Package 'BCA'

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Athletic

Intercollegiate Athletic Program Data Set

Description

The Athletic data set has 168 observations and 7 variables. The data come from a survey of stakeholders different (students, alumni, faculty, and athletic department employees) of a large US state university. The variables in the data set are conjoint analysis based relative importance weights for seven potential indicators of intercollegiate athletic program success.

Usage

data(Athletic)

Format

This data set contains the following variables:

- Win The importance of winning (won/loss record percentage) in determining the respondent's judgment of an intercollegiate athletic program's success.
- **Grad** The importance of student athlete graduation rates in determining the respondent's judgment of an intercollegiate athletic program's success.
- **Violat** The importance of NCAA rule violations in determining the respondent's judgment of an intercollegiate athletic program's success.
- Attnd The importance of home game attendance in determining the respondent's judgment of an intercollegiate athletic program's success.
- **Fem** The importance of gender equity (based on the ratio of female to male student athletes) in determining the respondent's judgment of an intercollegiate athletic program's success.
- **Teams** The importance of the number of different sports teams in determining the respondent's judgment of an intercollegiate athletic program's success.
- **Finan** The importance of the financial success of the program in determining the respondent's judgment of an intercollegiate athletic program's success.

bootCVD

Source

Wolfe, Richard A. and Daniel S. Putler (2002), "How Tight are the Ties that Bind Stakeholder Groups?, *Organizaton Science*, 13(January-February), 64-82.

bootCVD

Cluster Solution Diagnositics Using Bootstrap Replicates

Description

Provides a plot of both the Rand index and the Calinski-Harabas index for different numbers of clusters for a common underlying dataset using either the K-Means, K-Medians, or Neural Gas clusting algorithms based on a set of bootstrap replicates of the data.

Usage

```
bootCVD(x, k, nboot=100, nrep=1, method = c("kmn", "kmd", "neuralgas"),
    col1, col2, dsname)
bootCH(xdat, k_vals, clstr1, clstr2, cntrs1, cntrs2,
    method = c("kmn", "kmd", "neuralgas"))
bootPlot(fc, ch, col1="blue", col2="green")
```

Arguments

x	A numeric matrix of the data to be clustered.
k	An integer vector giving the set of clustering solutions to be examined.
nboot	The number of bootstrap replicates to use for the assessment.
nrep	The number of each set of initial cluster seeds on which to base a solution.
method	The clustering method, one of "kmn" (K-Means), "kmd" (K-Medians), and "neuralgas" (neural gas).
col1	The color to use for the plot of the Rand index values.
col2	The color to use for the plot of the Calinski-Harabas index values.
dsname	The name of the dataset being used (used only for output purposes.
xdat	A numeric matrix of the data to be clustered.
k_vals	An integer vector giving the set of clustering solutions to be examined.
clstr1	The cluster assignments from a bootFlexclust object for one side of the Rand index paired comparisons.
clstr2	The cluster assignments from a bootFlexclust object for the other side of the Rand index paired comparisons.
cntrs1	The cluster centers from a bootFlexclust object for one side of the bootFlexclust Rand index paired comparisons.
cntrs2	The cluster centers from a bootFlexclust object for the other side of the boot- Flexclust Rand index paired comparisons.
fc	A bootFlexclust object.
ch	A matrix of Calinski-Harabas index values from bootCH.

Details

The Rand index provides a measure of cluster stability, with relatively higher values indicating relatively more stable clusters, and the the Calinski-Harabas index gives a ratio of cluster seperation to cluster homogeneity, with higher values of the index being comparatively more preferred. The use of bootstrap replicates addresses both potential randomness in both the sample data and the clustering algorithms.

Value

The functions bootCVD and bootPlot return invisibly. Their benefit is the side effect plot produced and the printed summary of the index values. The function bootCH a matrix of Calinski-Harabas index values, the rows are the replicates, and each column corresponds to a particular number of clusters for a solution.

Author(s)

Dan Putler

References

S. Dolnicar, F. Leisch (2010), Evaluation of Structure and Reproducibility of Cluster Solution Using the Bootstrap. *Marketing Letters*, 21:1.

F. Leisch (2006), A Toolbox for K-Centroids Cluster Analysis. *Computational Statistics and Data Analysis*, 51:2.

See Also

bootFlexclust

bpCent

Two Dimensional Biplot of a Clustering Solution

Description

Plot a biplot of a clustering solution on the current graphics device.

Usage

```
bpCent(pc, clsAsgn, data.pts = TRUE, centroids = TRUE,
    choices = 1:2, scale = 1, pc.biplot=FALSE, var.axes = TRUE, col,
    cex = rep(par("cex"), 2), xlabs = NULL, ylabs = NULL, expand=1, xlim = NULL,
    ylim = NULL, arrow.len = 0.1, main = NULL, sub = NULL, xlab = NULL,
    ylab = NULL, ...)
```

bpCent

Arguments

рс	The prcomp object of the data used in clustering.
clsAsgn	A vector containing the cluster assignment for each record in the clustering data.
data.pts	If TRUE the point for each record is plotted.
centroids	If TRUE the centroid for each cluster is plotted.
choices	length 2 vector specifying the components to plot.
scale	The variables scaled by lambda ^ scale and the observations are scaled by lambda ^ (1-scale), where lambda are the eigen values of the principal components solution. scale should be between 0 and 1.
pc.biplot	If true, then lambda = 1 and the observations are are scaled up the $sqrt(n)$ and the variables scaled down by $sqrt(n)$. In this case the inner product between variables approximate covariances, and the distances between observations approximate Mahalanobis distance. Gabriel refers to this as a "principal component biplot".
var.axes	If TRUE the second set of points have arrows representing them as (unscaled) axes.
col	A vector of length 2 giving the colours for the first and second set of points respectively (and the corresponding axes). If a single colour is specified it will be used for both sets. If missing the default colour is looked for in the palette: if there it and the next colour as used, otherwise the first two colours of the paletter are used.
cex	The character expansion factor used for labelling the points. The labels can be of different sizes for the two sets by supplying a vector of length two.
xlabs	A vector of character strings to label the first set of points: the default is to use the row dimname of x, or $1:n$ is the dimname is NULL.
ylabs	A vector of character strings to label the second set of points: the default is to use the row dimname of y, or 1:n is the dimname is NULL.
expand	An expansion factor to apply when plotting the second set of points relative to the first. This can be used to tweak the scaling of the two sets to a physically comparable scale.
arrow.len	The length of the arrow heads on the axes plotted in var.axes is true. The arrow head can be suppressed by arrow.len = 0 .
xlim, ylim main, sub, xlał	Limits for the x and y axes in the units of the first set of variables. b, ylab, graphical parameters.

Details

This function is a reworking of biplot in order to allow the cluster centroids of a clustering solution to be displayed as well as the individual data points. If both data.pts and centroids are set to FALSE then only the variable directional vectors are displayed.

Graphical parameters can also be given to biplot.

Side Effects

a plot is produced on the current graphics device.

See Also

biplot

bpCent3d

Three Dimensional Biplot of a Clustering Solution

Description

Plot a biplot of a clustering solution on the current graphics device.

Usage

```
bpCent3d(pc, clsAsgn, data.pts = TRUE, centroids = TRUE,
    choices = 1:3, scale = 1, pc.biplot = FALSE, var.axes = TRUE,
    col.score = "black", col.cntrs = "blue", col.load = "red",
    xlabs = NULL, ylabs = NULL, xlim = NULL, ylim = NULL, zlim = NULL,
    xlab = NULL, ylab = NULL, dim.lab = NULL, fov = 60)
```

Arguments

рс	The prcomp object of the data used in clustering.
clsAsgn	A vector containing the cluster assignment for each record in the clustering data.
data.pts	If TRUE the point for each record is plotted.
centroids	If TRUE the centroid for each cluster is plotted.
choices	length 2 vector specifying the components to plot.
scale	The variables scaled by lambda ^ scale and the observations are scaled by lambda ^ (1-scale), where lambda are the eigen values of the principal components solution. scale should be between 0 and 1.
pc.biplot	If TRUE, then lambda = 1 and the observations are scaled up the $sqrt(n)$ and the variables scaled down by $sqrt(n)$. In this case the inner product between variables approximate covariances, and the distances between observations approximate Mahalanobis distance. Gabriel refers to this as a "principal component biplot".
var.axes	If TRUE the second set of points have arrows representing them as (unscaled) axes.
col.score, col.	cntrs, col.load
	A vector of length 2 giving the colours for the first and second set of points respectively (and the corresponding axes). If a single colour is specified it will be used for both sets. If missing the default colour is looked for in the palette: if there it and the next colour as used, otherwise the first two colours of the paletter are used.

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xlabs	A vector of character strings to label the first set of points: the default is to use the row dimname of x , or $1:n$ is the dimname is NULL.
ylabs	A vector of character strings to label the second set of points: the default is to use the row dimname of y, or 1:n is the dimname is NULL.
xlim, ylim	, zlim
	Limits for the x, y, z axes in the units of the first set of variables.
xlab, ylab	, dim.lab, fov
	Graphical parameters for rgl.

Details

This function is a reworking of biplot in order to allow the cluster centroids of a clustering solution to be displayed as well as the individual data points, and extend it to three dimensions. The use of arrow heads is discouraged since their creation is very resource intensive. If both data.pts and centroids are set to FALSE then only the variable directional vectors are displayed.

Side Effects

A plot is produced on the current graphics device.

See Also

biplot

CCS

Charitable Giving Dataset

Description

The CCS data set has 1600 observations and 20 variables, and is from the data is from the Britsh Columbia and Yukon region of a nationwide Canadian charitable society. The data examine whether a donor joined the Monthly Giver Program, in which he or she opts to make a pre-arranged monthly donation to the charity by credit card. The actual monthly givers have been highly over-sampled. In the charity's database they represent one percent of donors, but they comprise 50 percent of this sample. This level of oversampling is common in data mining applications.

Usage

data(CCS)

Format

This data set contains the following variables:

MonthGive A factor indicating whether the donor joined the monthly giver program in the last two annual campaigns with levels:

No

Yes

Region A factor indicating where in British Columbia or the Yukon the donor resides with levels:

- R1 (Vancouver Island)
- R2 (Greater Vancouver)
- R3 (The Fraser Valley)
- R4 (The North Coast of BC)
- R5 (The Central and Southern Interior of BC)
- R6 (The Northern Interior of BC and the Yukon)

YearsGive The number of years the individual has given to the Canadian Cancer Society.

AveDonAmt The average dollar amount of past donations by the donor.

LastDonAmt The dollar amount of the last donation by the donor.

- **DonPerYear** The rate of donating to the Canadian Cancer Society measured as the number of donations per year by the donor.
- **NewDonor** A factor that indicates whether the individual is a first time donor with levels. No
 - Yes
- **Age20t29** The percentage of people between age 20 and 29 residing in the Enumeration Area in which the donor resides.
- **Age20t39** The percentage of people between age 20 and 39 residing in the Enumeration Area in which the donor resides.
- **Age60pls** The percentage of people 60 years of age or older residing in the Enumeration Area in which the donor resides.
- **Age70pls** The percentage of people 70 years of age or older residing in the Enumeration Area in which the donor resides.
- **Age80pls** The percentage of people 80 years of age or older residing in the Enumeration Area in which the donor resides.
- AdultAge The average age of adult residents in the Enumeration Area in which the donor resides.
- **SomeUnivP** The percentage of adults in the postal code in which the donor resides that have an educational attainment of some college or university attendance.
- **FinUnivP** The percentage of adults in the Enumeration Area in which the donor resides that have a university degree.
- hhlt2mem The percentage of households in the Enumeration Area in which the donor resides that have only one or two members present.
- **hh1mem** The percentage of households in the Enumeration Area in which the donor resides that have only a single member present.
- AveIncEA The average pre-tax income of households in the Enumeration Area in which the donor resides.
- **DwelValEA** The average dwelling value in the Enumeration Area in which the donor resides.
- **EngPrmLang** The percentage of households in the Enumeration Area in which the donor resides that have have English as their primary language.

Source

An anonymous charity.

create.samples

Description

Provides a character vector with possible values of "Estimation", "Validation" and "Holdout" that can then be used to assign observations of a data frame to estimation, validation, or (optionally) holdsout samples using the subset option of a variety of functions.

Usage

create.samples(x, est=0.34, val=0.33, rand.seed=1)

Arguments

x	A data frame.
est	The percentage of the total sample to allocate to the estimation sample. The value of est should range from zero to one
val	The percentage of the total sample to allocate to the validation sample. The value of val should range from zero to one
rand.seed	A parameter passed to set.seed for to specify the seed of the random number generator.

Details

The values of est and val should sum to a value between zero and one. If greater than one, an error is returned. If less than one, the remaining percentage of the sample is allocated to the holdout sample.

Value

A character vector with possible values of "Estimation", "Validation" and (optionally) "Holdout". The length of the vector equals the number of rows in the original data frame.

Author(s)

Dan Putler

See Also

set.seed

Examples

```
data(CCS)
# Create a new set of samples with 40 percent in each of the estimation and
# validation samples, and 20 percent in the holdout sample.
CCS$Sample <- create.samples(CCS, est=0.4, val=0.4)</pre>
```

Eggs

Retail Sales of Eggs

Description

The Eggs data set has 105 observations and 10 variables. The data contains information on weekly sales of eggs in Southern California over a two year period.

Usage

data(Eggs)

Format

This data set contains the following variables:

Week The observation week (1 to 105). This variable can be used as a time trend.

Month A factor that gives the name of the month in which the observation occured.

First.Week A factor indicating whether the observation fell on the first week of the month with levels:

No

Yes

Easter A factor that indicates whether the observation fell the week prior to the week containing Easter Sunday, the week containing Easter Sunday, the week containing Easter Sunday, or a non-Easter week with levels:

Non Easter Pre Easter

Easter

Post Easter

Cases Retail sales of eggs in cases.

Egg.Pr Average retail egg price in cents per dozen.

Beef.Pr Average retail price of 7-bone beef roast in cents per pound.

Pork.Pr Average retail price of strip bacon in cents per pound.

Chicken.Pr Average retail price of whole frying chicken in cents per pound.

Cereal.Pr Average retail price of Cheerios breakfast cereal in cents per pound.

Source

Putler (1992)

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jack.jill

Description

The jack.jill data set has 557 observations and 8 variables. The data contains information on children's apparel spending and household level demographic and socioeconomic information for a sample of households residing in Alberta and British Columbia.

Usage

data(jack.jill)

Format

This data set contains the following variables:

SPENDING Dollars spent on children's apparel over a one-year period.

CHILDREN Pre-tax income given as a factor with levels:

INCOME Pre-tax income given as a factor with levels:

\$0k-\$20k \$20k-\$30k \$30k-\$40k \$40k-\$50k \$50k-\$60k \$60k-\$75k \$75k-\$100k \$100k+

EMPOLYMENT The employment status of the female head of household with levels:

No female head Unemployed Part-time Full-time

AGE Age of the female head of household given as a factor with levels:

- No female head
- 29 and under
- 30 to 39
- 40 to 49
- 50 to 59
- 60 and over

- **EDUCATION** The educational attainment of the female head of household given as a factor with levels:
 - No female head Not stated Elementary or less Some or completed secondary Some post-secondary Post-secondary diploma University degree

OCCUPATION The occupation group of the female head of household with levels:

```
No female head
Blue collar
Pink collar
White collar
Other
Non-working or retired
```

BIRTHCNTRY The birth country of the female head of household with levels:

No female head Canada US, N&W Europe S&E Europe Asia and Oceania Other (Caribbean, Middle East, and Africa)

Source

Statistics Canada

lift.chart

Lift Charts to Compare Binary Predictive Models

Description

Provides either a total cumulative response or incremental response rate lift chart for the purposes of comparing the predictive capability of different binary predictive models.

Usage

```
lift.chart(modelList, data, targLevel, trueResp, type = "cumulative", sub = "")
```

Arguments

modelList A character vector containing the names of the different models to be compared. The selected models must have the same y variable that must be a binary factor, and have been estimated using the same data set.

lift.chart

data	The dataframe that constitues the comparison sample. If this dataframe is not the same as the dataframe used to estimated models, the dataframe must contain all the variables used in the models to be compared.
targLevel	The label for the level of the binary factor of interest. For example, in a database marketing application, this level could be "Yes" for a variable that takes on the values "Yes" and "No" to indicate if a customer responded favorably to a promotion offer.
trueResp	The true rate of the target level for the master database the estimation and com- parison dataframes were originally drawn from.
type	A character string that must either have the value of "cummulative" (to produce a total cummaltive response chart) or "incremental" (to produce an incremental response rate chart).
sub	A sub-title for the plot, typically to identify the sample used.

Details

Lift charts are a commonly used tool in business data mining applications. They are used to assess how well a model is able to predict a desirable (from an organization's point-of-view) response on the part of a customer compared to alternative estimated models and a benchmark model of approaching customers randomly. The total cummulative response chart shows the percentage of the total response the organization would receive from only contacting a given percentage (grouped by deciles) of its entire customer base. This chart is best for selecting between alternative models, and in predicting the revenues the organization will receive by contacting a given percentage of their customers that the model predicts are most likely to favorably respond. The incremental response rate chart provides the response rate among each of ten decile groups of the organization's customers, with the decile groups ordered by their estimated likelihood of a favorable response.

Value

The function returns invisibly. Its benefit is the side effect plot produced.

Author(s)

Dan Putler

Examples

```
library(rpart)
layout(matrix(c(1,2), 2, 1))
data(CCS)
CCS$Sample <- create.samples(CCS, est=0.4, val=0.4)
CCSEst <- CCS[CCS$Sample == "Estimation",]
CCS.glm <- glm(MonthGive ~ DonPerYear + LastDonAmt + Region + YearsGive,
   family=binomial(logit), data=CCSEst)
library(rpart)
CCS.rpart <- rpart(MonthGive ~ DonPerYear + LastDonAmt + Region + YearsGive,
   data=CCSEst, cp=0.0074)
CCSVal <- CCS[CCS$Sample == "Validation",]
lift.chart(c("CCS.glm", "CCS.rpart"), data=CCSVal, targLevel="Yes",</pre>
```

```
trueResp=0.01, type="cumulative", sub="Validation")
lift.chart(c("CCS.glm", "CCS.rpart"), data=CCSVal, targLevel="Yes",
    trueResp=0.01, type="incremental", sub="Validation")
```

relabel.factor Relabel Factor Levels

Description

Relabel the levels of factors to provide more descriptive names and reduce the number of factor levels.

Usage

```
relabel.factor(x, new.labels, old.labels=levels(x))
```

Arguments

x	A factor.
new.labels	The new factor level labels.
old.labels	The old factor level labels.

Details

The number of new factor labels/levels must be less than the number of labels/levels than the original factor.

Value

A factor whose length is equal to the old factor.

Author(s)

Dan Putler

Description

A minor modification of the car package's scatterplot function that makes enhanced scatterplots, with boxplots in the margins, a lowess smooth, smoothed conditional spread, outlier identification, and a regression line; sp is an abbreviation for scatterplot.

Usage

```
scatterplotBCA(x, ...)
## S3 method for class 'formula'
scatterplotBCA(x, data, subset, xlab, ylab, legend.title, legend.coords,
labels, ...)
## Default S3 method:
scatterplotBCA(x, y, smooth = TRUE, spread = !by.groups,
         span = 0.5, loess.threshold = 2, reg.line = lm,
         boxplots = if (by.groups) "" else "xy", xlab = deparse(substitute(x)),
         ylab = deparse(substitute(y)), las = par("las"), lwd = 2,
         lwd.smooth = lwd, lwd.spread = lwd, lty = 1, lty.smooth = lty,
         lty.spread = 2, labels, id.method = "mahal",
         id.n = if(id.method[1] == "identify") length(x) else 0, id.cex = 1,
         id.col = palette()[1], log = "", jitter = list(), xlim = NULL,
         ylim = NULL, cex = par("cex"), cex.axis = par("cex.axis"),
         cex.lab = par("cex.lab"), cex.main = par("cex.main"),
         cex.sub = par("cex.sub"), groups, by.groups = !missing(groups),
         legend.title = deparse(substitute(groups)), legend.coords,
         ellipse = FALSE, levels = c(0.5, 0.95), robust = TRUE, col = if
         (n.groups == 1) palette()[c(2, 1, 3)] else rep(palette(),
         length = n.groups), pch = 1:n.groups, legend.plot = !missing(groups),
         reset.par = TRUE, grid = TRUE, ...)
```

spBCA(...)

Arguments

х	vector of horizontal coordinates, or a "model" formula, of the form $y \sim x$ or (to plot by groups) $y \sim x \mid z$, where z evaluates to a factor or other variable dividing the data into groups. If x is a factor, then parallel boxplots are produced using the Boxplot function.
У	vector of vertical coordinates.
data	data frame within which to evaluate the formula.
subset	expression defining a subset of observations.

smooth	if TRUE (the default) a loess nonparametric regression line is drawn on the plot.
spread	if TRUE (the default when there are no groups), a smoother is applied to the root-mean-square positive and negative residuals from the loess line to display conditional spread and asymmetry.
span	span for the loess smoother.
loess.threshol	
	suppress the loess smoother if there are fewer than loess.threshold unique values (default, 5) of y.
reg.line	function to draw a regression line on the plot or FALSE not to plot a regression line.
boxplots	if "x" a boxplot for x is drawn below the plot; if "y" a boxplot for y is drawn to the left of the plot; if "xy" both boxplots are drawn; set to "" or FALSE to suppress both boxplots.
xlab	label for horizontal axis.
ylab	label for vertical axis.
las	if 0, ticks labels are drawn parallel to the axis; set to 1 for horizontal labels (see par).
lwd	width of linear-regression lines (default 1).
lwd.smooth	width for smooth regression lines (default is the same as 1wd).
lwd.spread	width for lines showing spread (default is the same as 1wd).
lty	type of linear-regression lines (default 1, solid line).
lty.smooth	type of smooth regression lines (default is the same as lty).
lty.spread	width for lines showing spread (default is 2, broken line).
id.method,id.n	
	Arguments for the labelling of points. The default is id.n=0 for labeling no points. See showLabels for details of these arguments. If the plot uses different colors for groups, then the id.col argument is ignored and label colors are determined by the col argument.
labels	a vector of point labels; if absent, the function tries to determine reasonable labels, and, failing that, will use observation numbers.
log	same as the log argument to plot, to produce log axes.
jitter	a list with elements x or y or both, specifying jitter factors for the horizontal and vertical coordinates of the points in the scatterplot. The jitter function is used to randomly perturb the points; specifying a factor of 1 produces the default jitter. Fitted lines are unaffected by the jitter.
xlim	the x limits (min, max) of the plot; if NULL, determined from the data.
ylim	the y limits (min, max) of the plot; if NULL, determined from the data.
groups	a factor or other variable dividing the data into groups; groups are plotted with different colors and plotting characters.
by.groups	if TRUE, regression lines are fit by groups.
legend.title	title for legend box; defaults to the name of the groups variable.

legend.coords	coordinates for placing legend; an be a list with components x and y to specify the coordinates of the upper-left-hand corner of the legend; or a quoted keyword, such as "topleft", recognized by legend.
ellipse	if TRUE data-concentration ellipses are plotted.
levels	level or levels at which concentration ellipses are plotted; the default is $c(.5, .95)$.
robust	if TRUE (the default) use the cov.trob function in the MASS package to calculate the center and covariance matrix for the data ellipses.
col	colors for lines and points; the default is taken from the color palette, with palette()[2] for nonparametric regression lines and palette()[1] for linear regression line and points if there are no groups, and successive colors for the groups if there are groups.
pch	plotting characters for points; default is the plotting characters in order (see par).
cex, cex.axis,	cex.lab, cex.main, cex.sub set sizes of various graphical elements; (see par).
legend.plot	if TRUE then a legend for the groups is plotted in the upper margin.
reset.par	if TRUE then plotting parameters are reset to their previous values when scatterplot exits; if FALSE then the mar and mfcol parameters are altered for the current plot- ting device. Set to FALSE if you want to add graphical elements (such as lines) to the plot.
	other arguments passed down and to plot.
grid	If TRUE, the default, a light-gray background grid is put on the graph

Value

If points are identified, their labels are returned; otherwise NULL is returned invisibly.

Author(s)

John Fox with modifications made by Dan Putler

See Also

scatterplot

scatterplotMatrixBCA Scatterplot Matrices

Description

A minor modification of the car package's scatterplotMatrix function that makes enhanced scatterplot matrices with univariate displays down the diagonal; spmBCA is an abbreviation for scatterplotMatrixBCA. This function just sets up a call to pairs with custom panel functions.

Usage

```
scatterplotMatrixBCA(x, ...)
## S3 method for class 'formula'
scatterplotMatrixBCA(x, data=NULL, subset, labels, ...)
## Default S3 method:
scatterplotMatrixBCA(x, var.labels = colnames(x), diagonal = c("density",
                 "boxplot", "histogram", "oned", "qqplot", "none"),
                 adjust = 1, nclass, plot.points = TRUE, smooth = TRUE,
                 spread = smooth && !by.groups, span = 0.5,
                 loess.threshold = 2, reg.line = lm, transform = FALSE,
                 family = c("bcPower", "yjPower"), ellipse = FALSE,
                 levels = c(0.5, 0.95), robust = TRUE, groups = NULL,
                 by.groups = FALSE, labels, id.method = "mahal", id.n =
                 0, id.cex = 1, id.col = palette()[1], col = if
                 (n.groups == 1) palette()[c(2, 1, 3)] else rep(palette(),
                 length = n.groups), pch = 1:n.groups, lwd = 2,
                 lwd.smooth = lwd, lwd.spread = lwd, lty = 1,
                 lty.smooth = lty, lty.spread = 2, cex = par("cex"),
                 cex.axis = par("cex.axis"), cex.labels = NULL,
                 cex.main = par("cex.main"), legend.plot =
                 length(levels(groups)) > 1, row1attop = TRUE, ...)
```

spmBCA(x, ...)

Arguments

х	a data matrix, numeric data frame, or a one-sided "model" formula, of the form $\sim x1 + x2 + \ldots + xk$ or $\sim x1 + x2 + \ldots + xk \mid z$ where z evaluates to a factor or other variable to divide the data into groups.	
data	for scatterplotMatrix.formula, a data frame within which to evaluate the formula.	
subset	expression defining a subset of observations.	
labels,id.method,id.n,id.cex,id.col		
	Arguments for the labelling of points. The default is id.n=0 for labeling no points. See showLabels for details of these arguments. If the plot uses different colors for groups, then the id.col argument is ignored and label colors are determined by the col argument.	
var.labels	variable labels (for the diagonal of the plot).	
diagonal	contents of the diagonal panels of the plot.	
adjust	relative bandwidth for density estimate, passed to density function.	
nclass	number of bins for histogram, passed to hist function.	
plot.points	if TRUE the points are plotted in each off-diagonal panel.	

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smooth	if TRUE a loess smooth is plotted in each off-diagonal panel.
spread	if TRUE (the default when not smoothing by groups), a smoother is applied to the root-mean-square positive and negative residuals from the loess line to display conditional spread and asymmetry.
span	span for loess smoother.
loess.threshol	
	suppress the loess smoother if there are fewer than loess.threshold unique values (default, 5) of the variable on the vertical axis.
reg.line	if not FALSE a line is plotted using the function given by this argument; e.g., using rlm in package MASS plots a robust-regression line.
transform	if TRUE, multivariate normalizing power transformations are computed with powerTransform, rounding the estimated powers to 'nice' values for plotting; if a vector of powers, one for each variable, these are applied prior to plotting. If there are groups and by.groups is TRUE, then the transformations are estimated <i>conditional</i> on the groups factor.
family	family of transformations to estimate: "bcPower" for the Box-Cox family or "yjPower" for the Yeo-Johnson family (see powerTransform).
ellipse	if TRUE data-concentration ellipses are plotted in the off-diagonal panels.
levels	levels or levels at which concentration ellipses are plotted; the default is c(.5, .9).
robust	if TRUE use the cov.trob function in the MASS package to calculate the center and covariance matrix for the data ellipses.
groups	a factor or other variable dividing the data into groups; groups are plotted with different colors and plotting characters.
by.groups	if TRUE, regression lines are fit by groups.
pch	plotting characters for points; default is the plotting characters in order (see par).
col	colors for lines and points; the default is taken from the color palette, with palette()[2] for nonparametric regression lines and palette()[1] for linear regression lines and points if there are no groups, and successive colors for the groups if there are groups.
lwd	width of linear-regression lines (default 1).
lwd.smooth	width for smooth regression lines (default is the same as lwd).
lwd.spread	width for lines showing spread (default is the same as lwd).
lty	type of linear-regression lines (default 1, solid line).
lty.smooth	type of smooth regression lines (default is the same as lty).
lty.spread	width for lines showing spread (default is 2, broken line).
cex, cex.axis,	cex.labels, cex.main set sizes of various graphical elements (see par).
legend.plot	if TRUE then a legend for the groups is plotted in the first diagonal cell.
row1attop	If TRUE (the default) the first row is at the top, as in a matrix, as opposed to at the bottom, as in graph (argument suggested by Richard Heiberger).
	arguments to pass down.

Value

NULL. This function is used for its side effect: producing a plot.

Author(s)

John Fox with minor modifications by Dan Putler

See Also

scatterplotMatrix

score

Score a Database based on a Predictive Model

Description

Provides either an integer vector that contains the "desirability" rank of a case in a data set, the fitted probability of a desired response, or the fitted probability adjusted for the true response rate based on the fitted values of a predictive model.

Usage

```
rankScore(model, data, targLevel)
rawProbScore(model, data, targLevel)
adjProbScore(model, data, targLevel, trueResp)
```

Arguments

model	A character string containing the name of the model to use to score the database.
data	A data frame of the database to be scored. All the predictor variables of the model need to be amoung the variables of the data frame.
targLevel	The "desired" level of the y variable factor as a character string.
trueResp	The true "desired" response rate for the overall population of interest.

Details

Only binomial glm, binary rpart, and binary nnet models can be used as the basis of scoring a database.

Value

An integer vector that indicates the rank order desirability (a value of 1 means most desirable) of the correponding case of the database being scored or a probability measure bounded between zero and one.

Author(s)

Dan Putler

SD.clv

Description

Provides a wrapper to several function calls in the clv package needed to construct the SD index value for a clustering solution. The number of clusters that has the lowest value of the SD index represents the "best" solution under the criteria used to construct the SD index.

Usage

SD.clv(x, clus, alpha)

Arguments

Х	A numeric matrix of data, or an object that can be coerced to such a matrix (such as a numeric vector or a dataframe with all numeric columns) used to construct the clustering solution.
clus	The cluster to which each row of x was assigned.
alpha	A weight to be placed on the average scattering of the clustering solution.

Details

The SD index corresponds to the weighted sum of the average "scattering" of points within clusters and the inverse of the total seperation between clusters. The average scattering measure is based on the average sum of the squared differences between a clusters centroid all the points in a cluster, while total seperation is measured by the sum of the squared distance between cluster centroids. A solution with a low average scattering and a low value of the inverse total seperation is considered to be better than a solution with higher levels of these two measures.

Value

A scalar SD index value for the clustering solution.

Author(s)

Dan Putler

References

M. Haldiki, Y. Batistakis, M. Vazirgiannis (2001), On Clustering Validation Techniques, *Journal of Intelligent Information Systems*, 17:2/3.

See Also

clv.SD

Examples

```
data(iris)
iris.data <- iris[,1:4]
irisC3 <- kmeans(iris.data, centers=3, nstart=10)
SD.clv(iris.data, clus=irisC3$cluster, alpha=0.1)</pre>
```

```
SDIndex
```

A Plot of SD Index Values for K-Means Clustering Solutions

Description

Provides a plot of SD cluster validation index values for different numbers of k-means clusters for a common underlying dataset. The number of clusters that has the lowest value of the SD index represents the "best" solution under the criteria used to construct the SD index.

Usage

SDIndex(x, minClust, maxClust, iter.max=10, num.seeds=10)

Arguments

х	A numeric matrix of data, or an object that can be coerced to such a matrix (such as a numeric vector or a dataframe with all numeric columns).
minClust	The minimum number of clusters to be considered for a solution.
maxClust	The maximum number of clusters to be considered for a solution.
iter.max	The maximum number of iterations allowed for a solution.
num.seeds	The number of different starting random seeds to use for a solution with a given number of clusters.

Details

The SD index corresponds to the weighted sum of the average "scattering" of points within clusters and the inverse of the total seperation between clusters. The average scattering measure is based on the average sum of the squared differences between a clusters centroid all the points in a cluster, while total seperation is measured by the sum of the squared distance between cluster centroids. A solution with a low average scattering and a low value of the inverse total seperation is considered to be better than a solution with higher levels of these two measures.

Value

The function returns invisibly. Its benefit is the side effect plot produced.

Author(s)

Dan Putler

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variable.summary

References

M. Haldiki, Y. Batistakis, M. Vazirgiannis (2001), On Clustering Validation Techniques, *Journal of Intelligent Information Systems*, 17:2/3.

See Also

KMeans, SD.clv

Examples

```
data(iris)
iris.data <- iris[,1:4]
SDIndex(iris.data, minClust=2, maxClust=6, iter.max=10, num.seeds=10)</pre>
```

variable.summary Basic summary information of the variables of a data frame

Description

The function returns a data frame where, the row names correspond to the variable names, and a set of columns with summary information for each variable. Its purpose is to allow the user to quickly scan the data frame for potentially problematic variables.

Usage

```
variable.summary(dframe)
```

Arguments

dframe A data frame.

Value

The returned data frame contains the variables Class (numeric, integer, factor, or character), missing values), Levels (the levels of a factor variable, or NA for non-factor variables), Min.Level.Size (the number of cases for the smallest level of a factor, or NA for a non-factor), Mean (the mean of non-missing cases for a numeric or integer variable, or NA for factor and character variables), and SD (the standard deviation of non-missing cases for a numeric or integer variable, or NA for factor and character variables).

Author(s)

Dan Putler

Examples

data(CCS)
variable.summary(CCS)

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