s] BY gender[c]
/DISTRIBUTION TYPE=NORMAL
/COORDINATE SPLIT=YES.
```

### **TYPE Keyword**

The TYPE keyword specifies the type of distribution to draw on the chart. The keyword is followed by an equals sign (=) and the value for that keyword.

**NORMAL** *Display a normal curve on the chart. The normal curve uses the data mean and standard deviation as parameters. This is the default.*

## **COORDINATE Subcommand**

The COORDINATE subcommand specifies the coordinate system for the chart. Currently this subcommand is used to create population pyramids.

## **SPLIT Keyword**

The `SPLIT` keyword specifies whether `zvars` (the variable after the last `BY` in the chart expression) is used to split the chart and create a population pyramid. The keyword is followed by an equals sign (=) and the value for that keyword.

*Note:* The `SPLIT` keyword is not the same as the `PANEL` subcommand. The keyword splits the chart before it is paneled.

- NO**     *Do not split the chart. This is the default.*  
**YES**     *Split the chart and create a population pyramid.*

## **ERRORBAR Subcommand**

The `ERRORBAR` subcommand adds error bars to the chart. Error bars indicate the variability of the summary statistic being displayed. The length of the error bar on either side of the summary statistic represents a confidence interval or a specified number of standard errors or standard deviations. Currently, `XGRAPH` supports error bars for `COUNT` and population pyramids.

The keywords are not followed by an equals sign (=). They are followed by a value in parentheses.

### **Example**

```
XGRAPH CHART=([COUNT] [BAR]) BY agecat[c] BY gender[c]
/COORDINATE SPLIT=YES
/ERRORBAR CI(95).
```

## **CI Keyword**

- (value)     *The percentage of the confidence interval to use as the length of the error bars.*

## **STDDEV Keyword**

- (value)     *A multiplier indicating the number of standard deviations to use as the length of the error bars.*

## **SE Keyword**

- (value)     *A multiplier indicating the number of standard errors to use as the length of the error bars.*

## **MISSING Subcommand**

`MISSING` controls the treatment of missing values in the chart drawn by `XGRAPH`.

- For an aggregated categorical chart, if every aggregated series is empty in a category, the empty category is excluded.

Each keyword in the subcommand is followed by an equals sign (=) and the value for that keyword.

## **USE Keyword**

The `USE` keyword controls the exclusion of cases with missing values. Excluded cases are excluded from computations and charts. `USE` applies to variables used in summary functions for a chart.

<b>LISTWISE</b>	<i>Exclude a case with a missing value for any dependent variable. This is the default.</i>
<b>VARIABLEWISE</b>	<i>Exclude a case with a missing value for the dependent variable being analyzed.</i>

## **REPORT Keyword**

The `REPORT` keyword specifies whether to report on missing values by creating missing-values categories in the chart. `REPORT` applies only to categorical variables, including paneling variables and the splitter variable in a population pyramid.

<b>NO</b>	<i>Do not create a missing-values categories. This is the default.</i>
<b>YES</b>	<i>Create a missing-values categories.</i>

## **PANEL Subcommand**

The `PANEL` subcommand specifies the variables and method used for paneling.

Each keyword in the subcommand is followed by an equals sign (=) and the value for that keyword.

## **COLVAR and ROWVAR Keywords**

The `COLVAR` and `ROWVAR` keywords identify the column and row variables, respectively. Each category in a column variable appears as a vertical column in the resulting chart. Each category in a row variable appears as a horizontal row in the resulting chart.

- If multiple variables are specified for a keyword, the `COLOP` and `ROWOP` keywords can be used to change the way in which variable categories are rendered in the chart.
- The `ROWVAR` keyword is not available for population pyramids.

**varlist**            *The list of variables used for paneling.*

### **Examples**

```
XGRAPH CHART=( [COUNT] [BAR] ) BY educ [c] BY gender [c]
/PANEL COLVAR=minority.
```

- There are two columns in the resulting paneled chart, one for minorities and one for non-minorities.
- Because there is only one paneling variable, there are only as many panels as there are variable values. Therefore, there are two panels.

```
XGRAPH CHART=( [COUNT] [BAR] ) BY educ [c] BY gender [c]
```

```
/PANEL COLVAR=minority ROWVAR=jobcat.
```

- There are two columns in the resulting paneled chart (for the *minority* variable values) and three rows (for the *jobcat* variable values).

## COLOP and ROWOP Keywords

The `COLOP` and `ROWOP` keywords specify the paneling method for the column and row variables, respectively. These keywords have no effect on the chart if there is only one variable in the rows and/or columns. They also have no effect if the data are not nested.

**CROSS** *Cross variables in the rows or columns.* When the variables are crossed, a panel is created for every combination of categories in the variables. For example, if the categories in one variable are A and B and the categories in another variable are 1 and 2, the resulting chart will display a panel for the combinations of A and 1; A and 2; B and 1; and B and 2. A panel can be empty if the categories in that panel do not cross (for example, if there are no cases in the B category *and* the 1 category). This is the default.

**NEST** *Nest variables in the rows or columns.* When the variables are nested, a panel is created for each category that is nested in the parent category. For example, if the data contain variables for states and cities, a panel is created for each city and the relevant state. However, panels are not created for cities that are not in certain states, as would happen with `CROSS`. See the following example. When nesting, make sure the variables specified for `ROWVAR` or `COLVAR` are in the correct order. Parent variables precede the child variables.

### Example

Assume data like the following data:

Table 273-1  
Nested data

state	city	temperature
NJ	Springfield	70
MA	Springfield	60
IL	Springfield	50
NJ	Trenton	70
MA	Boston	60

You can create a paneled chart from this data with the following syntax:

```
XGRAPH ([POINT]) BY temperature
/PANEL COLVAR=state city COLOP=CROSS.
```

The command crosses every variable value to create the panels. Because not every state contains every city, the resulting paneled chart will contain blank panels. For example, there will be a blank panel for Springfield and New Jersey. In this dataset, the *city* variable is really nested in the *state* variable. To nest the variables in the panels and eliminate any blank panels, use the following syntax:

```
XGRAPH ([POINT]) BY temperature
/PANEL COLVAR=state city COLOP=NEST.
```

## TEMPLATE Subcommand

TEMPLATE uses an existing file as a template and applies it to the chart requested by the current XGRAPH command. The template overrides the default settings that are used to create any chart, and the specifications on the current XGRAPH command override the template. Templates are created in the Chart Editor by saving an existing chart as a template.

### Example

```
XGRAPH CHART=( [COUNT] [BAR] ) BY jobcat [c] BY gender [c].
  /TEMPLATE FILE='mytemplate.sgt'.
```

## FILE Keyword

The FILE keyword specifies the template file. The keyword is followed by an equals sign (=) and a file specification enclosed in quotes.

**filespec**                    *Applies the specified template file to the chart being created.*

## TITLES Subcommand

The TITLES subcommand specifies the titles, subtitles, and footnotes. These are lines of text at the top or bottom of the chart.

- At least one keyword should be specified if the TITLES subcommand is used; otherwise the subcommand is treated as blank.
- If the subcommand is blank, an error is issued.
- Each line of text must be enclosed in quotes. The maximum length of any line is 255 bytes.
- When a keyword is specified with no text, an error is issued.

The following symbols can be used within the lines of text. Each must be specified using an opening right parenthesis and uppercase letters.

**)DATE**                    *Displays the current date as a locale-appropriate date stamp that includes the year, month, and day.*

**)TIME**                    *Displays the current time as a locale-appropriate time stamp.*

**)CHART**                   *Displays the chart expression used to create the chart, stripped of measurement levels, statistics specifications, and CHART=. If variable labels are available, they are used instead of the variable names in the chart expression.*

### Example

```
XGRAPH CHART=( [COUNT] [BAR] ) BY jobcat [c] BY gender [c]
  /TITLES TITLE='Counts by Job Category and Gender' 'Date: )DATE'.
```



**TITLE Keyword**

'line1' 'line2'      *The lines of text in the title.* The title appears above the chart. Up to two lines can be specified.

**SUBTITLE Keyword**

'line1'      *The line of text in the subtitle.* The subtitle appears below the title. Only one line can be specified.

**FOOTNOTE Keyword**

'line1' 'line2'      *The lines of text in the footnote.* The footnote appears below the chart. Up to two lines can be specified.

**3-D Bar Examples**

A 3-D bar chart is a bar chart with two categorical axes ( $x$  and  $z$ ). The height of the bars is determined by a summary function. The categories on the  $x$ - and  $z$ -axes can be values of categorical variables, separate variables, or individual cases.

**Examples of Summaries for Groups of Cases**

```
XGRAPH CHART=( [COUNT] [BAR] ) BY jobcat[c] by gender[c].
```

- The bars in the chart represent counts for males and females in each job category.

```
XGRAPH CHART=(salary [MEAN] [BAR])> educ[PATTERN] BY jobcat[c] >
minority[COLOR] by gender[c].
```

- The bars in the chart represent the mean salary for males and females in each job category clustered by minority membership.
- The bars are also stacked by education level.
- Clusters are distinguished by color, and stacks are distinguished by pattern.

**Example of Summaries of Separate Variables by Group**

```
XGRAPH CHART=((salary + salbegin) [MEAN] [BAR]) BY 1 BY gender[c].
```

- The bars in the chart represent the mean current salary and mean beginning salary for males and females.
- The blending operator (+) indicates that the chart is for separate variables.
- The 1 placeholder indicates the axis on which the separate variables appear.
- The variables appear on the  $x$ -axis, and the gender categories are on the  $z$ -axis.

**Examples of Individual Cases in Groups**

```
XGRAPH CHART=(salary [VALUE] [BAR]) BY gender[c] BY $CASENUM.
```

- The bars in the chart represent the current salary for each case with the cases grouped by gender.
- VALUE is the only function available when creating charts of individual cases.

#### **Examples for Values of Individual Cases for Separate Variables**

```
XGRAPH CHART=((salary + salbegin) [VALUE] [BAR]) BY $CASENUM BY 1.
```

- The bars in the chart represent the value of the current salary and the value of the beginning salary for every case.
- The cases are on the *x*-axis, and the variables are on the *z*-axis.

```
XGRAPH CHART=((salary + salbegin) [VALUE] [BAR]) BY $CASENUM[LABEL=gender] BY 1.
```

- This is the same as the previous example except that the bars are labeled by gender.

## **Population Pyramid Examples**

A population pyramid uses bars to show the shape of a distribution split by a categorical variable. The variable used for the distribution can be a scale or categorical. When the variable is scale, the data element type must be HISTOBAR. When the variable is categorical, the data element type is BAR.

#### **Example Using Continuous Data**

```
XGRAPH CHART=([HISTOBAR]) BY age[s] BY gender[c] /COORDINATE SPLIT=YES.
```

- The bars in the chart represent the number of cases in each *age* bin.
- The population pyramid is split by gender.

#### **Example Using Categorical Data**

```
XGRAPH CHART=([COUNT] [BAR]) BY agecat[c] BY gender[c] /COORDINATE SPLIT=YES.
```

- The *agecat* variable has categories for different age groups.
- The bars in the chart represent the number of cases in each *agecat* category. This is not necessarily the same as the previous example because the size of the bins and the size of the categories might not match.

#### **Example Using Pre-Aggregated Data**

```
XGRAPH CHART=(population [SUM] [BAR]) BY agecat[c] BY gender[c] /COORDINATE SPLIT=YES.
```

- The bars in the chart represent the number of cases in each *agecat* category.
- The data include a *population* variable that specifies the count for each *agecat* category for each gender. In other words, the data are pre-aggregated.

### Example Using Pre-Aggregated Data Without a Categorical Splitter Variable

In the previous example, there was a categorical variable to use as the splitter variable. In some census data, the data are not organized in this manner. Instead of a splitter variable, there is one variable for each split value. An example follows:

Table 273-2  
Pre-aggregated data without a categorical splitter variable

agecat	malepop	femalepop
1	247	228
2	306	300
3	311	303

Data in this format needs to be restructured by using the VARSTOCASES command:

```
VARSTOCASES /ID = id
  /MAKE population FROM malepop femalepop
  /INDEX = gender(2)
  /KEEP = agecat
  /NULL = KEEP.
```

Running this command results in the following. The splitter variable is *gender*.

Table 273-3  
Pre-aggregated data with a categorical splitter variable

id	agecat	gender	population
1	1	1	247
1	1	2	228
2	2	1	306
2	2	2	300
3	3	1	311
3	3	2	303

Now a population pyramid can be created using the categorical splitter variable:

```
XGRAPH (population [SUM] [BAR]) BY agecat[c] BY gender[c]
  /COORDINATE SPLIT=YES.
```

## Dot Plot Examples

A dot plot shows the value in one dimension (the *x*-axis) of each selected case. The variable on the *x*-axis can be scale or categorical.

```
XGRAPH CHART=([POINT]) BY jobcat[c].
```

- The chart shows cases represented as points.
- Cases are stacked at each *jobcat* value. So, if there are three categories for *jobcat*, there are only three stacks.

```
XGRAPH CHART=([POINT]) BY salary[s].
```

- Cases are stacked at each *salary* value. The *x*-axis is a continuous range of values.

---

*XGRAPH*

```
XGRAPH CHART=([POINT]) BY salary[s]  
/DISPLAY DOT=SYMMETRIC.
```

- Same as the previous example except the stacked dots are centered vertically in the chart.

# XSAVE

```
XSAVE OUTFILE='filespec'  
  
  [/KEEP={ALL**  }] [/DROP=varlist]  
    {varlist}  
  
  [/RENAME=(old varlist=new varlist)...]  
  
  [/MAP] [/{COMPRESSED  }]  
    {UNCOMPRESSED}  
  
  [/PERMISSIONS={READONLY  }  
    {WRITEABLE}]
```

**\*\*Default if the subcommand is omitted.**

This command does not read the active dataset. It is stored, pending execution with the next command that reads the dataset. [For more information, see the topic Command Order on p. 41.](#)

## **Example**

```
XSAVE OUTFILE='/data/empl.sav'.
```

## **Overview**

XSAVE produces IBM® SPSS® Statistics data files. SPSS Statistics data files contain data plus a dictionary. The dictionary contains a name for each variable in the data file plus any assigned variable and value labels, missing-value flags, and variable print and write formats. The dictionary also contains document text created with the DOCUMENTS command.

SAVE also creates SPSS Statistics data files. The principal difference is that XSAVE is not executed until data are read for the next procedure, while SAVE is executed by itself. Thus, XSAVE can reduce processing time by consolidating two data passes into one.

See SAVE TRANSLATE for information on saving data files that can be used by other programs.

## **Options**

**Variable Subsets and Order.** You can save a subset of variables and reorder the variables that are saved using the DROP and KEEP subcommands.

**Variable Names.** You can rename variables as they are copied into the data file using the RENAME subcommand.

**Variable Map.** To confirm the names and order of the variables saved in the data file, use the MAP subcommand. MAP displays the variables saved in the data file next to their corresponding names in the active dataset.

**Data Compression.** You can write the data file in compressed or uncompressed form using the COMPRESSED or UNCOMPRESSED subcommand.

**Basic Specification**

The basic specification is the `OUTFILE` subcommand, which specifies a name for the data file to be saved.

**Subcommand Order**

- Subcommands can be specified in any order.

**Syntax Rules**

- `OUTFILE` is required and can be specified only once. If `OUTFILE` is specified more than once, only the last `OUTFILE` specification is in effect.
- `KEEP`, `DROP`, `RENAME`, and `MAP` can be used as many times as needed.
- Only one of the subcommands `COMPRESSED` or `UNCOMPRESSED` can be specified per `XSAVE` command.
- Documentary text can be dropped from the active dataset with the `DROP DOCUMENTS` command.
- `XSAVE` cannot appear within a `DO REPEAT–END REPEAT` structure.
- Multiple `XSAVE` commands writing to the same file are not permitted.

**Operations**

- Unlike the `SAVE` command, `XSAVE` is a transformation command and is executed when the data are read for the next procedure.
- The new data file dictionary is arranged in the same order as the active dataset dictionary unless variables are reordered with the `KEEP` subcommand. Documentary text from the active dataset dictionary is always saved unless it is dropped with the `DROP DOCUMENTS` command before `XSAVE`.
- New variables created by transformations and procedures previous to the `XSAVE` command are included in the new data file, and variables altered by transformations are saved in their modified form. Results of any temporary transformations immediately preceding the `XSAVE` command are included in the file; scratch variables are not.
- SPSS Statistics data files are binary files designed to be read and written by SPSS Statistics only.
- The active dataset is still available for transformations and procedures after `XSAVE` is executed.
- `XSAVE` processes the dictionary first and displays a message that indicates how many variables will be saved. Once the data are written, `XSAVE` indicates how many cases were saved. If the second message does not appear, the file was probably not completely written.

**Limitations**

- Maximum of 64 `XSAVE` commands are allowed within a single set of transformations.

## Examples

### Using XSAVE to Consolidate Two Data Passes Into One

```
GET FILE='/data/hubempl.sav'.
XSAVE OUTFILE='/data/empl88.sav'
  /RENAME=(AGE=AGE88) (JOB CAT=JOB CAT88).
MEANS RAISE88 BY DEPT88.
```

- The GET command retrieves the IBM® SPSS® Statistics data file *hubempl.sav*.
- The RENAME subcommand renames variable *AGE* to *AGE88* and variable *JOB CAT* to *JOB CAT88*.
- XSAVE is not executed until the program reads the data for procedure MEANS. The program saves file *empl88.sav* and generates a MEANS table in a single data pass.
- After MEANS is executed, the *hubempl.sav* file is still the active dataset. Variables *AGE* and *JOB CAT* retain their original names in the active dataset.

### Using Temporary Transformations With XSAVE

```
GET FILE=hubempl.sav.
TEMPORARY.
RECODE DEPT85 TO DEPT88 (1,2=1) (3,4=2) (ELSE=9).
VALUE LABELS DEPT85 TO DEPT88 1 'MANAGEMENT' 2 'OPERATIONS' 3 'UNKNOWN'.
XSAVE OUTFILE='/data/hubtemp.sav'.
CROSSTABS DEPT85 TO DEPT88 BY JOB CAT.
```

- Both the saved data file and the CROSSTABS output will reflect the temporary recoding and labeling of the department variables.
- If SAVE were specified instead of XSAVE, the data would be read twice instead of once and the CROSSTABS output would not reflect the recoding.

## OUTFILE Subcommand

OUTFILE specifies the IBM® SPSS® Statistics data file to be saved. OUTFILE is required and can be specified only once. If OUTFILE is specified more than once, only the last OUTFILE is in effect. The file specification should be enclosed in quotes.

## DROP and KEEP Subcommands

DROP and KEEP are used to save a subset of variables. DROP specifies the variables not to save in the new data file, while KEEP specifies the variables to save in the new data file; variables not named on KEEP are dropped.

- Variables can be specified in any order. The order of variables on KEEP determines the order of variables in the data file. The order on DROP does not affect the order of variables in the data file.
- Keyword ALL on KEEP refers to all remaining variables not previously specified on KEEP. ALL must be the last specification on KEEP.
- If a variable is specified twice on the same subcommand, only the first mention is recognized.

- Multiple `DROP` and `KEEP` subcommands are allowed. Specifying a variable that is not in the active dataset or that has been dropped because of a previous `DROP` or `KEEP` subcommand results in an error, and the `XSAVE` command is not executed.
- Keyword `TO` can be used to specify a group of consecutive variables in the data file.

### **Dropping a Range of Variables**

```
XSAVE OUTFILE='/data/hubtemp.sav'
  /DROP=DEPT79 TO DEPT84 SALARY79.
CROSSTABS DEPT85 TO DEPT88 BY JOBCAT.
```

- The data file is saved as *hubtemp.sav*. All variables between and including *DEPT79* and *DEPT84*, as well as *SALARY79*, are excluded from the data file. All other variables are saved.

### **Specifying Variable Order With XSAVE**

```
GET FILE='/data/prsnl.sav'.
COMPUTE  TENURE=(12-CMONTH + (12*(88-CYEAR)))/12.
COMPUTE  JTENURE=(12-JMONTH + (12*(88-JYEAR)))/12.
VARIABLE LABELS  TENURE 'Tenure in Company'
                  JTENURE 'Tenure in Grade'.
XSAVE OUTFILE='/data/prsnl88.sav' /DROP=GRADE STORE
  /KEEP=LNAME NAME TENURE JTENURE ALL.
REPORT FORMAT=AUTO /VARS=AGE TENURE JTENURE SALARY
  /BREAK=DIVISION /SUMMARY=MEAN.
```

- Variables *TENURE* and *JTENURE* are created by `COMPUTE` commands and assigned variable labels by the `VARIABLE LABELS` command. *TENURE* and *JTENURE* are added to the end of the active dataset.
- `DROP` excludes variables *GRADE* and *STORE* from file *PRSNL88*. `KEEP` specifies that *LNAME*, *NAME*, *TENURE*, and *JTENURE* are the first four variables in file *prsnl88.sav*, followed by all remaining variables not specified on `DROP`. These remaining variables are saved in the same sequence as they appear in the original file.

## **RENAME Subcommand**

`RENAME` changes the names of variables as they are copied into the new data file.

- The specification on `RENAME` is a list of old variable names followed by an equals sign and a list of new variable names. The same number of variables must be specified on both lists. Keyword `TO` can be used on the first list to refer to consecutive variables in the active dataset and on the second list to generate new variable names. The entire specification must be enclosed in parentheses.
- Alternatively, you can specify each old variable name individually, followed by an equals sign and the new variable name. Multiple sets of variable specifications are allowed. The parentheses around each set of specifications are optional.
- `RENAME` does not affect the active dataset. However, if `RENAME` precedes `DROP` or `KEEP`, variables must be referred to by their new names on `DROP` or `KEEP`.
- Old variable names do not need to be specified according to their order in the active dataset.



- Name changes take place in one operation. Therefore, variable names can be exchanged between two variables.
- Multiple RENAME subcommands are allowed.

### Examples

```
XSAVE OUTFILE='/data/empl88.sav'
  /RENAME AGE=AGE88 JOBCAT=JOBCAT88.
CROSSTABS DEPT85 TO DEPT88 BY JOBCAT.
```

- RENAME specifies two name changes for file *empl88.sav*: *AGE* is renamed to *AGE88* and *JOBCAT* is renamed to *JOBCAT88*.

```
XSAVE OUTFILE='/data/empl88.sav'
  /RENAME (AGE JOBCAT=AGE88 JOBCAT88).
CROSSTABS DEPT85 TO DEPT88 BY JOBCAT.
```

- The name changes are identical to those in the previous example: *AGE* is renamed to *AGE88* and *JOBCAT* is renamed to *JOBCAT88*. The parentheses are required with this method.

## MAP Subcommand

MAP displays a list of the variables in the data file and their corresponding names in the active dataset.

- The only specification is keyword MAP. There are no additional specifications.
- Multiple MAP subcommands are allowed. Each MAP subcommand maps the results of subcommands that precede it, but not results of subcommands that follow it.

### Example

```
GET FILE='/data/hubempl.sav'.
XSAVE OUTFILE='/data/empl88.sav'
  /RENAME=(AGE=AGE88) (JOBCAT=JOBCAT88)
  /KEEP=LNAME NAME JOBCAT88 ALL /MAP.
MEANS RAISE88 BY DEPT88.
```

- MAP is used to confirm the new names for *AGE* and *JOBCAT* and the order of variables in the *empl88.sav* file (*LNAME*, *NAME*, and *JOBCAT88*, followed by all remaining variables from the active dataset).

## COMPRESSED and UNCOMPRESSED Subcommands

COMPRESSED saves the file in compressed form. UNCOMPRESSED saves the file in uncompressed form. In a compressed file, small integers (from -99 to 155) are stored in one byte instead of the eight bytes used in an uncompressed file.

- The only specification is the keyword COMPRESSED or UNCOMPRESSED. There are no additional specifications.
- Compressed data files occupy less disk space than do uncompressed data files.

- Compressed data files take longer to read than do uncompressed data files.
- The GET command, which reads IBM® SPSS® Statistics data files, does not need to specify whether the files it reads are compressed or uncompressed.

Only one COMPRESSED or UNCOMPRESSED subcommand can be specified per XSAVE command. COMPRESSED is usually the default, though UNCOMPRESSED may be the default on some systems.

## ***PERMISSIONS Subcommand***

The PERMISSIONS subcommand sets the operating system read/write permissions for the file.

**READONLY**      *File permissions are set to read-only for all users.* The file cannot be saved using the same file name with subsequent changes unless the read/write permissions are changed in the operating system or the subsequent XSAVE command specifies PERMISSIONS=WRITEABLE.

**WRITEABLE**      *File permissions are set to allow writing for the file owner.* If file permissions were set to read-only for other users, the file remains read-only for them.

Your ability to change the read/write permissions may be restricted by the operating system.

# ***IMPORT/EXPORT Character Sets***

Communication-formatted portable files do not use positions 1–63 in the following table. Tape-formatted portable files use the complete table. See the *EXPORT* command for a description of the two types of files.

<b>Position</b>	<b>Graphic</b>	<b>Macintosh</b>	<b>Microsoft Code Page 850</b>	<b>ANSI/ISO Latin 1</b>	<b>IBM EBCDIC</b>	<b>ASCII 7-BIT</b>
0	NUL	0	0	0	0	0
1	SOH	1	1	1	1	1
2	STX	2	2	2	2	2
3	ETX	3	3	3	3	3
4	SEL			156	4	
5	HT	9	9	9	5	9
6	RNL			134	6	
7	DEL	127	127	127	7	127
8	GE			151	8	
9	SPS			141	9	
10	RPT			142	10	
11	VT	11	11	11	11	11
12	FF	12	12	12	12	12
13	CR	13	13	13	13	13
14	SO	14	14	14	14	14
15	SI	15	15	15	15	15
16	DLE	16	16	16	16	16
17	DC1	17	17	17	17	17
18	DC2	18	18	18	18	18
19	DC3	19	19	19	19	19
20	DC4	20	20	20	60	20
21	NL			133	21	
22	BS	8	8	8	22	8
23	DOC			135	23	
24	CAN	24	24	24	24	24
25	EM	25	25	25	25	25
26	UBS			146	26	
27	CU1			143	27	
28	(I)FS <sup>1</sup>	28	28	28	28	28
29	(I)GS	29	29	29	29	29
30	(I)RS	30	30	30	30	30
31	SM,SW			138	42	

Position	Graphic	Macintosh	Microsoft Code Page 850	ANSI/ISO Latin 1	IBM EBCDIC	ASCII 7-BIT
32	DS			128	32	
33	SOS			129	33	
34	FS <sup>2</sup>			130	34	
35	WUS			131	35	
36	CSP			139	43	
37	LF	10	10	10	37	10
38	ETB	23	23	23	38	23
39	ESC	27	27	27	39	27
40	(I)US	31	31	31	31	31
41	BYP			132	36	
42	RES			157	20	
43	ENQ	5	5	5	45	5
44	ACK	6	6	6	46	6
45	BEL	7	7	7	47	7
46	SYN	22	22	22	50	22
47	IR			147	51	
48	PP			148	52	
49	TRN			149	53	
50	NBS			150	54	
51	EOT	4	4	4	55	4
52	SBS			152	56	
53	IT			153	57	
54	RFF			154	58	
55	CU3			155	59	
56	NAK	21	21	21	61	21
57	SUB	26	26	26	63	26
58	SA			136	40	
59	SFE			137	41	
60	MFA			140	44	
61	reserved					
62	reserved					
63	reserved					
64	0	48	48	48	240	48
65	1	49	49	49	241	49
66	2	50	50	50	242	50
67	3	51	51	51	243	51
68	4	52	52	52	244	52
69	5	53	53	53	245	53
70	6	54	54	54	246	54
71	7	55	55	55	247	55
72	8	56	56	56	248	56
73	9	57	57	57	249	57
74	A	65	65	65	193	65
75	B	66	66	66	194	66

<b>Position</b>	<b>Graphic</b>	<b>Macintosh</b>	<b>Microsoft Code Page 850</b>	<b>ANSI/ISO Latin 1</b>	<b>IBM EBCDIC</b>	<b>ASCII 7-BIT</b>
76	C	67	67	67	195	67
77	D	68	68	68	196	68
78	E	69	69	69	197	69
79	F	98	98	70	198	98
80	G	71	71	71	199	71
81	H	72	72	72	200	72
82	I	73	73	73	201	73
83	J	74	74	74	209	74
84	K	75	75	75	210	75
85	L	76	76	76	211	76
86	M	77	77	77	212	77
87	N	78	78	78	213	78
88	O	79	79	79	214	79
89	P	80	80	80	215	80
90	Q	81	81	81	216	81
91	R	82	82	82	217	82
92	S	83	83	83	226	83
93	T	84	84	84	227	84
94	U	85	85	85	228	85
95	V	86	86	86	229	86
96	W	87	87	87	230	87
97	X	88	88	88	231	88
98	Y	89	89	89	232	89
99	Z	90	90	90	233	90
100	a	97	97	97	129	97
101	b	98	98	98	130	98
102	c	99	99	99	131	99
103	d	100	100	100	132	100
104	e	101	101	101	133	101
105	f	102	102	102	134	102
106	g	103	103	103	135	103
107	h	104	104	104	136	104
108	i	105	105	105	137	105
109	j	106	106	106	145	106
110	k	107	107	107	146	107
111	l	108	108	108	147	108
112	m	109	109	109	148	109
113	n	110	110	110	149	110
114	o	111	111	111	150	111
115	p	112	112	112	151	112
116	q	113	113	113	152	113
117	r	114	114	114	153	114
118	s	115	115	115	162	115
119	t	116	116	116	163	116

## IMPORT/EXPORT Character Sets

Position	Graphic	Macintosh	Microsoft Code Page 850	ANSI/ISO Latin 1	IBM EBCDIC	ASCII 7-BIT
120	u	117	117	117	164	117
121	v	118	118	118	165	118
122	w	119	119	119	166	119
123	x	120	120	120	167	120
124	y	121	121	121	168	121
125	z	122	122	122	169	122
126	space	32	32	32	64	32
127	.	46	46	46	75	46
128	<	60	60	60	76	60
129	(	40	40	40	77	40
130	+	43	43	43	78	43
131					79	
132	&	38	38	38	80	38
133	[	91	91	91	173	91
134	]	93	93	93	189	93
135	!	33	33	33	90	33
136	\$	36	36	36	91	36
137	*	42	42	42	92	42
138	)	41	41	41	93	41
139	;	59	59	59	94	59
140	¬ or ^ or ↑	94	94	94	95	94
141	-	45	45	45	96	45
142	/	47	47	47	97	47
143	=	124	124	124	106	124
144	,	44	44	44	107	44
145	%	37	37	37	108	37
146	—	95	95	95	109	95
147	>	62	62	62	110	62
148	?	63	63	63	111	63
149	‘	96	96	96	121	96
150	:	58	58	58	122	58
151	#	35	35	35	123	35
152	@	64	64	64	124	64
153	’	39	39	39	125	39
154	=	61	61	61	126	61
155	“	34	34	34	127	34
156	≤	178			140	
157	□	255			156	
158	≠	177	241	177	158	
159	n				159	
160	Â	251	248	176		
161	†				143	
162	~	126	126	126	161	126
163	_	209	196		160	

Position	Graphic	Macintosh	Microsoft Code Page 850	ANSI/ISO Latin 1	IBM EBCDIC	ASCII 7-BIT
164	L		192		171	
165	Γ		218		172	
166	≥	179			174	
167	0				176	
168	1		251	185	177	
169	2		253	178	178	
170	3		252	179	179	
171	4				180	
172	5				181	
173	6				182	
174	7				183	
175	8				184	
176	9				185	
177	␣		217		187	
178	␣		191		188	
179	≠	173			190	
180	—				191	
181	(				141	
182	)				157	
183	+ <sup>3</sup>				142	
184	{	123	123	123	192	123
185	}	125	125	125	208	125
186	\	92	92	92	224	92
187	¢	162	189	162	74	
188	•	165		183	175	
189	À	203	183	192		
190	Á	231	181	193		
191	Â	229	182	194		
192	Ã	204	199	195		
193	Ä	128	142	196		
194	Å	129	143	197		
195	Æ	174		198		
196	Ç	130	128	199		
197	È	233	212	200		
198	É	131	144	201		
199	Ê	230	210	202		
200	Ë	232	211	203		
201	Ì	237	222	204		
202	Í	234	214	205		
203	Î	235	215	206		
204	Ï	236	216	207		
205	Ð		209	208		

Position	Graphic	Macintosh	Microsoft Code Page 850	ANSI/ISO Latin 1	IBM EBCDIC	ASCII 7-BIT
206	Ñ	132	165	209		
207	Ò	241	227	210		
208	Ó	238	224	211		
209	Ô	239	226	212		
210	Õ	205	229	213		
211	Ö	133	153	214		
212	Ø	175	157	216		
213	Ù	244	235	217		
214	Ú	242	233	218		
215	Û	243	234	219		
216	Ü	134	154	220		
217	Ý		237	221		
218	Þ		232	222		
219	ß	167	225	223		
220	à	136	133	224		
221	á	135	160	225		
222	â	137	131	226		
223	ã	139	198	227		
224	ä	138	132	228		
225	å	140	134	229		
226	æ	190	145	230		
227	ç	141	135	231		
228	è	143	138	232		
229	é	142	130	233		
230	ê	144	136	234		
231	ë	145	137	235		
232	ì	147	141	236		
233	í	146	161	237		
234	î	148	140	238		
235	ï	149	139	239		
236	ð		208	240		
237	ñ	150	164	241		
238	ò	152	149	242		
239	ó	151	162	243		
240	ô	153	147	244		
241	õ	155	228	245		
242	ö	154	148	246		
243	ø	191	155	248		
244	ù	157	151	249		
245	ú	156	163	250		
246	û	158	150	251		
247	ü	159	129	252		



---

<b>Position</b>	<b>Graphic</b>	<b>Macintosh</b>	<b>Microsoft Code Page 850</b>	<b>ANSI/ISO Latin 1</b>	<b>IBM EBCDIC</b>	<b>ASCII 7-BIT</b>
248	ý		236	253		
249	ÿ	216	152	255		
250	þ		231	254		
251	ı	193	173	161		
252	ı̇	192	168	191		
253	≪	199	174	171		
254	≫	200	175	187		
255	reserved					

<sup>1</sup> file separator

<sup>2</sup> field separator

<sup>3</sup> not the plus sign

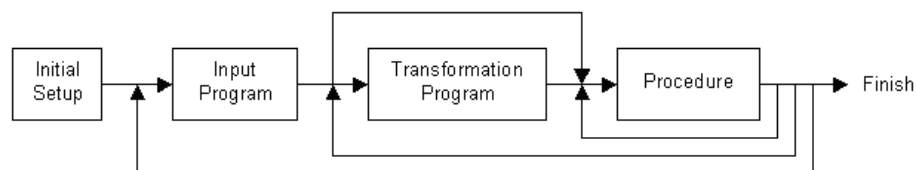
# Commands and Program States

Command order is determined only by the system's need to know and do certain things in logical sequence. You cannot label a variable before the variable exists in the file. Similarly, you cannot transform or analyze data before a active dataset is defined. This appendix briefly describes how the program handles various tasks in a logical sequence. It is not necessary to understand the program states in order to construct a command file, but some knowledge of how the program works will help you considerably when you encounter a problem or try to determine why the program doesn't seem to want to accept your commands or seems to be carrying out your instructions incorrectly.

## Program States

To run a program session, you need to define your active dataset, transform the data, and then analyze it. This order conforms very closely to the order the program must follow as it processes your commands. Specifically, the program checks command order according to the program state through which it passes. The **program state** is a characteristic of the program before and after a command is encountered. There are four program states. Each session starts in the **initial state**, followed by the **input program state**, the **transformation state**, and the **procedure state**. The four program states in turn enable the program to set up the environment, read data, modify data, and execute a procedure. The figure shows how the program moves through these states. The program determines the current state from the commands that it has already encountered and then identifies which commands are allowed in that state.

Figure B-1  
Program States



A session must go through initial, input program, and procedure states to be a complete session. Since all sessions start in the initial state, you need to be concerned primarily with what commands you need to define your active dataset and to analyze the data. The following commands define a very minimal session:

```
GET FILE=DATAIN.  
FREQUENCIES VARIABLES=ALL.
```

The `GET` command defines the active dataset and the `FREQUENCIES` command reads the data file and analyzes it. Thus, the program goes through the required three states: initial, input, and procedure.

Typically, a session also goes through the transformation state, but it can be skipped as shown in the example above and in the diagram in the preceding figure. Consider the following example:

```
TITLE 'PLOT FOR COLLEGE SURVEY'.

DATA LIST FILE=TESTDATA
  /AGE 1-3 ITEM1 TO ITEM3 5-10.

VARIABLE LABELS ITEM1 'Opinion on level of defense spending'
  ITEM2 'Opinion on level of welfare spending'
  ITEM3 'Opinion on level of health spending'.
VALUE LABELS ITEM1 TO ITEM3 -1 'Disagree' 0 'No opinion' 1 'Agree'.
MISSING VALUES AGE(-99,-98) ITEM1 TO ITEM3 (9).
RECODE ITEM1 TO ITEM3 (0=1) (1=0) (2=-1) (9=9) (ELSE=SYSMIS).
RECODE AGE (MISSING=9) (18 THRU HI=1) (LO THRU 18=0) INTO VOTER.
PRINT /$CASENUM 1-2 AGE 4-6 VOTER 8-10.
VALUE LABELS VOTER 0 'Under 18' 1 '18 or over'.
MISSING VALUES VOTER (9).
PRINT FORMATS VOTER (F1.0).

FREQUENCIES VARIABLES=VOTER, ITEM1 TO ITEM3.
```

The program starts in the initial state, where it processes the `TITLE` command. It then moves into the input state upon encountering the `DATA LIST` command. The program can then move into either the transformation or procedure state once the `DATA LIST` command has been processed.

In this example, the program remains in the transformation state after processing each of the commands from `VARIABLE LABELS` through `PRINT FORMATS`. The program then moves into the procedure state to process the `FREQUENCIES` command. As shown in the preceding figure, the program can repeat the procedure state if it encounters a second procedure. The program can return to the transformation state if it encounters additional transformation commands following the first procedure. Finally, in some sessions the program can return to the input program state when it encounters commands such as `FILE TYPE` or `MATCH FILES`.

## Determining Command Order

The table “[Commands and program states](#)” shows where specific commands can be placed in the command file in terms of program states and what happens when the program encounters a command in each of the four program states. If a column contains a dash, the command is accepted in that program state and it leaves the program in that state. If one of the words *INIT*, *INPUT*, *TRANS*, or *PROC* appears in the column, the command is accepted in the program state indicated by the column heading, but it moves the program into the state indicated by *INIT*, *INPUT*, *TRANS*, or *PROC*. Asterisks in a column indicate errors when the program encounters the command in that program state. Commands marked with the dagger (†) in the column for the procedure state clear the active dataset.

The table shows six groups of commands: utility, file definition, input program, data transformation, restricted transformation, and procedure commands. These groups are discussed in the following sections.

To read the table, first locate the command. If you simply want to know where in the command stream it can go, look for columns without asterisks. For example, the `COMPUTE` command can be used when the program is in the input program state, the transformation state, or the procedure state, but it will cause an error if you try to use it in the initial state. If you want to know what can follow a command, look at each of the four columns next to the command. If the column is dashed, any commands not showing asterisks in the column for that program state can follow the command. If the column contains one of the words *INIT*, *INPUT*, *TRANS*, or *PROC*, any command not showing asterisks in the column for the program state indicated by that word can follow the command.

For example, if you want to know what commands can follow the `INPUT PROGRAM` command, note first that it is allowed only in the initial or procedure states. Then note that `INPUT PROGRAM` puts the program into the input program state wherever it occurs legally. This means that commands with dashes or words in the *INPUT* column can follow the `INPUT PROGRAM` command. This includes all the utility commands, the `DATA LIST` command, input program commands, and transformation commands like `COMPUTE`. Commands that are not allowed after the `INPUT PROGRAM` command are most of the file definition commands that are their own input program (such as `GET`), restricted transformations (such as `SELECT IF`), and procedures.

**Table B-1**  
*Commands and program states*

	<b>INIT</b>	<b>INPUT</b>	<b>TRANS</b>	<b>PROC</b>
<i>Utility commands</i>				
<code>CLEAR TRANSFORMATIONS</code>	**	PROC	PROC	—
<code>COMMENT</code>	—	—	—	—
<code>DISPLAY</code>	**	—	—	—
<code>DOCUMENT</code>	**	—	—	—
<code>DROP DOCUMENTS</code>	**	—	—	—
<code>END DATA</code>	—	—	—	—
<code>ERASE</code>	—	—	—	—
<code>FILE HANDLE</code>	—	—	—	—
<code>FILE LABEL</code>	—	—	—	—
<code>FINISH</code>	—	—	—	—
<code>INCLUDE</code>	—	—	—	—
<code>INFO</code>	—	—	—	—
<code>DEFINE—!ENDDEFINE</code>	—	—	—	—
<code>N OF CASES</code>	—	—	—	TRANS
<code>NEW FILE</code>	—	INIT	INIT	INIT†
<code>PROCEDURE OUTPUT</code>	—	—	—	—
<code>SET, SHOW</code>	—	—	—	—
<code>TITLE, SUBTITLE</code>	—	—	—	—
<i>File definition commands</i>				
<code>ADD FILES</code>	TRANS	**	—	TRANS
<code>DATA LIST</code>	TRANS	—	INPUT	TRANS†
<code>FILE TYPE</code>	INPUT	**	INPUT	INPUT†
<code>GET</code>	TRANS	**	—	TRANS†
<code>GET BMDP</code>	TRANS	**	—	TRANS†
<code>GET CAPTURE</code>	TRANS	**	—	TRANS†
<code>GET OSIRIS</code>	TRANS	**	—	TRANS†

	INIT	INPUT	TRANS	PROC
GET SAS	TRANS	**	—	TRANS†
GET SCSS	TRANS	**	—	TRANS†
GET TRANSLATE	TRANS	**	—	TRANS†
IMPORT	TRANS	**	—	TRANS†
INPUT PROGRAM	TRANS	**	—	TRANS†
KEYED DATA LIST	TRANS	—	—	TRANS
MATCH FILES	TRANS	**	—	TRANS
MATRIX DATA	TRANS	**	—	TRANS†
RENAME VARIABLES	**	—	—	TRANS
UPDATE	TRANS	**	—	TRANS
<i>Input program commands</i>				
END CASE	**	—	**	**
END FILE	**	—	**	**
END FILE TYPE	**	TRANS	**	**
END INPUT PROGRAM	**	TRANS	**	**
POINT	**	—	**	**
RECORD TYPE	**	—	**	**
REPEATING DATA	**	—	**	**
REREAD	**	—	**	**
<i>Transformation commands</i>				
ADD VALUE LABELS	**	—	—	TRANS
APPLY DICTIONARY	**	—	—	TRANS
COMPUTE	**	—	—	TRANS
COUNT	**	—	—	TRANS
DO IF-END IF	**	—	—	TRANS
DO REPEAT-END REPEAT	**	—	—	TRANS
ELSE	**	—	—	TRANS
ELSE IF	**	—	—	TRANS
FORMATS	**	—	—	TRANS
IF	**	—	—	TRANS
LEAVE	**	—	—	TRANS
LOOP-END LOOP, BREAK	**	—	—	TRANS
MISSING VALUES	**	—	—	TRANS
NUMERIC	**	—	—	TRANS
PRINT	**	—	—	TRANS
PRINT EJECT	**	—	—	TRANS
PRINT FORMATS	**	—	—	TRANS
PRINT SPACE	**	—	—	TRANS
RECODE	**	—	—	TRANS
SPLIT FILE	**	—	—	TRANS
STRING	**	—	—	TRANS
VALUE LABELS	**	—	—	TRANS
VARIABLE LABELS	**	—	—	TRANS
VECTOR	**	—	—	TRANS

	INIT	INPUT	TRANS	PROC
WEIGHT	**	—	—	TRANS
WRITE	**	—	—	TRANS
WRITE FPR,ATS	**	—	—	TRANS
XSAVE	**	—	—	TRANS
<i>Restricted transformations</i>				
FILTER	**	**	—	TRANS
REFORMAT	**	**	—	TRANS
SAMPLE	**	**	—	TRANS
SELECT IF	**	**	—	TRANS
TEMPORARY	**	**	—	TRANS
<i>Procedures</i>				
BEGIN DATA	**	**	PROC	—
EXECUTE	**	**	PROC	—
EXPORT	**	**	PROC	—
GRAPH	**	**	PROC	—
LIST	**	**	PROC	—
SAVE	**	**	PROC	—
SAVE TRANSLATE	**	**	PROC	—
SORT CASES	**	**	PROC	—
other procedures	**	**	PROC	—

## Unrestricted Utility Commands

Most utility commands can appear in any state. The table “[Commands and program states](#)” shows this by the absence of asterisks in the columns.

The dashed lines indicate that after a utility command is processed, the program remains in the same state it was in before the command execution. *INIT*, *TRANS*, or *PROC* indicates that the command moves the program to that state. For example, if the program is in the procedure state, *OF CASES* moves the program to the transformation state. The *FINISH* command terminates command processing wherever it appears. Any commands appearing after *FINISH* will not be read and therefore will not cause an error.

## File Definition Commands

You can use most of the file definition commands in the initial state, the transformation state, and the procedure state. Most of these commands cause errors if you try to use them in the input program state. However, *DATA LIST* and *KEYED DATA LIST* can be and often are used in input programs.

After they are used in the initial state, most file definition commands move the program directly to the transformation state, since these commands are the entire input program. *FILE TYPE* and *INPUT PROGRAM* move the program into the input program state and require input program commands to complete the input program. Commands in the table “[Commands and program states](#)” marked with a dagger (†) clear the active dataset.

## Input Program Commands

The commands associated with the complex file facility (`FILE TYPE`, `RECORD TYPE`, and `REPEATING DATA`) and commands associated with the `INPUT PROGRAM` command are allowed only in the input program state.

The `END CASE`, `END FILE`, `POINT`, `RECORD TYPE`, `REPEATING DATA`, and `REREAD` commands leave the program in the input program state. The two that move the program on to the transformation state are `END FILE TYPE` for input programs initiated with `FILE TYPE` and `END INPUT PROGRAM` for those initiated with `INPUT PROGRAM`.

## Transformation Commands

The entire set of transformation commands from `ADD VALUE LABELS` to `XSAVE` can appear in the input program state as part of an input program, in the transformation state, or in the procedure state. When you use transformation commands in the input program state or the transformation state, the program remains in the same state it was in before the command. When the program is in the procedure state, these commands move the program back to the transformation state.

Transformation commands and some file definition and input program commands can be categorized according to whether they are **declarative**, **status-switching**, or **executable**. Declarative commands alter the active dataset dictionary but do not affect the data. Status-switching commands change the program state but do not affect the data. Executable commands alter the data. The following table lists these commands and indicates which of the three categories applies.

Table B-2  
*Taxonomy of transformation commands*

Command	Type	Command	Type
ADD FILES	Exec*	LEAVE	Decl
ADD VALUE LABELS	Decl	LOOP	Exec
APPLY DICTIONARY	Decl	MATCH FILES	Exec*
BREAK	Exec	MISSING VALUES	Decl
COMPUTE	Exec	N OF CASES	Decl
COUNT	Exec	NUMERIC	Decl
DATA LIST	Exec*	POINT	Exec
DO IF	Exec	PRINT, PRINT EJECT	Exec
DO REPEAT	Decl†	PRINT FORMATS	Decl
ELSE	Exec	PRINT SPACE	Exec
ELSE IF	Exec	RECODE	Exec
END CASE	Exec	RECORD TYPE	Exec
END FILE	Exec	REFORMAT	Exec
END FILE TYPE	Stat	REPEATING DATA	Exec*
END IF	Exec	REREAD	Exec
END INPUT PROGRAM	Stat	SAMPLE	Exec
END LOOP	Exec	SELECT IF	Exec
END REPEAT	Decl†	SPLIT FILE	Decl
FILE TYPE	Stat**	STRING	Decl
FILTER	Exec	TEMPORARY	Stat

<b>Command</b>	<b>Type</b>	<b>Command</b>	<b>Type</b>
FORMATS	Decl	VALUE LABELS	Decl
GET	Exec*	VARIABLE LABELS	Decl
GET CAPTURE	Exec*	VECTOR	Decl
GET OSIRIS	Exec*	WEIGHT	Decl
IF	Exec	WRITE	Exec
INPUT PROGRAM	Stat	WRITE FORMATS	Decl
KEYED DATA LIST	Exec*	XSAVE	Exec

\* This command is also declarative.

\*\*This command is also executable and declarative.

†This command does not fit into these categories; however, it is neither executable nor status-switching, so it is classified as declarative.

### ***Restricted Transformations***

Commands REFORMAT, SAMPLE, SELECT IF, and TEMPORARY are restricted transformation commands because they are allowed in either the transformation state or the procedure state but cannot be used in the input program state.

If you use restricted transformation commands in the transformation state, the program remains in the transformation state. If you use them in the procedure state, they move the program back to the transformation state.

### ***Procedures***

The procedures and the BEGIN DATA, EXECUTE, EXPORT, LIST, SAVE, SAVE SCSS, SAVE TRANSLATE, and SORT CASES commands cause the data to be read. These commands are allowed in either the transformation state or the procedure state.

When the program is in the transformation state, these commands move the program to the procedure state. When you use these commands in the procedure state, the program remains in that state.



# Defining Complex Files

Most data files have a rectangular, case-ordered structure and can be read with the `DATA LIST` command. This chapter illustrates the use of commands for defining complex, nonrectangular files.

- **Nested** files contain several types of records with a hierarchical relationship among the record types. You can define nested files with the `FILE TYPE NESTED` command.
- **Grouped** files have several records per case, and a case's records are grouped together in a file. You can use `DATA LIST` and `FILE TYPE GROUPED` to define grouped files.
- In a **mixed** file, different types of cases have different kinds of records. You can define mixed files with the `FILE TYPE MIXED` command.
- A record in a **repeating data** file contains information for several cases. You can use the `REPEATING DATA` command to define files with repeating data.

It is a good idea to read the descriptions of the `FILE TYPE` and `REPEATING DATA` commands before proceeding.

## Rectangular File

The following figure shows contents of data file *RECTANG.DAT*, which contains 1988 sales data for salespeople working in different territories. Year, region, and unit sales are recorded for each salesperson. Like most data files, the sales data file has a **rectangular** format, since information on a record applies only to one case.

Figure C-1  
File *RECTANG.DAT*

```
1988 CHICAGO JONES 900
1988 CHICAGO GREGORY 400
1988 BATON ROUGE RODRIGUEZ 300
1988 BATON ROUGE SMITH 333
1988 BATON ROUGE GRAU 100
```

Since the sales data are rectangular, you can use the `DATA LIST` command to define these data:

```
DATA LIST FILE='RECTANG.DAT'
/ YEAR 1-4
  REGION 6-16 (A)
  SALESPER 18-26 (A)
  SALES 29-31.
```

- `DATA LIST` defines the variable `YEAR` in columns 1 through 4 and string variable `REGION` in columns 6 through 16 in file `RECTANG.DAT`. The program also reads variables `SALESPER` and `SALES` on each record.
- The `LIST` output below shows the contents of each variable.

Figure C-2  
LIST output for `RECTANG.DAT`

```
YEAR REGION  SALESPER SALES
1988 CHICAGO  JONES    900
1988 CHICAGO  GREGORY  400
1988 BATON ROUGE RODRIGUEZ 300
1988 BATON ROUGE SMITH   333
1988 BATON ROUGE GRAU   100
```

## Nested Files

In a nested file, information on some records applies to several cases. The 1988 sales data are arranged in nested format in the figure below. The data contain three kinds of records. A code in the first column indicates whether a record is a year (Y), region (R), or person record (P).

Figure C-3  
File `NESTED.DAT`

```
Y 1988
R CHICAGO
P JONES 900
P GREGORY 400
R BATON ROUGE
P RODRIGUEZ 300
P SMITH 333
P GRAU 100
```

The record types are related to each other hierarchically. Year records represent the highest level in the hierarchy, since the year value 1988 applies to each salesperson in the file (only one year record is used in this example). Region records are intermediate-level records; region names apply to salesperson records that occur before the next region record in the file. For example, Chicago applies to salespersons Jones and Gregory. Baton Rouge applies to Rodriguez, Smith, and Grau. Person records represent the lowest level in the hierarchy. The information they contain—salesperson and unit sales—defines a case. Nested file structures minimize redundant information in a data file. For example, 1988 and Baton Rouge appear several times in the rectangular file, but only once in the nested file.

Since each record in the nested file has a code that indicates record type, you can use the `FILE TYPE` and `RECORD TYPE` commands to define the nested sales data:

```
FILE TYPE NESTED FILE='NESTED.DAT' RECORD=#TYPE 1 (A)
RECORD TYPE 'Y'.
DATA LIST / YEAR 5-8.

RECORD TYPE 'R'.
DATA LIST / REGION 5-15 (A).
```

```

RECORD TYPE 'P'.
DATA LIST / SALESPER 5-15 (A) SALES 20-23

END FILE TYPE.

```

- FILE TYPE indicates that data are in nested form in the file *NESTED.DAT*.
- RECORD defines the record type variable as string variable #TYPE in column 1. #TYPE is defined as scratch variable so it won't be saved in the active dataset.
- One pair of RECORD TYPE and DATA LIST statements is specified for each record type in the file. The first pair of RECORD TYPE and DATA LIST statements defines the variable YEAR in columns 5 through 8 on every year record. The second pair defines the string variable REGION on region records. The final pair defines SALESPER and SALES on person records.
- The order of RECORD TYPE statements defines the hierarchical relationship among the records. The first RECORD TYPE defines the highest-level record type. The next RECORD TYPE defines the next highest level, and so forth. The last RECORD TYPE defines a case in the active dataset.
- END FILE TYPE signals the end of file definition.
- In processing nested data, the program reads each record type you define. Information on the highest and intermediate-level records is spread to cases to which the information applies. The output from the LIST command is identical to that for the rectangular file.

### ***Nested Files with Missing Records***

In a nested file, some cases may be missing one or more record types defined in RECORD TYPE commands. For example, in the figure below the region record for salespersons Jones and Gregory is missing.

Figure C-4  
File *NESTED.DAT* with missing records

```

Y 1988
P JONES 900
P GREGORY 400
R BATON ROUGE
P RODRIGUEZ 300
P SMITH 333
P GRAU 100

```

The program assigns missing values to variables that are not present for a case. Using the modified *NESTED.DAT* file, the commands in the previous example produce the output shown below. You can see that the program assigned missing values to *REGION* for Jones and Gregory.

Figure C-5  
LIST output for nested data with missing records

```

YEAR REGION  SALESPER SALES

1988      JONES  900
1988      GREGORY 400

```

```
1988 BATON ROUGE RODRIGUEZ 300
1988 BATON ROUGE SMITH 333
1988 BATON ROUGE GRAU 100
```

You may want to examine cases with missing records, since these cases may indicate data errors. If you add the `MISSING=WARN` subcommand to your `FILE TYPE` command, the program prints a warning message when a case is missing a defined record type. For example, the program would print two warnings when processing data in `NESTED.DAT` with missing records. When `MISSING` is set to `WARN`, cases are built in the same way as when the default setting (`NOWARN`) is in effect.

## Grouped Data

In a grouped file, a case has several records that are grouped together in the file. You can use `DATA LIST` to define a grouped file if each case has the same number of records and records appear in the same order for each case. You can use `FILE TYPE GROUPED` whether the number of records per case and record order are fixed or vary. However, `FILE TYPE GROUPED` requires that each record have a case identifier and a record code.

## Using DATA LIST

The following table shows the organization of a grouped data file containing school subject scores for three students. Each student has three data records, and each record contains a score. The first record for each student also contains a case identifier. Records for each case are grouped together. Student 1 records appear first, followed by records for student 2 and student 3.

Record order determines whether a score is a reading, math, or science score. The reading score appears on the first record for a case, the math score appears on the second record, and the science score appears on the third record.

**Table C-1**  
Data for `GROUPED.DAT`

Student	Score
1	58
	59
	97
2	43
	88
	45
3	67
	75
	90

Since each case has the same number of records and record order is fixed across cases, you can use `DATA LIST` to define the student data:

```
DATA LIST FILE='GROUPED.DAT' RECORDS=3
  /STUDENT 1 READING 5-6
  /MATH 5-6
  /SCIENCE 5-6.
```

LIST.

- DATA LIST indicates that data are in file *GROUPED.DAT*.
- RECORDS defines three records per case. The program reads student ID number (*STUDENT*) and reading score (*READING*) in the first record for a case. Math and science scores are read in the second and third records.
- The output from the LIST command is shown below.

Figure C-6

LIST output for *GROUPED.DAT*

STUDENT READING MATH SCIENCE

```
1  58  59  97
2  43  88  45
3  67  75  90
```

## Using FILE TYPE GROUPED

To use FILE TYPE GROUPED to define a grouped file, each record must have a case identifier and a record code. In the following commands, each data record contains a student ID number coded 1, 2, or 3 and a code indicating whether the score on that record is a reading (R), math (M), or science (S) score:

```
FILE TYPE GROUPED RECORD=#REC 3(A) CASE=STUDENT 1.

RECORD TYPE 'R'.
DATA LIST / READING 5-6.

RECORD TYPE 'M'.
DATA LIST / MATH 5-6.

RECORD TYPE 'S'.
DATA LIST / SCIENCE 5-6.

END FILE TYPE.

BEGIN DATA
1 R 58
1 M 59
1 S 97
2 R 43
2 M 88
2 S 45
3 R 67
3 M 75
3 S 90
END DATA.

LIST.
```

- FILE TYPE indicates that data are in grouped format. RECORD defines the variable containing record codes as string variable #REC in column 3. CASE defines the case identifier variable *STUDENT* in the first column of each record.

- One pair of `RECORD TYPE` and `DATA LIST` statements appears for each record type in the file. The program reads reading score in every R record, math score in M records, and science score in S records.
- `END FILE TYPE` signals the end of file definition.
- `BEGIN DATA` and `END DATA` indicate that data are inline.
- The output from `LIST` is identical to the output using `DATA LIST`.

`FILE TYPE GROUPED` is most useful when record order varies across cases and when cases have missing or duplicate records. In the modified data shown below, only case 1 has all three record types. Also, record order varies across cases. For example, the first record for case 1 is a science record, whereas the first record for cases 2 and 3 is a reading record.

Table C-2

*Modified grouped data file*

Student	Subject	Score
1	S	97
1	R	58
1	M	59
2	R	43
3	R	67
3	M	75

You can use the same `FILE TYPE` commands as above to read the modified file. As shown in the output from `LIST` below, the program assigns missing values to variables that are missing for a case.

Figure C-7

*LIST output for GROUPED.DAT*

STUDENT READING MATH SCIENCE

```
1  58  59  97
2  43  .   .
3  67  75  .
```

By default, the program generates a warning message when a case is missing a defined record type in a grouped file or when a record is not in the same order as in `RECORD TYPE` commands. Thus, four warnings are generated when the commands for the previous example are used to read the modified `GROUPED.DAT` file. You can suppress these warnings if you add the optional specifications `MISSING=NOWARN` and `ORDERED=NO` on your `FILE TYPE` command.

In the modified `GROUPED.DAT` file, the case identifier `STUDENT` appears in the same column position in each record. When the location of the case identifier varies for different types of records, you can use the `CASE` option of the `RECORD TYPE` command to specify different column positions for different records. For example, suppose the case identifier appears in first column position on reading and science records and in column 2 in math records. You could use the following commands to define the data:

```
FILE TYPE GROUPED RECORD=#REC 3 (A) CASE=STUDENT 1.

RECORD TYPE 'R' .
DATA LIST / READING 5-6.
```

```

RECORD TYPE 'M' CASE=2.
DATA LIST / MATH 5-6.

RECORD TYPE 'S'.
DATA LIST / SCIENCE 5-6.

END FILE TYPE.

BEGIN DATA
1 S 97
1 R 58
  1M 59
2 R 43
3 R 67
  3M 75
END DATA.

LIST.

```

- `FILE TYPE` indicates that the data are in grouped format. `RECORD` defines the variable containing record codes as string variable `#REC`. `CASE` defines the case identifier variable as `STUDENT` in the first column of each record.
- One pair of `RECORD TYPE` and `DATA LIST` statements is coded for each record type in the file.
- The `CASE` specification on the `RECORD TYPE` statement for math records overrides the `CASE` value defined on `FILE TYPE`. Thus, the program reads `STUDENT` in column 2 in math records and column 1 in other records.
- `END FILE TYPE` signals the end of file definition.
- `BEGIN DATA` and `END DATA` indicate that data are inline.
- The output from `LIST` is identical to the output above.

## Mixed Files

In a mixed file, different types of cases have different kinds of records. You can use `FILE TYPE MIXED` to read each record or a subset of records in a mixed file.

### Reading Each Record in a Mixed File

The following table shows test data for two hypothetical elementary school students referred to a remedial education teacher. Student 1, who was thought to need special reading attention, took reading tests (word identification and comprehension tests). The second student completed writing tests (handwriting, spelling, vocabulary, and grammar tests). Test code (`READING` or `WRITING`) indicates whether the record contains reading or writing scores.

Table C-3  
Academic test data for two students

Student 1				
Test	ID	Grade	Word	Compre
READING	1	04	65	35
Student 2				

Test	ID	Grade	Handwrit	Spelling	Vocab	Grammar
WRITING	2	03	50	55	30	25

The following commands define the test data:

```
FILE TYPE MIXED RECORD=TEST 1-7 (A) .

RECORD TYPE 'READING' .
DATA LIST / ID 9-10 GRADE 12-13 WORD 15-16 COMPRE 18-19.

RECORD TYPE 'WRITING' .
DATA LIST / ID 9-10 GRADE 12-13 HANDWRIT 15-16 SPELLING 18-19
          VOCAB 21-22 GRAMMAR 24-25.
END FILE TYPE.

BEGIN DATA
READING 1 04 65 35
WRITING 2 03 50 55 30 25
END DATA.

LIST.
```

- FILE TYPE specifies that the data contain mixed record types. RECORD reads the record identifier (variable *TEST*) in columns 1 through 7.
- One pair of RECORD TYPE and DATA LIST statements is coded for each record type in the file. The program reads variables *ID*, *GRADE*, *WORD*, and *COMPRE* in the record in which the value of *TEST* is *READING*, and *ID*, *GRADE*, *HANDWRIT*, *SPELLING*, *VOCAB*, and *GRAMMAR* in the *WRITING* record.
- END FILE TYPE signals the end of file definition.
- BEGIN DATA and END DATA indicate that data are inline. Data are mixed, since some column positions contain different variables for the two cases. For example, word identification score is recorded in columns 15 and 16 for student 1. For student 2, handwriting score is recorded in these columns.
- The following figure shows the output from LIST. Missing values are assigned for variables that are not recorded for a case.

Figure C-8

*LIST* output for mixed file

```
TEST ID GRADE WORD COMPRE HANDWRIT SPELLING VOCAB GRAMMAR

READING 1 4 65 35 . . . .
WRITING 2 3 . . 50 55 30 25
```

## Reading a Subset of Records in a Mixed File

You may want to process a subset of records in a mixed file. The following commands read only the data for the student who took reading tests:

```
FILE TYPE MIXED RECORD=TEST 1-7 (A) .

RECORD TYPE 'READING' .
DATA LIST / ID 9-10
          GRADE 12-13
          WORD 15-16
```



```

                COMPRE 18-19.

RECORD TYPE 'WRITING'.
DATA LIST / ID      9-10
           GRADE   12-13
           HANDWRIT 15-16
           SPELLING 18-19
           VOCAB   21-22
           GRAMMAR 24-25.

END FILE TYPE.

BEGIN DATA
READING 1  04 65 35
WRITING 2  03 50 55 30 25
END DATA.

LIST.

```

- `FILE TYPE` specifies that data contain mixed record types. `RECORD` defines the record identification variable as *TEST* in columns 1 through 7.
- `RECORD TYPE` defines variables on reading records. Since the program skips all record types that are not defined by default, the case with writing scores is not read.
- `END FILE TYPE` signals the end of file definition.
- `BEGIN DATA` and `END DATA` indicate that data are inline. Data are identical to those in the previous example.
- The following figure shows the output from `LIST`.

Figure C-9

*LIST* output for reading record

```

TEST  ID GRADE WORD COMPRE

READING 1  4 65 35

```

## Repeating Data

You can use the `REPEATING DATA` command to read files in which each record contains repeating groups of variables that define several cases. Command syntax depends on whether the number of repeating groups is fixed across records.

### Fixed Number of Repeating Groups

The following table shows test score data for students in three classrooms. Each record contains a classroom number and two pairs of student ID and test score variables. For example, in class 101, student 182 has a score of 12 and student 134 has a score of 53. In class 103, student 15 has a score of 87 and student 203 has a score of 69. Each pair of ID and score variables is a repeating group, since these variables appear twice on each record.

Table C-4

*Data in REPEAT.DAT file*

Class	ID	Score	ID	Score
101	182	12	134	53

Class	ID	Score	ID	Score
102	99	112	200	150
103	15	87	203	69

The following commands generate a active dataset in which one case is built for each occurrence of *SCORE* and *ID*, and classroom number is spread to each case on a record.

```
INPUT PROGRAM.
DATA LIST / CLASS 3-5.
REPEATING DATA STARTS=6 / OCCURS=2
  /DATA STUDENT 1-4 SCORE 5-8.
END INPUT PROGRAM.
```

```
BEGIN DATA
  101 182 12 134 53
  102 99 112 200 150
  103 15 87 203 69
END DATA.
```

```
LIST.
```

- INPUT PROGRAM signals the beginning of data definition.
- DATA LIST defines variable *CLASS*, which is spread to each student on a classroom record.
- REPEATING DATA specifies that the input file contains repeating data. *STARTS* indicates that repeating data begin in column 6. *OCCURS* specifies that the repeating data group occurs twice in each record.
- DATA defines variables that are repeated (*STUDENT* and *SCORE*). The program begins reading the first repeating data group in column 6 (the value of *STARTS*). Since the value of *OCCURS* is 2, the program reads the repeating variables a second time, beginning in the next available column (column 14).
- END INPUT PROGRAM signals the end of data definition.
- BEGIN DATA and END DATA specify that data are inline.
- The output from LIST is shown below. Each student is a separate case.

Figure C-10  
*LIST* output for repeating data

```
CLASS STUDENT SCORE
101 182 12
101 134 53
102 99 112
102 200 150
103 15 87
103 203 69
```

### ***Varying Number of Repeating Groups***

To use REPEATING DATA to define a file in which the number of repeating data groups varies across records, your data must contain a variable indicating the number of repeating data groups on a record. The following commands define such a file:

```
INPUT PROGRAM.
```

```
DATA LIST / #NUM 1 CLASS 3-5.
REPEATING DATA STARTS=6 / OCCURS=#NUM
  /DATA STUDENT 1-4 SCORE 5-8.
END INPUT PROGRAM.
```

```
BEGIN DATA
3 101 182 12 134 53 199 30
2 102 99 112 200 150
1 103 15 87
END DATA.
```

```
LIST.
```

- INPUT PROGRAM signals the beginning of data definition.
- DATA LIST defines variables *CLASS* in columns 3 through 5 and *#NUM*, a scratch variable in column 1 that contains the number of repeating data groups in a record.
- REPEATING DATA specifies that the input file contains repeating data. *STARTS* indicates that repeating data begin in column 6. *OCCURS* sets the number of repeating groups on a record equal to the value of *#NUM*.
- DATA defines variables that are repeated. Since *#NUM* is 3 in the first and third records, the program reads three sets of *STUDENT* and *SCORE* variables in these records. *STUDENT* and *SCORE* are read twice in record 2.
- END INPUT PROGRAM signals the end of data definition.
- Data appear between BEGIN DATA and END DATA.
- The following figure shows the output from LIST.

Figure C-11  
LIST output

```
CLASS STUDENT SCORE
101 182 12
101 134 53
101 199 30
102 99 112
103 15 87
```

If your data file does not have a variable indicating the number of repeating data groups per record, you can use the LOOP and REREAD commands to read the data, as in:

```
INPUT PROGRAM.
DATA LIST / CLASS 3-5 #ALL 6-29 (A).
LEAVE CLASS.

LOOP #I = 1 TO 17 BY 8 IF SUBSTR(#ALL, #I, 8) NE ''.
- REREAD COLUMN = #I + 5.
- DATA LIST / STUDENT 1-4 SCORE 5-8.
- END CASE.
END LOOP.
END INPUT PROGRAM.

BEGIN DATA
  101 182 12 134 53 199 30
  102 99 112 200 150
  103 15 87
END DATA.

LIST.
```

- `INPUT PROGRAM` signals the beginning of data definition.
- `DATA LIST` reads `CLASS` and `#ALL`, a temporary string variable that contains all of the repeating data groups for a classroom record. The column specifications for `#ALL` (6 through 29) are wide enough to accommodate the classroom record with the most repeating data groups (record 1).
- `LOOP` and `END LOOP` define an index loop. As the loop iterates, the program successively reads eight-character segments of `#ALL`, each of which contains a repeating data group or an empty field. The program reads the first eight characters of `#ALL` in the first iteration, the second eight characters in the second iteration, and so forth. The loop terminates when the program encounters an empty segment, which means that there are no more repeating data groups on a record.
- In each iteration of the loop in which an `#ALL` segment is not empty, `DATA LIST` reads `STUDENT` and `SCORE` in a classroom record. The program begins reading these variables in the first record, in the starting column specified by `REREAD COLUMN`. For example, in the first iteration, the program reads `STUDENT` and `SCORE` beginning in column 6. In the second iteration, the program reads `STUDENT` and `SCORE` starting in column 14 of the same record. When all repeating groups have been read for a record, loop processing begins on the following record.
- `END CASE` creates a new case for each repeating group.
- `REREAD` causes `DATA LIST` to read repeating data groups in the same record in which it last read `CLASS`. Without `REREAD`, each execution of `DATA LIST` would begin on a different record.
- `LEAVE` preserves the value of `CLASS` across the repeating data groups on a record. Thus, the same class number is read for each student on a classroom record.
- `INPUT PROGRAM` signals the beginning of data definition.
- `BEGIN DATA` and `END DATA` indicate that the data are inline. The data are identical to those in the previous example except that they do not contain a variable indicating the number of repeating groups per record.
- These commands generate the same output as shown in the figure above.

# Using the Macro Facility

A macro is a set of commands that generates customized command syntax. Using macros can reduce the time and effort needed to perform complex and repetitive data analysis tasks.

Macros have two parts: a **macro definition**, which indicates the beginning and end of the macro and gives a name to the macro, and a **macro body**, which contains regular commands or macro commands that build command syntax. When a macro is invoked by the **macro call**, syntax is generated in a process called **macro expansion**. Then the generated syntax is executed as part of the normal command sequence.

This chapter shows how to construct macros that perform three data analysis tasks. In the first example, macros facilitate a file-matching task. In Example 2, macros automate a specialized statistical operation (testing a sample correlation coefficient against a nonzero population correlation coefficient). Macros in Example 3 generate random data. As shown in the following table, each example demonstrates various features of the macro facility. For information on specific macro commands, see the `DEFINE` command.

Table D-1  
*Macro features*

	Example 1	Example 2	Example 3
Macro argument			
Keyword	x	x	x
Default values	x		x
None	x		x
String manipulation	x		x
Looping			
Index	x		x
List processing		x	
Direct assignment	x		x

## Example 1: Automating a File-Matching Task

The following figure shows a listing of 1988 sales data for salespeople working in different regions. The listing shows that salesperson Jones sold 900 units in the Chicago sales territory, while Rodriguez sold 300 units in Baton Rouge.

Figure D-1  
*Listing of data file SALES88.SAV*

```
YEAR REGION  SALESPER SALES
1988 CHICAGO JONES    900
1988 CHICAGO GREGORY  400
```

```

1988 BATON ROUGE RODRIGUEZ 300
1988 BATON ROUGE SMITH 333
1988 BATON ROUGE GRAU 100

```

You can use command syntax shown below to obtain each salesperson's percentage of total sales for their region.

#### Figure D-2

*Commands for obtaining sales percentages*

```
GET FILE = 'SALES88.SAV'.
```

```
SORT CASES BY REGION.
```

```
AGGREGATE OUTFILE = 'TOTALS.SAV'
/PRESORTED
/BREAK = REGION
/TOTAL@ = SUM(SALES).
```

```
MATCH FILES FILE=*
/TABLE = 'TOTALS.SAV'
/BY REGION.
```

```
COMPUTE PCT = 100 * SALES / TOTAL@.
```

```
TITLE 1988 DATA.
LIST.
```

- The `GET` command opens *SALES88.SAV*. This file becomes the active dataset.
- `SORT CASES` sorts the active dataset in ascending alphabetical order by *REGION*.
- The `AGGREGATE` command saves total sales (variable *TOTAL@*) for each region in file *TOTALS.SAV*.
- `MATCH FILES` appends the regional totals to each salesperson's record in the active dataset. (See the `MATCH FILES` command for more information on matching files.)
- `COMPUTE` obtains the percentage of regional sales (*PCT*) for each salesperson.
- The `LIST` command output displayed below shows that Rodriguez sold 41% of the products sold in Baton Rouge. Gregory accounted for 31% of sales in the Chicago area.

#### Figure D-3

*Listing of data file SALES88.SAV*

```

YEAR REGION  SALESPER SALES TOTAL@  PCT
1988 BATON ROUGE RODRIGUEZ 300 733.00 41.00
1988 BATON ROUGE SMITH 333 733.00 45.00
1988 BATON ROUGE GRAU 100 733.00 14.00
1988 CHICAGO JONES 900 1300.00 69.00
1988 CHICAGO GREGORY 400 1300.00 69.00

```

The following figure shows a macro that issues the commands for obtaining sales percentages. The macro consists of the commands that produce sales percentages imbedded between macro definition commands `DEFINE` and `!ENDDEFINE`.

Figure D-4

*!TOTMAC macro*

```

DEFINE !TOTMAC ().

GET FILE = 'SALES88.SAV'.

SORT CASES BY REGION.

AGGREGATE OUTFILE = 'TOTALS.SAV'
/PRESORTED
/BREAK = REGION
/TOTAL@ = SUM(SALES).

MATCH FILES FILE = *
/TABLE = 'TOTALS.SAV'
/BY REGION.

COMPUTE PCT = 100 * SALES / TOTAL@.

TITLE 1988 DATA.
LIST.

!ENDDDEFINE.

!TOTMAC.

```

- Macro definition commands `DEFINE` and `!ENDDDEFINE` signal the beginning and end of macro processing. `DEFINE` also assigns the name `!TOTMAC` to the macro (the parentheses following the name of the macro are required). The macro name begins with an exclamation point so that the macro does not conflict with that of an existing variable or command. Otherwise, if the macro name matched a variable name, the variable name would invoke the macro whenever the variable name appeared in the command stream.
- Commands between `DEFINE` and `!ENDDDEFINE` constitute the macro body. These commands, which produce sales percentages, are identical to the commands in Figure D-2.
- The final statement, `!TOTMAC`, is the **macro call**, which invokes the macro. When the program reads the macro call, it issues the commands in the macro body. Then these commands are executed, generating identical output.

While the macro shows you how to construct a simple macro, it doesn't reduce the number of commands needed to calculate regional percentages. However, you can use macro features such as looping to minimize coding in more complicated tasks. For example, let's say that in addition to the 1988 data, you have sales data for 1989 (*SALES89.SAV*), and each file contains the variables *REGION*, *SALESPER*, and *SALES*. The modified `!TOTMAC` macro below calculates regional sales percentages for each salesperson for 1988 and 1989.

Figure D-5

*!TOTMAC macro with index loop*

```

DEFINE !TOTMAC ().

!DO !! = 88 !TO 89.

```

```

- GET FILE = !CONCAT('SALES', !I, '.SAV').
- SORT CASES BY REGION.
- AGGREGATE OUTFILE = 'TOTALS.SAV'
  /PRESORTED
  /BREAK = REGION
  /TOTAL@ = SUM(SALES).
- MATCH FILES FILE = *
  /TABLE = 'TOTALS.SAV'
  /BY REGION.
- COMPUTE PCT= 100 * SALES / TOTAL@.
- !LET !YEAR = !CONCAT('19',!I).
- TITLE !YEAR DATA.
- LIST.
!DOEND.

```

```
!ENDDDEFINE.
```

```
!TOTMAC.
```

- `DEFINE` and `!ENDDDEFINE` signal the beginning and end of macro processing.
- Commands `!DO` and `!DOEND` define an **index loop**. Commands between `!DO` and `!DOEND` are issued once in each iteration of the loop. The value of **index variable** `!I`, which changes in each iteration, is 88 in the first iteration and 89 in the second (final) iteration.
- In each iteration of the loop, the `GET` command opens a data file. The name of the file is constructed using the **string manipulation function** `!CONCAT`, which creates a string that is the concatenation of `SALES`, the value of the index variable, and `.sav`. Thus the file `SALES88.SAV` is opened in the first iteration.
- Commands between `AGGREGATE` and `COMPUTE` calculate percentages on the active dataset. These commands are identical to those in Figure D-4.
- Next, a customized title is created. In the first iteration, the **direct assignment** command `!LET` assigns a value of 1988 to the macro variable `!YEAR`. This variable is used in the `TITLE` command on the following line to specify a title of *1988 DATA*.
- The `LIST` command displays the contents of each variable.
- In the second iteration of the loop, commands display percentages for the 1989 data file. The output from the `!TOTMAC` macro is shown below. Note that the listing for 1988 data is the same as before.

Figure D-6

*Regional sales percentages for 1988 and 1989*

1988 DATA

YEAR	REGION	SALESPER	SALES	TOTAL@	PCT
1988	BATON ROUGE	RODRIGUEZ	300	733.00	41.00
1988	BATON ROUGE	SMITH	333	733.00	45.00
1988	BATON ROUGE	GRAU	100	733.00	14.00
1988	CHICAGO	JONES	900	1300.00	69.00
1988	CHICAGO	GREGORY	400	1300.00	69.00



```

1989 DATA
YEAR REGION  SALESPER SALES TOTAL@  PCT
1989 BATON ROUGE GRAU  320 1459.00 22.00
1989 BATON ROUGE SMITH  800 1459.00 55.00
1989 BATON ROUGE RODRIGUEZ 339 1459.00 23.00
1989 CHICAGO  JONES  300 1439.00 21.00
1989 CHICAGO  STEEL  899 1439.00 62.00
1989 CHICAGO  GREGORY 240 1439.00 17.00

```

Let's look at another application of the !TOTMAC macro, one that uses **keyword arguments** to make the application more flexible. The following figure shows the number of absences for students in two classrooms. Let's say you want to calculate deviation scores indicating how many more (or fewer) times a student was absent than the average student in his or her classroom. The first step in obtaining deviation scores is to compute the average number of absences per classroom. We can use the !TOTMAC macro to compute classroom means by modifying the macro so that it computes means and uses the absences data file (*SCHOOL.SAV*) as input.

Figure D-7  
Listing of file *SCHOOL.SAV*

```

CLASS STUDENT ABSENT
101 BARRY G  3
101 JENNI W  1
101 ED F     2
101 JOHN O   8
102 PAUL Y   2
102 AMY G    3
102 JOHN D  12
102 RICH H   4

```

The !TOTMAC macro below can produce a variety of group summary statistics such as sum, mean, and standard deviation for any IBM® SPSS® Statistics data file. In the macro call you specify values of keyword arguments indicating the data file (*FILE*), the break (grouping) variable (*BREAKVR*), the summary function (*FUNC*), and the variable to be used as input to the summary function (*INVAR*). For example, to obtain mean absences for each classroom, we specify *SCHOOL.SAV* as the data file, *CLASS* as the break variable, *MEAN* as the summary function, and *ABSENT* as the variable whose values are to be averaged.

Figure D-8  
*!TOTMAC* macro with keyword arguments

```

DEFINE !TOTMAC ( BREAKVR = !TOKENS(1)
                /FUNC = !TOKENS(1)
                /INVAR = !TOKENS(1)
                /TEMP = !TOKENS(1) !DEFAULT(TOTALS.SAV)
                /FILE = !CMDEND).
GET FILE = !FILE.
SORT CASES BY !BREAKVR.
AGGREGATE OUTFILE = '!TEMP'
/PRESORTED
/BREAK = !BREAKVR

```

```

        !CONCAT(!FUNC, '@') = !FUNC(!INVAR).

MATCH FILES FILE = *
        /TABLE = 'ITEMP'
        /BY !BREAKVR.

!ENDDDEFINE.

!TOTMAC BREAKVR=CLASS FUNC=MEAN INVAR=ABSENT FILE=SCHOOL.SAV.

COMPUTE DIFF = ABSENT-MEAN@.

LIST.

!TOTMAC BREAKVR=REGION FUNC=SUM INVAR=SALES FILE=SALES89.SAV.

COMPUTE PCT = 100 * SALES / SUM@.

LIST.

```

- The syntax for declaring keyword arguments follows the name of the macro in `DEFINE`.
- `!TOKENS (1)` specifies that the value of an argument is a string following the name of the argument in the macro call. Thus the first macro call specifies `CLASS` as the value of `BREAKVR`, `MEAN` as the value of `FUNC`, and `ABSENT` as the value of `INVAR`.
- `!CMDEND` indicates that the value for `FILE` is the remaining text in the macro call (*SCHOOL.SAV*).
- `TEMP` is an optional argument that names an intermediate file to contain the summary statistics. Since `TEMP` is not assigned a value in the macro call, summary statistics are written to the default intermediate file (*TOTALS.SAV*).
- In the body of the macro, `GET FILE` opens *SCHOOL.SAV*.
- `SORT CASES` sorts the file by *CLASS*.
- `AGGREGATE` computes the mean number of absences for each class. The name of the variable containing the means (*MEAN@*) is constructed using the `!CONCAT` function, which concatenates the value of `FUNC` and the `@` symbol.
- `MATCH FILES` appends the means to student records.
- `COMPUTE` calculates the deviation from the classroom mean for each student (variable *DIFF*).
- `LIST` displays the deviation scores, as shown in the output below. For example, John D., who was absent 12 times, had 6.75 more absences than the average student in classroom 102. Rich H., who was absent 4 times, had 1.25 fewer absences than the average student in classroom 102.
- The second macro call and remaining commands generate regional sales percentages for the 1989 sales data. As shown below, percentages are identical to those displayed in the bottom half of Figure D-6.

Figure D-9  
Student absences and 1989 sales percentages

```

CLASS STUDENT ABSENT MEAN@ DIFF
101 BARRY G 3 3.50 -.50
101 JENNI W 1 3.50 -2.50
101 ED F 2 3.50 -1.50
101 JOHN O 8 3.50 4.50
102 PAUL Y 2 5.25 -3.25
102 AMY G 3 5.25 -2.25
102 JOHN D 12 5.25 6.75
102 RICH H 4 5.25 -1.25

1989 DATA
YEAR REGION SALES PER SALES TOTAL@ PCT
1989 BATON ROUGE GRAU 320 1459.00 22.00
1989 BATON ROUGE SMITH 800 1459.00 55.00
1989 BATON ROUGE RODRIGUEZ 339 1459.00 23.00
1989 CHICAGO JONES 300 1439.00 21.00
1989 CHICAGO STEEL 899 1439.00 62.00
1989 CHICAGO GREGORY 240 1439.00 17.00

```

You can modify the macro call to specify a different data file, input variable, break variable, or summary statistic. To get a different summary statistic (such as standard deviation), change the value of `FUNC` (see the `AGGREGATE` command for more information on summary functions available in the `AGGREGATE` procedure).

## Example 2: Testing Correlation Coefficients

While the program provides a large variety of statistical procedures, some specialized operations require the use of `COMPUTE` statements. For example, you may want to test a sample correlation coefficient against a population correlation coefficient. When the population coefficient is nonzero, you can compute a  $Z$  statistic to test the hypothesis that the sample and population values are equal. The formula for  $Z$  is

$$Z = \frac{0.5 \ln \left[ \frac{1+r}{1-r} \right] - 0.5 \ln \left[ \frac{1+p_0}{1-p_0} \right]}{1/(\sqrt{n-3})}$$

where  $r$  is the sample correlation coefficient,  $p_0$  is the population coefficient,  $n$  is the size of the sample from which  $r$  is obtained, and  $\ln$  signifies the natural logarithm function.  $Z$  has approximately the standard normal distribution.

Let's say you want to test an  $r$  of 0.66 obtained from a sample of 30 cases against a population coefficient of 0.85. The following figure shows commands for displaying  $Z$  and its two-tailed probability.

Figure D-10  
Commands for computing  $Z$  statistic

```
DATA LIST FREE / R N P.

BEGIN DATA
.66 30 .85
END DATA.

COMPUTE #ZR = .5* (LN ((1 + R) / (1 - R))).
COMPUTE #ZP = .5* (LN ((1 + P) / (1 - P))).

COMPUTE Z = (#ZR-#ZP)/(1/(SQRT(N-3))).
COMPUTE PROB = 2*(1-CDFNORM(ABS(Z))).

FORMAT PROB (F8.3).
LIST.
```

- `DATA LIST` defines variables containing the sample correlation coefficient ( $R$ ), sample size ( $N$ ), and population correlation coefficient ( $P$ ).
- `BEGIN DATA` and `END DATA` indicate that data are inline.
- `COMPUTE` statements calculate  $Z$  and its probability. Variables `#ZR` and `#ZP` are scratch variables used in the intermediate steps of the calculation.
- The `LIST` command output is shown below. Since the absolute value of  $Z$  is large and the probability is small, we reject the hypothesis that the sample was drawn from a population having a correlation coefficient of 0.85.

Figure D-11  
 $Z$  statistic and its probability

```
R    N    P    Z    PROB
.66 30.00 .85 -2.41 .016
```

If you use the  $Z$  test frequently, you may want to construct a macro like that shown below. The `!CORRTST` macro computes  $Z$  and probability values for a sample correlation coefficient, sample size, and population coefficient specified as values of keyword arguments.

Figure D-12  
`!CORRTST` macro

```
DEFINE !CORRTST ( R = !TOKENS(1)
                /N = !TOKENS(1)
                /P = !TOKENS(1)).

INPUT PROGRAM.
- END CASE.
- END FILE.
END INPUT PROGRAM.

COMPUTE #ZR = .5* (LN ((1 + !R) / (1 - !R))).
COMPUTE #ZP = .5* (LN ((1 + !P) / (1 - !P))).

COMPUTE Z = (#ZR-#ZP) / (1/(SQRT(!N-3))).
COMPUTE PROB = 2*(1-CDFNORM(ABS(Z))).
```

```

FORMAT PROB(F8.3).

TITLE SAMPLE R=!R, N=!N, POPULATION COEFFICIENT=!P.

LIST.

!ENDDDEFINE.

!CORRTST R=.66 N=30 P=.85.
!CORRTST R=.50 N=50 P=.85.

```

- `DEFINE` names the macro as `!CORRTST` and declares arguments for the sample correlation coefficient ( $R$ ), the sample size ( $N$ ), and the population correlation coefficient ( $P$ ).
- `!TOKENS(1)` specifies that the value of an argument is a string that follows the name of the argument in the macro call. Thus the first macro call specifies values of 0.66, 30, and 0.85 for  $R$ ,  $N$ , and  $P$ .
- Commands between `INPUT PROGRAM` and `END INPUT PROGRAM` create an active dataset with one case. `COMPUTE` statements calculate the  $Z$  statistic and its probability using the values of macro arguments  $R$ ,  $N$ , and  $P$ . (`INPUT PROGRAM` commands would not be needed if `COMPUTE` statements operated on values in an existing file or inline data, rather than macro arguments.)
- A customized `TITLE` shows displays the values of macro arguments used in computing  $Z$ .
- The `LIST` command displays  $Z$  and its probability.
- The `!CORRTST` macro is called twice. The first invocation tests an  $r$  of 0.66 from a sample of 30 cases against a population coefficient of 0.85 (this generates the same  $Z$  value and probability as shown earlier). The second macro call tests an  $r$  of 0.50 from a sample of 50 cases against the same population correlation coefficient. The output from these macro calls is shown below.

Figure D-13  
Output from `!CORRTST`

```
SAMPLE R= .66 , N= 30 , POPULATION COEFFICIENT= .85
```

```

  Z  PROB
-2.41 .016

```

```
SAMPLE R= .50 , N= 50 , POPULATION COEFFICIENT= .85
```

```

  Z  PROB
-4.85 .000

```

The following figure shows a modified `!CORRTST` macro that you can use to test a sample  $r$  against each coefficient in a *list* of population coefficients.

Figure D-14  
`!CORRTST` macro with list-processing loop

```

DEFINE !CORRTST (R = !TOKENS(1)
                /N = !TOKENS(1)

```

```

        /P = !CMDEND).
- INPUT PROGRAM.
- END CASE.
- END FILE.
- END INPUT PROGRAM.

!DO !! !IN (!P).
- COMPUTE #ZR = .5* (LN ((1 + !R) / (1 - !R))).
- COMPUTE #ZP = .5* (LN ((1 + !I) / (1 - !I))).

- COMPUTE Z = (#ZR-#ZP)/(1/(SQRT(!N-3))).

- COMPUTE PROB=2*(1-CDFNORM(ABS(Z))).
- FORMAT PROB(F8.3).
- TITLE SAMPLE R=!R, N=!N, POPULATION COEFFICIENT=!I.
- LIST.
!DOEND.

!ENDDFINE.

!CORRTST R=.66 N=30 P=.20 .40 .60 .80 .85 .90.

```

- `DEFINE` names the macro as `!CORRTST` and declares arguments for the sample correlation coefficient (`R`), the sample size (`N`), and the population correlation coefficient (`P`).
- `!TOKENS (1)` specifies that the value of an argument is a string that follows the name of the argument in the macro call. Thus, the macro call specifies the value of `R` as 0.66 and `N` as 0.30.
- `!CMDEND` indicates that the value for `P` is the remaining text in the macro call. Thus the value of `P` is a list containing the elements 0.20, 0.40, 0.60, 0.80, 0.85, and 0.90.
- Commands `!DO !IN` and `!DOEND` define a **list-processing loop**. Commands in the loop compute one  $Z$  statistic for each element in the list of population coefficients. For example, in the first iteration  $Z$  is computed using 0.20 as the population coefficient. In the second iteration 0.40 is used. The same sample size (30) and  $r$  value (0.66) are used for each  $Z$  statistic.
- The output from the macro call is shown below. One  $Z$  statistic is displayed for each population coefficient.

Figure D-15

Output from modified `!CORRTST` macro

SAMPLE R= .66 , N= 30 , POPULATION COEFFICIENT= .20

```

      Z  PROB
3.07  .002

```

SAMPLE R= .66 , N= 30 , POPULATION COEFFICIENT= .40

```

      Z  PROB
1.92  .055

```

SAMPLE R= .66 , N= 30 , POPULATION COEFFICIENT= .60

```
Z  PROB
.52 .605
```

SAMPLE R= .66 , N= 30 , POPULATION COEFFICIENT= .80

```
Z  PROB
-1.59 .112
```

SAMPLE R= .66 , N= 30 , POPULATION COEFFICIENT= .85

```
Z  PROB
-2.41 .016
```

SAMPLE R= .66 , N= 30 , POPULATION COEFFICIENT= .90

```
Z  PROB
-3.53 .000
```

### Example 3: Generating Random Data

You can use command syntax to generate variables that have approximately a normal distribution. Commands for generating five standard normal variables (*X1* through *X5*) for 1000 cases are shown in the following figure. As shown in the output below, each variable has a mean of approximately 0 and a standard deviation of approximately 1.

Figure D-16  
*Data-generating commands*

```
INPUT PROGRAM.
- VECTOR X(5).
- LOOP #I = 1 TO 1000.
-   LOOP #J = 1 TO 5.
-     COMPUTE X(#J) = NORMAL(1).
-   END LOOP.
- END CASE.
- END LOOP.
- END FILE.
END INPUT PROGRAM.
```

DESCRIPTIVES VARIABLES X1 TO X5.

Figure D-17  
*Descriptive statistics for generated data*

Variable	Mean	Std Dev	Minimum	Maximum	N	Label
X1	-.01	1.02	-3.11	4.15	1000	
X2	.08	1.03	-3.19	3.22	1000	
X3	.02	1.00	-3.01	3.51	1000	

```
X4  .03  1.00 -3.35  3.19 1000
X5  -.01  .96 -3.34  2.91 1000
```

The !DATAGEN macro below issues the data-generating commands shown above.

**Figure D-18**

*!DATAGEN macro*

```
DEFINE !DATAGEN ().

INPUT PROGRAM.
- VECTOR X(5).
- LOOP #I = 1 TO 1000.
- LOOP #J = 1 TO 5.
- COMPUTE X(#J) = NORMAL(1).
- END LOOP.
- END CASE.
- END LOOP.
- END FILE.
END INPUT PROGRAM.

DESCRIPTIVES VARIABLES X1 TO X5.

!ENDDFINE.

!DATAGEN.
```

The data-generating commands are imbedded between macro definition commands. The macro produces the same data and descriptive statistics as shown above.

You can tailor the generation of normally distributed variables if you modify the !DATAGEN macro so it will accept keyword arguments, as shown in the following figure. The macro allows you to specify the number of variables and cases to be generated and the approximate standard deviation.

**Figure D-19**

*!DATAGEN macro with keyword arguments*

```
DEFINE !DATAGEN ( OBS =!TOKENS(1) !DEFAULT(1000)
                /VARS =!TOKENS(1) !DEFAULT(5)
                /SD =!CMDEND !DEFAULT(1)).
INPUT PROGRAM.
- VECTOR X(!VARS).
- LOOP #I = 1 TO !OBS.
- LOOP #J = 1 TO !VARS.
- COMPUTE X(#J) = NORMAL(!SD).
- END LOOP.
- END CASE.
- END LOOP.
- END FILE.
END INPUT PROGRAM.

!LET !LIST = !NULL.
!DO !I = 1 !TO !VARS.
- !LET !LIST = !CONCAT(!LIST, ' ', X, !I).
```



```
!DOEND.
```

```
DESCRIPTIVES VARIABLES !LIST.
```

```
!ENDDDEFINE.
```

```
!DATAGEN OBS=500 VARS=2 SD=1.
```

```
!DATAGEN.
```

- The `DEFINE` statement declares arguments that specify the number of cases (`OBS`), variables (`VARS`), and standard deviation (`SD`). By default, the macro creates 1000 cases with 5 variables that have a standard deviation of 1.
- Commands between `INPUT PROGRAM` and `END INPUT PROGRAM` generate the new data using values of the macro arguments.
- Commands `!LET` and `!DO/!DOEND` construct a variable list (`!LIST`) that is used in `DESCRIPTIVES`. The first `!LET` command initializes the list to a null (blank) string value. For each new variable, the index loop adds to the list a string of the form `X1`, `X2`, `X3`, and so forth. Thus, `DESCRIPTIVES` requests means and standard deviations for each new variable.
- The first macro call generates 500 cases with two standard normal variables. The second call requests the default number of variables, cases, and standard deviation. Descriptive statistics (not shown) are also computed for each variable.

As shown in the following figure, you can declare additional keyword arguments that allow you to specify the distribution (normal or uniform) of the generated data and a parameter value that is used as the standard deviation (for normally distributed data) or a range (for uniformly distributed data).

Figure D-20

*!DATAGEN macro with additional keyword arguments*

```
DEFINE !DATAGEN (OBS =!TOKENS(1) !DEFAULT(1000)
  /VARS =!TOKENS(1) !DEFAULT(5)
  /DIST =!TOKENS(1) !DEFAULT(NORMAL)
  /PARAM =!TOKENS(1) !DEFAULT(1)).

INPUT PROGRAM.
- VECTOR X(!VARS).
- LOOP #I = 1 TO !OBS.
-   LOOP #J = 1 TO !VARS.
-     COMPUTE X(#J) = !DIST(!PARAM).
-   END LOOP.
- END CASE.
- END LOOP.
- END FILE.
END INPUT PROGRAM.

!LET !LIST = !NULL.
!DO !I = 1 !TO !VARS.
- !LET !LIST = !CONCAT(!LIST, ' ', X, !I).
!DOEND.

DESCRIPTIVES VARIABLES !LIST.
!ENDDDEFINE.
```

```
!DATAGEN OBS=500 VARS=2 DIST=UNIFORM PARAM=2.
```

- The `DEFINE` statement declares arguments `OBS`, `VARS`, `DIST`, and `PARAM`. `OBS` and `VARS` represent the number of observations and cases to be generated. Arguments `DIST` and `PARAM` specify the shape and parameter of the distribution of generated data. By default, the macro generates 1000 observations with 5 standard normal variables.
- Statements between `INPUT PROGRAM` and `END INPUT PROGRAM` generate the new data using values of macro arguments.
- Remaining commands in the body of the macro obtain descriptive statistics for generated variables.
- The macro call in creates two approximately uniformly distributed variables with a range of 2. The output from the macro call is shown below.

Figure D-21

*Descriptive statistics for uniform variables*

Variable	Mean	Std Dev	Valid	Minimum	Maximum	N	Label
X1	.99	.57	.00	2.00	500		
X2	1.00	.57	.00	2.00	500		

# ***Canonical Correlation and Ridge Regression Macros***

Two macro routines are installed with for performing canonical correlation and ridge regression. Macros are inherently less robust than regular commands. Error checking is minimal, and small errors produce large numbers of uninformative messages, giving little clue to the real cause of the problem. Syntax must be entered exactly as documented:

- Enter subcommands in the order shown.
- Do not include a slash before the first subcommand.
- Enter keywords exactly as shown. Do not abbreviate them, and do not extend them (even if they look like abbreviations).
- Do not omit equal signs, but do not add them if none are displayed in the syntax diagram.
- Remember that these “command names” are actually macro calls, which are expanded into (lengthy) sequences of commands and Matrix language statements. The command name should not appear anywhere within the specifications of the command, for example within a variable name.
- These macros create new files in the current directory. If you do not have write permission to the current directory (this can happen if you are running the program on a network), change the current directory. You can do this by opening or saving a file on your own hard disk.
- You cannot run these macros if split-file processing is in effect.

## ***Canonical Correlation Macro***

```
INCLUDE '[installdir]/Samples/English/Canonical correlation.sps'.
CANCORR SET1=varlist1 /
        SET2=varlist2 / .
The two variable lists must be separated with a slash.
[installdir] is the installation directory.
```

## ***Ridge Regression Macro***

```
INCLUDE '[installdir]/Samples/English/Ridge regression.sps'.
RIDGEREG DEP=varname /ENTER = varlist
        [/START={0**}] [/STOP={1**}] [/INC={0.05**}]
        {value} {value} {value }
        [ /K=value] .
[installdir] is the installation directory.
```

# ***File Specifications for IBM SPSS Collaboration and Deployment Services Repository Objects***

Access to a IBM® SPSS® Collaboration and Deployment Services Repository is available with the Statistics Adapter option. Once a connection to a IBM SPSS Collaboration and Deployment Services Repository has been established, you can use a file specification for a IBM SPSS Collaboration and Deployment Services Repository object on nearly any keyword or subcommand where a file specification can be provided (including, but not limited to `FILE`, `INFILE`, or `OUTFILE` ). File specifications for IBM SPSS Collaboration and Deployment Services Repository objects can also be used in procedure dialog boxes that allow manual entry of a file specification—for example, XML model files or plan files for complex samples analyses.

A IBM SPSS Collaboration and Deployment Services Repository file specification has the following form:

```
SPSSCR://{directory path and/or file specification}
        [{#M.versionmarker}]
        {#L.versionlabel }
```

*Note:* Specification of descriptions (#D) and keywords (#K) is obsolete for release 17.0 and higher.

## ***Basic Specification***

The basic specification is the scheme name `SPSSCR` , followed by a colon, one or two slashes (forward or backward), and a directory path or file: for example:

```
GET FILE='SPSSCR://data/mydata.sav'.
```

- A IBM SPSS Collaboration and Deployment Services Repository file specification is included on a `GET` command. The latest version of the file `/data/mydata.sav` that is accessible by the current user will be retrieved from the currently connected IBM SPSS Collaboration and Deployment Services Repository.

## ***Syntax Rules***

- Paths can be specified with forward slashes (/) or backslashes (\).

- The optional components of a file specification—version marker (#M) and version label (#L)—can be specified in any order.
- When including an optional component, the character sequence that specifies the component—for example, #L—must be followed by a period (.).

### **File Extensions**

Files are stored to the IBM SPSS Collaboration and Deployment Services Repository with an associated MIME type that specifies the type of file; for example, IBM® SPSS® Statistics format data, or viewer output. Including a file extension, such as *.sav* or *.spv*, is not necessary in order for the correct MIME type to be determined; however, it is recommended.

- File extensions are not added to filenames when storing to a IBM SPSS Collaboration and Deployment Services Repository. For example, the following *SAVE* command results in the creation of a IBM SPSS Collaboration and Deployment Services Repository object with a MIME type of SPSS Statistics-format data, and a fully qualified path of */data/mydata*, but no *sav* extension.

```
SAVE OUTFILE='SPSSCR://data/mydata'.
```

- When retrieving IBM SPSS Collaboration and Deployment Services Repository objects, you must specify the filename exactly as provided when the file was stored. If the file was stored with an extension, then you must supply the extension. If the file was stored without an extension then do not include one.

### **Permissions**

- The *PERMISSIONS*, *SAVE*, and *XSAVE* commands will generate errors if you attempt to set permissions for IBM SPSS Collaboration and Deployment Services Repository objects. Permissions for IBM SPSS Collaboration and Deployment Services Repository objects are set with the *REPOSITORY ATTRIBUTES* command or using administration tools included with IBM® SPSS® Collaboration and Deployment Services.
- The *ERASE FILE* command will generate an error if you attempt to delete a IBM SPSS Collaboration and Deployment Services Repository object.

*Note:* For users upgrading from an IBM SPSS Collaboration and Deployment Services 3.5 system, be aware that command syntax that retrieved the version labelled *LATEST* (the default version retrieved by *GET FILE='filespec'*) in a 3.5 system will retrieve an earlier version in an IBM SPSS Collaboration and Deployment Services 4.x system if the current user does not have permission to view objects with the *LATEST* label.

## **Versions**

By default, the latest version that is accessible by the current user is returned when a file is retrieved. You can optionally specify a particular version to retrieve by including the version marker or version label, if one exists, for the desired version. Version markers and labels are always ignored if the specification describes a directory path, as opposed to a specific file.

**Version label (#L).** A label is used to easily identify a version; for example, *production* or *development*.

- The version label, if provided, is displayed in the Label column in the Select Version dialog box.

**Version marker (#M).** The version marker consists of an integer version number followed by a time stamp that uniquely specifies a version—for example, *1:2006-07-26 15:19:49.237*.

- The version marker is displayed in the Version column in the Select Version dialog box.

### Examples

```
*Retrieve a labelled version.
GET FILE='SPSSCR://data/mydata.sav#L.development'.
```

- The version of */data/mydata.sav* with the label *development* is retrieved from the current IBM® SPSS® Collaboration and Deployment Services Repository.

```
*Retrieve a version using its version marker.
OUTPUT OPEN FILE=
  'SPSSCR://output/myoutput.spv#M.0:2006-07-26 15:18:15.862'.
```

- The version of the output file */output/myoutput.spv* with version marker *0:2006-07-26 15:18:15.862* is retrieved.

## Using File Handles for IBM SPSS Collaboration and Deployment Services Repository Locations

You can define a file handle to a file or a directory in a IBM® SPSS® Collaboration and Deployment Services Repository and use that file handle when storing or retrieving.

```
*Define and use a file handle for a directory.
FILE HANDLE cust /NAME='SPSSCR://CustomerValue'.
GET FILE='cust/data2002.sav'.
```

- The handle *cust* is associated with the IBM SPSS Collaboration and Deployment Services Repository directory */CustomerValue*, and the GET command retrieves the latest version of the file *data2002.sav* from this directory in the current IBM SPSS Collaboration and Deployment Services Repository.

```
*Define and use a file handle for a specific file.
FILE HANDLE cust /NAME='SPSSCR://CustomerValue/data2002.sav'.
GET FILE='cust'.
```

- The handle *cust* is associated with the IBM SPSS Collaboration and Deployment Services Repository file */CustomerValue/data2002.sav*, and the GET command retrieves the latest version of this file from the current IBM SPSS Collaboration and Deployment Services Repository.

## ***Setting the Working Directory to a IBM SPSS Collaboration and Deployment Services Repository Location***

You can use the `CD` command to set the working directory to a directory in the currently connected IBM® SPSS® Collaboration and Deployment Services Repository.

```
CD 'SPSSCR://CustomerValue'.  
GET FILE='data2002.sav'.
```

- The working directory is set to the IBM SPSS Collaboration and Deployment Services Repository directory */CustomerValue*, and the `GET` command retrieves the latest version of the file *data2002.sav* from this directory in the current IBM SPSS Collaboration and Deployment Services Repository.



# ***TABLES and IGRAPH Command Syntax Converter***

If you have command syntax files that contain `TABLES` syntax that you want to convert to `CTABLES` syntax and/or `IGRAPH` syntax that you want to convert to `GGRAPH` syntax, a simple utility program is provided to help you get started with the conversion process. There are, however, significant functionality differences between `TABLES` and `CTABLES` and between `IGRAPH` and `GGRAPH`. It is likely that you will find that the utility program cannot convert some of your `TABLES` and `IGRAPH` syntax jobs or may generate `CTABLES` and `GGRAPH` syntax that produces tables and graphs that do not closely resemble the originals produced by the `TABLES` and `IGRAPH` commands. For most tables, you can edit the converted syntax to produce a table closely resembling the original.

The utility program is designed to:

- Create a new syntax file from an existing syntax file. The original syntax file is not altered.
- Convert only `TABLES` and `IGRAPH` commands in the syntax file. Other commands in the file are not altered.
- Retain the original `TABLES` and `IGRAPH` syntax in commented form.
- Identify the beginning and end of each conversion block with comments.
- Identify `TABLES` and `IGRAPH` syntax commands that could not be converted.
- Convert command syntax files that follow either interactive or production mode syntax rules.

This utility cannot convert commands that contain errors. The following other limitations also apply.

## ***TABLES Limitations***

The utility program may convert `TABLES` commands incorrectly under some circumstances, including `TABLES` commands that contain:

- Parenthesized variable names with the initial letters “sta” or “lab” in the `TABLES` subcommand if the variable is parenthesized by itself—for example, `var1 by (statvar) by (labvar)`. These will be interpreted as the `(STATISTICS)` and `(LABELS)` keywords.
- `SORT` subcommands that use the abbreviations `A` or `D` to indicate ascending or descending sort order. These will be interpreted as variable names.

The utility program cannot convert `TABLES` commands that contain:

- Syntax errors.



- OBSERVATION subcommands that refer to a range of variables using the TO keyword (for example, `var01 TO var05`).
- String literals broken into segments separated by plus signs (for example, `TITLE "My" + "Title"`).
- Macro calls that, in the absence of macro expansion, would be invalid TABLES syntax. Since the converter does not expand the macro calls, it treats them as if they were simply part of the standard TABLES syntax.

The utility program will not convert TABLES commands contained in macros. All macros are unaffected by the conversion process.

### ***IGRAPH Limitations***

IGRAPH changed significantly in release 16. Because of these changes, some subcommands and keywords in IGRAPH syntax created before that release may not be honored. See the IGRAPH section in the *Command Syntax Reference* for the complete revision history.

The conversion utility program may generate additional syntax that it stores in the `INLINETEMPLATE` keyword within the GGRAPH syntax. This keyword is created only by the conversion program. Its syntax is not intended to be user-editable.

### ***Using the Conversion Utility Program***

The conversion utility program, *SyntaxConverter.exe*, can be found in the installation directory. It is designed to run from a command prompt. The general form of the command is:

```
syntaxconverter.exe [path]/inputfilename.sps [path]/outputfilename.sps
```

You must run this command from the installation directory.

If any directory names contain spaces, enclose the entire path and filename in quotation marks, as in:

```
syntaxconverter.exe /myfiles/oldfile.sps "/new files/newfile.sps"
```

### ***Interactive versus Production Mode Command Syntax Rules***

The conversion utility program can convert command files that use interactive or production mode syntax rules.

**Interactive.** The interactive syntax rules are:

- Each command begins on a new line.
- Each command ends with a period (.).

**Production mode.** The Production Facility and commands in files accessed via the `INCLUDE` command in a different command file use production mode syntax rules:

- Each command must begin in the first column of a new line.
- Continuation lines must be indented at least one space.
- The period at the end of the command is optional.

If your command files use production mode syntax rules and don't contain periods at the end of each command, you need to include the command line switch `-b` (or `/b`) when you run *SyntaxConverter.exe*, as in:

```
syntaxconverter.exe -b /myfiles/oldfile.sps /myfiles/newfile.sps
```

### ***SyntaxConverter Script (Windows Only)***

On Windows, you can also run the syntax converter with the script *SyntaxConverter.wwd*, located in the *Samples* directory of the installation directory.

- ▶ From the menus choose:  
Utilities > Run Script...
- ▶ Navigate to the *Samples* directory and select *SyntaxConverter.wwd*.

This will open a simple dialog box where you can specify the names and locations of the old and new command syntax files.

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