

# Package: clustrd (via r-universe)

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**Type** Package

**Title** Methods for Joint Dimension Reduction and Clustering

**Description** A class of methods that combine dimension reduction and clustering of continuous, categorical or mixed-type data (Markos, Iodice D'Enza and van de Velden 2019; <[DOI:10.18637/jss.v091.i10](https://doi.org/10.18637/jss.v091.i10)>). For continuous data, the package contains implementations of factorial K-means (Vichi and Kiers 2001; <[DOI:10.1016/S0167-9473\(00\)00064-5](https://doi.org/10.1016/S0167-9473(00)00064-5)>) and reduced K-means (De Soete and Carroll 1994; <[DOI:10.1007/978-3-642-51175-2\\_24](https://doi.org/10.1007/978-3-642-51175-2_24)>); both methods that combine principal component analysis with K-means clustering. For categorical data, the package provides MCA K-means (Hwang, Dillon and Takane 2006; <[DOI:10.1007/s11336-004-1173-x](https://doi.org/10.1007/s11336-004-1173-x)>), i-FCB (Iodice D'Enza and Palumbo 2013, <[DOI:10.1007/s00180-012-0329-x](https://doi.org/10.1007/s00180-012-0329-x)>) and Cluster Correspondence Analysis (van de Velden, Iodice D'Enza and Palumbo 2017; <[DOI:10.1007/s11336-016-9514-0](https://doi.org/10.1007/s11336-016-9514-0)>), which combine multiple correspondence analysis with K-means. For mixed-type data, it provides mixed Reduced K-means and mixed Factorial K-means (van de Velden, Iodice D'Enza and Markos 2019; <[DOI:10.1002/wics.1456](https://doi.org/10.1002/wics.1456)>), which combine PCA for mixed-type data with K-means.

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clustrd-package	<i>Methods for Joint Dimension Reduction and Clustering</i>
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## Description

A class of methods that combine dimension reduction and clustering of continuous, categorical or mixed-type data (Markos, Iodice D’Enza and van de Velden 2019; <DOI:10.18637/jss.v091.i10>). For continuous data, the package contains implementations of factorial K-means (Vichi and Kiers 2001; <DOI:10.1016/S0167-9473(00)00064-5>) and reduced K-means (De Soete and Carroll 1994; <DOI:10.1007/978-3-642-51175-2\_24>); both methods that combine principal component analysis with K-means clustering. For categorical data, the package provides MCA K-means (Hwang, Dillon and Takane 2006; <DOI:10.1007/s11336-004-1173-x>), i-FCB (Iodice D’Enza and Palumbo 2013, <DOI:10.1007/s00180-012-0329-x>) and Cluster Correspondence Analysis (van de Velden, Iodice D’Enza and Palumbo 2017; <DOI:10.1007/s11336-016-9514-0>), which combine multiple correspondence analysis with K-means. For mixed-type data, it provides mixed Reduced K-means and mixed Factorial K-means (van de Velden, Iodice D’Enza and Markos 2019; <DOI:10.1002/wics.1456>), which combine PCA for mixed-type data with K-means.

**Details**

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**Author(s)**

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**References**

Markos, A., Iodice D'Enza, A., & van de Velden, M. (2019). Beyond Tandem Analysis: Joint Dimension Reduction and Clustering in R. *Journal of Statistical Software*, 91(10), 1–24. doi:10.18637/jss.v091.i10.

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bribery

*Bribery cases in Russia*

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**Description**

The data set refers to a collection of 55 articles on bribery cases from central Russian newspapers 1999-2000 (Mirkin, 2005). The variables reflect the following five-fold structure of bribery situations: two interacting sides - the office and the client, their interaction, the corrupt service rendered, and the environment in which it all occurs. These structural aspects can be characterized by 11 variables that have been manually recovered from the newspaper articles.

**Usage**

```
data("bribery")
```

**Format**

A data frame with 55 observations on 11 categorical variables.

Of Type of Office

Cl Level of Client

Serv Type of service: obstruction of justice, favours, cover-up, change of category, extortion of money for rendering free services

Occ Frequency of occurrence

Init Who initiated the bribery act

Brib Bribe Level in \$

Typ Type of corruption  
 Net Corruption network  
 Con Condition of corruption  
 Bran Branch at which the corrupt service occurred  
 Pun Punishment

## References

Mirkin, B. (2005). *Clustering for data mining: a data recovery approach*. Chapman and Hall/CRC.

## Examples

```
data(bribery)
```

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clusmca	<i>Joint dimension reduction and clustering of categorical data.</i>
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## Description

This function implements MCA K-means (Hwang, Dillon and Takane, 2006), i-FCB (Iodice D'Enza and Palumbo, 2013) and Cluster Correspondence Analysis (van de Velden, Iodice D'Enza and Palumbo, 2017). The methods combine variants of Correspondence Analysis for dimension reduction with K-means for clustering.

## Usage

```
clusmca(data, nclus, ndim, method=c("clusCA","iFCB","MCAk"),
  alphak = .5, nstart = 100, smartStart = NULL, gamma = TRUE,
  inboot = FALSE, seed = NULL)
```

```
## S3 method for class 'clusmca'
print(x, ...)
```

```
## S3 method for class 'clusmca'
summary(object, ...)
```

```
## S3 method for class 'clusmca'
fitted(object, mth = c("centers", "classes"), ...)
```

## Arguments

data	Dataset with categorical variables
nclus	Number of clusters (nclus = 1 returns the MCA solution; see Details)
ndim	Dimensionality of the solution

method	Specifies the method. Options are MCAk for MCA K-means, iFCB for Iterative Factorial Clustering of Binary variables and clusCA for Cluster Correspondence Analysis (default = "clusCA")
alphak	Non-negative scalar to adjust for the relative importance of MCA (alphak = 1) and K-means (alphak = 0) in the solution (default = .5). Works only in combination with method = "MCAk"
nstart	Number of random starts (default = 100)
smartStart	If NULL then a random cluster membership vector is generated. Alternatively, a cluster membership vector can be provided as a starting solution
gamma	Scaling parameter that leads to similar spread in the object and variable scores (default = TRUE)
seed	An integer that is used as argument by set.seed() for offsetting the random number generator when smartStart = NULL. The default value is NULL.
inboot	Used internally in the bootstrap functions to perform bootstrapping on the indicator matrix.
x	For the print method, a class of clusmca
object	For the summary method, a class of clusmca
mth	For the fitted method, a character string that specifies the type of fitted value to return: "centers" for the observations center vector, or "class" for the observations cluster membership value
...	Not used

### Details

For the K-means part, the algorithm of Hartigan-Wong is used by default.

The hidden print and summary methods print out some key components of an object of class clusmca.

The hidden fitted method returns cluster fitted values. If method is "classes", this is a vector of cluster membership (the cluster component of the "clusmca" object). If method is "centers", this is a matrix where each row is the cluster center for the observation. The rownames of the matrix are the cluster membership values.

When nclus = 1 the function returns the MCA solution with objects in principal and variables in standard coordinates. plot(object) shows the corresponding asymmetric biplot.

### Value

obscoord	Object scores
attcoord	Attribute scores
centroid	Cluster centroids
cluster	Cluster membership
criterion	Optimal value of the objective criterion
size	The number of objects in each cluster
nstart	A copy of nstart in the return object
odata	A copy of data in the return object

## References

- Hwang, H., Dillon, W. R., and Takane, Y. (2006). An extension of multiple correspondence analysis for identifying heterogenous subgroups of respondents. *Psychometrika*, 71, 161-171.
- Iodice D'Enza, A., and Palumbo, F. (2013). Iterative factor clustering of binary data. *Computational Statistics*, 28(2), 789-807.
- van de Velden M., Iodice D' Enza, A., and Palumbo, F. (2017). Cluster correspondence analysis. *Psychometrika*, 82(1), 158-185.

## See Also

[cluspca](#), [cluspcamix](#), [tuneclus](#)

## Examples

```
data(cmc)
# Preprocessing: values of wife's age and number of children were categorized
# into three groups based on quartiles
cmc$W_AGE = ordered(cut(cmc$W_AGE, c(16,26,39,49), include.lowest = TRUE))
levels(cmc$W_AGE) = c("16-26", "27-39", "40-49")
cmc$NCHILD = ordered(cut(cmc$NCHILD, c(0,1,4,17), right = FALSE))
levels(cmc$NCHILD) = c("0", "1-4", "5 and above")

#Cluster Correspondence Analysis solution with 3 clusters in 2 dimensions
#after 10 random starts
outclusCA = clusmca(cmc, 3, 2, method = "clusCA", nstart = 10, seed = 1234)
outclusCA
#Scatterplot (dimensions 1 and 2)
plot(outclusCA)

#MCA K-means solution with 3 clusters in 2 dimensions after 10 random starts
outMCAk = clusmca(cmc, 3, 2, method = "MCAk", nstart = 10, seed = 1234)
outMCAk
#Scatterplot (dimensions 1 and 2)
plot(outMCAk)

#nclus = 1 just gives the MCA solution
#outMCA = clusmca(cmc, 1, 2)
#outMCA
#Scatterplot (dimensions 1 and 2)
#asymmetric biplot with scaling gamma = TRUE
#plot(outMCA)
```

---

cluspca

*Joint dimension reduction and clustering of continuous data.*

---

## Description

This function implements Factorial K-means (Vichi and Kiers, 2001) and Reduced K-means (De Soete and Carroll, 1994), as well as a compromise version of these two methods. The methods combine Principal Component Analysis for dimension reduction with K-means for clustering.

**Usage**

```
cluspca(data, nclus, ndim, alpha = NULL, method = c("RKM", "FKM"),
center = TRUE, scale = TRUE, rotation = "none", nstart = 100,
smartStart = NULL, seed = NULL)

## S3 method for class 'cluspca'
print(x, ...)

## S3 method for class 'cluspca'
summary(object, ...)

## S3 method for class 'cluspca'
fitted(object, mth = c("centers", "classes"), ...)
```

**Arguments**

<code>data</code>	Dataset with metric variables
<code>nclus</code>	Number of clusters ( <code>nclus = 1</code> returns the PCA solution)
<code>ndim</code>	Dimensionality of the solution
<code>method</code>	Specifies the method. Options are RKM for reduced K-means and FKM for factorial K-means (default = "RKM")
<code>alpha</code>	Adjusts for the relative importance of RKM and FKM in the objective function; <code>alpha = 0.5</code> leads to reduced K-means, <code>alpha = 0</code> to factorial K-means, and <code>alpha = 1</code> reduces to the tandem approach (PCA followed by K-means)
<code>center</code>	A logical value indicating whether the variables should be shifted to be zero centered (default = TRUE)
<code>scale</code>	A logical value indicating whether the variables should be scaled to have unit variance before the analysis takes place (default = TRUE)
<code>rotation</code>	Specifies the method used to rotate the factors. Options are none for no rotation, <code>varimax</code> for varimax rotation with Kaiser normalization and <code>promax</code> for promax rotation (default = "none")
<code>nstart</code>	Number of starts (default = 100)
<code>smartStart</code>	If NULL then a random cluster membership vector is generated. Alternatively, a cluster membership vector can be provided as a starting solution
<code>seed</code>	An integer that is used as argument by <code>set.seed()</code> for offsetting the random number generator when <code>smartStart = NULL</code> . The default value is NULL.
<code>x</code>	For the print method, a class of <code>clusmca</code>
<code>object</code>	For the summary method, a class of <code>clusmca</code>
<code>mth</code>	For the fitted method, a character string that specifies the type of fitted value to return: "centers" for the observations center vector, or "class" for the observations cluster membership value
<code>...</code>	Not used

## Details

For the K-means part, the algorithm of Hartigan-Wong is used by default.

The hidden `print` and `summary` methods print out some key components of an object of class `cluspca`.

The hidden `fitted` method returns cluster fitted values. If method is `"classes"`, this is a vector of cluster membership (the cluster component of the `"cluspca"` object). If method is `"centers"`, this is a matrix where each row is the cluster center for the observation. The rownames of the matrix are the cluster membership values.

When `nclus = 1` the function returns the PCA solution and `plot(object)` shows the corresponding biplot.

## Value

<code>obscoord</code>	Object scores
<code>attcoord</code>	Variable scores
<code>centroid</code>	Cluster centroids
<code>cluster</code>	Cluster membership
<code>criterion</code>	Optimal value of the objective function
<code>size</code>	The number of objects in each cluster
<code>scale</code>	A copy of scale in the return object
<code>center</code>	A copy of center in the return object
<code>nstart</code>	A copy of nstart in the return object
<code>odata</code>	A copy of data in the return object

## References

De Soete, G., and Carroll, J. D. (1994). K-means clustering in a low-dimensional Euclidean space. In Diday E. et al. (Eds.), *New Approaches in Classification and Data Analysis*, Heidelberg: Springer, 212-219.

Vichi, M., and Kiers, H.A.L. (2001). Factorial K-means analysis for two-way data. *Computational Statistics and Data Analysis*, 37, 49-64.

## See Also

[clusmca](#), [cluspcamix](#), [tuneclus](#)

## Examples

```
#Reduced K-means with 3 clusters in 2 dimensions after 10 random starts
data(macro)
outRKM = cluspca(macro, 3, 2, method = "RKM", rotation = "varimax", scale = FALSE, nstart = 10)
summary(outRKM)
#Scatterplot (dimensions 1 and 2) and cluster description plot
plot(outRKM, cludesc = TRUE)
```



```

#Factorial K-means with 3 clusters in 2 dimensions
#with a Reduced K-means starting solution
data(macro)
outFKM = cluspca(macro, 3, 2, method = "FKM", rotation = "varimax",
scale = FALSE, smartStart = outRKM$cluster)
outFKM
#Scatterplot (dimensions 1 and 2) and cluster description plot
plot(outFKM, cludesc = TRUE)

#To get the Tandem approach (PCA(SVD) + K-means)
outTandem = cluspca(macro, 3, 2, alpha = 1, seed = 1234)
plot(outTandem)

#nclus = 1 just gives the PCA solution
#outPCA = cluspca(macro, 1, 2)
#outPCA
#Scatterplot (dimensions 1 and 2)
#plot(outPCA)

```

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cluspcamix

*Joint dimension reduction and clustering of mixed-type data.*


---

## Description

This function implements clustering and dimension reduction for mixed-type variables, i.e., categorical and metric (see, Yamamoto & Hwang, 2014; van de Velden, Iodice D'Enza, & Markos 2019; Vichi, Vicari, & Kiers, 2019). This framework includes Mixed Reduced K-means and Mixed Factorial K-means, as well as a compromise of these two methods. The methods combine Principal Component Analysis of mixed-data for dimension reduction with K-means for clustering.

## Usage

```

cluspcamix(data, nclus, ndim, method=c("mixedRKM", "mixedFKM"),
center = TRUE, scale = TRUE, alpha=NULL, rotation="none",
nstart = 100, smartStart=NULL, seed=NULL, inboot = FALSE)

## S3 method for class 'cluspcamix'
print(x, ...)

## S3 method for class 'cluspcamix'
summary(object, ...)

## S3 method for class 'cluspcamix'
fitted(object, mth = c("centers", "classes"), ...)

```

## Arguments

data                    Dataset with categorical and metric variables

nclus	Number of clusters (nclus = 1 returns the PCAMIX solution)
ndim	Dimensionality of the solution
method	Specifies the method. Options are mixedRKM for mixed reduced K-means and mixedFKM for mixed factorial K-means (default = "mixedRKM")
center	A logical value indicating whether the variables should be shifted to be zero centered (default = TRUE)
scale	A logical value indicating whether the variables should be scaled to have unit variance before the analysis takes place (default = TRUE)
alpha	Adjusts for the relative importance of Mixed RKM and Mixed FKM in the objective function; alpha = 0.5 leads to mixed reduced K-means, alpha = 0 to mixed factorial K-means, and alpha = 1 reduces to the tandem approach (PCAMIX followed by K-means)
rotation	Specifies the method used to rotate the factors. Options are none for no rotation, varimax for varimax rotation with Kaiser normalization and promax for promax rotation (default = "none")
nstart	Number of random starts (default = 100)
smartStart	If NULL then a random cluster membership vector is generated. Alternatively, a cluster membership vector can be provided as a starting solution
seed	An integer that is used as argument by set.seed() for offsetting the random number generator when smartStart = NULL. The default value is NULL.
inboot	Used internally in the bootstrap functions to perform bootstrapping on the indicator matrix.
x	For the print method, a class of cluspcamix
object	For the summary method, a class of cluspcamix
mth	For the fitted method, a character string that specifies the type of fitted value to return: "centers" for the observations center vector, or "class" for the observations cluster membership value
...	Not used

## Details

For the K-means part, the algorithm of Hartigan-Wong is used by default.

The hidden print and summary methods print out some key components of an object of class cluspcamix.

The hidden fitted method returns cluster fitted values. If method is "classes", this is a vector of cluster membership (the cluster component of the "cluspcamix" object). If method is "centers", this is a matrix where each row is the cluster center for the observation. The rownames of the matrix are the cluster membership values.

When nclus = 1 the function returns the solution of PCAMIX and plot(object) shows the corresponding biplot.

**Value**

obscoord	Object scores
attcoord	Variable scores
centroid	Cluster centroids
cluster	Cluster membership
criterion	Optimal value of the objective criterion
size	The number of objects in each cluster
scale	A copy of scale in the return object
center	A copy of center in the return object
nstart	A copy of nstart in the return object
odata	A copy of data in the return object

**References**

van de Velden, M., Iodice D'Enza, A., & Markos, A. (2019). Distance-based clustering of mixed data. *Wiley Interdisciplinary Reviews: Computational Statistics*, e1456.

Vichi, M., Vicari, D., & Kiers, H.A.L. (2019). Clustering and dimension reduction for mixed variables. *Behaviormetrika*. doi:10.1007/s41237-018-0068-6.

Yamamoto, M., & Hwang, H. (2014). A general formulation of cluster analysis with dimension reduction and subspace separation. *Behaviormetrika*, 41, 115-129.

**See Also**

[cluspca](#), [clusmca](#), [tuneclus](#)

**Examples**

```
data(diamond)
#Mixed Reduced K-means solution with 3 clusters in 2 dimensions
#after 10 random starts
outmixedRKM = cluspcamix(diamond, 3, 2, method = "mixedRKM", nstart = 10, seed = 1234)
outmixedRKM
#A graph with the categories and a biplot of the continuous variables (dimensions 1 and 2)
plot(outmixedRKM)

#Tandem analysis: PCAMIX or FAMD followed by K-means solution
#with 3 clusters in 2 dimensions after 10 random starts
outTandem = cluspcamix(diamond, 3, 2, alpha = 1, nstart = 10, seed = 1234)
outTandem
#Scatterplot (dimensions 1 and 2)
plot(outTandem)

#nclus = 1 just gives the PCAMIX or FAMD solution
#outPCAMIX = cluspcamix(diamond, 1, 2)
#outPCAMIX
#Biplot (dimensions 1 and 2)
#plot(outPCAMIX)
```

---

cmc

*Contraceptive Choice in Indonesia*

---

### Description

Data of married women in Indonesia who were not pregnant (or did not know they were pregnant) at the time of the survey. The dataset contains demographic and socio-economic characteristics of the women along with their preferred method of contraception (no use, long-term methods, short-term methods).

### Usage

```
data(cmc)
```

### Format

A data frame containing 1,437 observations on the following 10 variables.

W\_AGE wife's age in years.

W\_EDU ordered factor indicating wife's education, with levels "low", "2", "3" and "high".

H\_EDU ordered factor indicating wife's education, with levels "low", "2", "3" and "high".

NCHILD number of children.

W\_REL factor indicating wife's religion, with levels "non-Islam" and "Islam".

W\_WORK factor indicating if the wife is working.

H\_OCC factor indicating husband's occupation, with levels "1", "2", "3" and "4". The labels are not known.

SOL ordered factor indicating the standard of living index with levels "low", "2", "3" and "high".

MEDEXP factor indicating media exposure, with levels "good" and "not good".

CM factor indicating the contraceptive method used, with levels "no-use", "long-term" and "short-term".

### Source

This dataset is part of the 1987 National Indonesia Contraceptive Prevalence Survey and was created by Tjen-Sien Lim. It has been taken from the UCI Machine Learning Repository at <http://archive.ics.uci.edu/ml/>.

### References

Lim, T.-S., Loh, W.-Y. & Shih, Y.-S. (1999). A Comparison of Prediction Accuracy, Complexity, and Training Time of Thirty-three Old and New Classification Algorithms. *Machine Learning*, 40(3), 203-228.

### Examples

```
data(cmc)
```

---

diamond

*Diamond Stone Pricing*

---

### Description

Data on 308 diamond stones sold in Singapore. The main attributes are diamond weight, colour, clarity, certification body and price in Singapore \$. The weight of a diamond stone is indicated in terms of carat units. Since stones may be divided into 3 clusters due to their size, namely small (less than 0.5 carats), medium (0.5 to less than 1 carat) and large (1 carat and over), following Chu (2001), three binary variables have been built representing the three caratage ranges, and three quantitative variables (denoted Small, Medium, Large) have been derived by multiplying such binary variables by carats. So, the "Small" variable has nonzero values (i.e., the carat values) only for the smallest diamonds (less than 0.5 carats), and likewise for the other two variables. Thus, these variables are weighted binary variables. The colour of a diamond is graded from D (completely colourless), E, F, G, ..., to I (almost colorless). Clarity refers to the diamond's internal and external imperfections. Clarity is graded on a scale from IF (internally flawless), to very very slightly imperfect (VVS1 or VVS2), and very slightly imperfect, VS1 or VS2. Three certification bodies were used: New York based Gemmological Institute of America (GIA), Antwerp based International Gemmological Institute (IGI) and Hoge Raad Voor Diamant (HRD).

### Usage

```
data(diamond)
```

### Format

A data frame with 308 observations on the following 7 variables.

**Small** weighted binary variable with nonzero values (i.e., the carat values) for diamonds with less than 0.5 carats.

**Medium** weighted binary variable with nonzero values (i.e., the carat values) for diamonds from 0.5 to less than 1 carat.

**Large** weighted binary variable with nonzero values (i.e., the carat values) for diamonds from 1 carat and over.

**Colour** the color of the diamond with a factor with levels (D, E, F, G, H, I).

**Clarity** the clarity of the diamond with a factor with levels (IF, VVS1, VVS2, VS1, VS2).

**Certification** the certification body with a factor with levels (GIA, IGI, HRD).

**Price** the price of a diamond in Singapore \$.

### References

Chu, S. (2001). Pricing the C's of Diamond Stones, *Journal of Statistics Education*, 9(2).

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global_bootclus	<i>Global stability assessment of Joint Dimension Reduction and Clustering methods by bootstrapping.</i>
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### Description

Runs joint dimension and clustering algorithms repeatedly for different numbers of clusters on bootstrap replica of the original data and returns corresponding cluster assignments, and cluster agreement indices comparing pairs of partitions.

### Usage

```
global_bootclus(data, nclusrange = 3:4, ndim = NULL,
method = c("RKM", "FKM", "mixedRKM", "mixedFKM", "clusCA", "MCAk", "iFCB"),
nboot = 10, alpha = NULL, alphak = NULL, center = TRUE,
scale = TRUE, nstart = 100, smartStart = NULL, seed = NULL)
```

### Arguments

data	Continuous, Categorical or Mixed data set
nclusrange	An integer or an integer vector with the number of clusters or a range of numbers of clusters (should be greater than one)
ndim	Dimensionality of the solution; if NULL it is set to nclus - 1
method	Specifies the method. Options are RKM for Reduced K-means, FKM for Factorial K-means, mixedRKM for Mixed Reduced K-means, mixedFKM for Mixed Factorial K-means, MCAk for MCA K-means, iFCB for Iterative Factorial Clustering of Binary variables and clusCA for Cluster Correspondence Analysis.
nboot	Number of bootstrap pairs of partitions
alpha	Adjusts for the relative importance of (mixed) RKM and FKM in the objective function; alpha = 1 reduces to PCA/PCAMIX, alpha = 0.5 to (mixed) reduced K-means, and alpha = 0 to (mixed) factorial K-means
alphak	Non-negative scalar to adjust for the relative importance of MCA (alphak = 1) and K-means (alphak = 0) in the solution (default = .5). Works only in combination with method = "MCAk"
center	A logical value indicating whether the metric variables should be shifted to be zero centered (default = TRUE)
scale	A logical value indicating whether the metric variables should be scaled to have unit variance before the analysis takes place (default = TRUE)
nstart	Number of random starts (default = 100)
smartStart	If NULL then a random cluster membership vector is generated. Alternatively, a cluster membership vector can be provided as a starting solution
seed	An integer that is used as argument by set.seed() for offsetting the random number generator when smartStart = NULL. The default value is NULL.

## Details

The algorithm for assessing global cluster stability is similar to that in Dolnicar and Leisch (2010) and can be summarized in three steps:

*Step 1. Resampling:* Draw bootstrap samples  $S_i$  and  $T_i$  of size  $n$  from the data and use the original data,  $X$ , as evaluation set  $E_i = X$ . Apply the clustering method of choice to  $S_i$  and  $T_i$  and obtain  $C^{S_i}$  and  $C^{T_i}$ .

*Step 2. Mapping:* Assign each observation  $x_i$  to the closest centers of  $C^{S_i}$  and  $C^{T_i}$  using Euclidean distance, resulting in partitions  $C^{XS_i}$  and  $C^{XT_i}$ , where  $C^{XS_i}$  is the partition of the original data,  $X$ , predicted from clustering bootstrap sample  $S_i$  (same for  $T_i$  and  $C^{XT_i}$ ).

*Step 3. Evaluation:* Use the Adjusted Rand Index (ARI, Hubert & Arabie, 1985) or the Measure of Concordance (MOC, Pfitzner 2008) as measure of agreement and stability.

Inspect the distributions of ARI/MOC to assess the global reproducibility of the clustering solutions.

While  $nboot = 100$  is recommended, smaller run numbers could give quite informative results as well, if computation times become too high.

Note that the stability of a clustering solution is assessed, but stability is not the only important validity criterion - clustering solutions obtained by very inflexible clustering methods may be stable but not valid, as discussed in Hennig (2007).

## Value

nclusrange	An integer or an integer vector with the number of clusters or a range of numbers of clusters
clust1	Partitions, $C^{XS_i}$ of the original data, $X$ , predicted from clustering bootstrap sample $S_i$ (see Details)
clust2	Partitions, $C^{XT_i}$ of the original data, $X$ , predicted from clustering bootstrap sample $T_i$ (see Details)
index1	Indices of the original data rows in bootstrap sample $S_i$
index2	Indices of the original data rows in bootstrap sample $T_i$
rand	Adjusted Rand Index values
moc	Measure of Concordance values

## References

Hennig, C. (2007). Cluster-wise assessment of cluster stability. *Computational Statistics and Data Analysis*, 52, 258-271.

Pfitzner, D., Leibbrandt, R., & Powers, D. (2009). Characterization and evaluation of similarity measures for pairs of clusterings. *Knowledge and Information Systems*, 19(3), 361-394.

Dolnicar, S., & Leisch, F. (2010). Evaluation of structure and reproducibility of cluster solutions using the bootstrap. *Marketing Letters*, 21(1), 83-101.

## See Also

[local\\_bootclus](#)

## Examples

```
## 3 bootstrap replicates and nstart = 1 for speed in example,
## use at least 20 replicates for real applications
data(diamond)
boot_mixedRKM = global_bootclus(diamond[,-7], nclusrange = 3:4,
method = "mixedRKM", nboot = 3, nstart = 1, seed = 1234)

boxplot(boot_mixedRKM$rand, xlab = "Number of clusters", ylab =
"adjusted Rand Index")

## 5 bootstrap replicates and nstart = 10 for speed in example,
## use more for real applications
#data(macro)
#boot_RKM = global_bootclus(macro, nclusrange = 2:5,
#method = "RKM", nboot = 5, nstart = 10, seed = 1234)

#boxplot(boot_RKM$rand, xlab = "Number of clusters", ylab =
#"adjusted Rand Index")

## 5 bootstrap replicates and nstart = 1 for speed in example,
## use more for real applications
#data(bribery)
#boot_cluCA = global_bootclus(bribery, nclusrange = 2:5,
#method = "cluCA", nboot = 5, nstart = 1, seed = 1234)

#boxplot(boot_cluCA$rand, xlab = "Number of clusters", ylab =
#"adjusted Rand Index")
```

---

hsq

*Humor Styles*


---

## Description

The dataset was collected with an interactive online version of the Humor Styles Questionnaire (HSQ) which assesses four independent ways in which people express and appreciate humor (Martin et al. 2003): affiliative (items with prefix AF), defined as the benign uses of humor to enhance one's relationships with others; self-enhancing (SE), indicating uses of humor to enhance the self; aggressive (AG), the use of humor to enhance the self at the expense of others; self-defeating (SD), the use of humor to enhance relationships at the expense of oneself. The main part of the questionnaire consisted of 32 statements rated from 1 to 5 according to the respondents' level of agreement. The number of respondents is 993.

## Usage

```
data("hsq")
```



**Format**

A data frame with 993 observations on 32 Likert-type variables (statements) with 5 response categories, ranging from 1 (strong agreement) to 5 (strong disagreement).

- AF1 I usually don't laugh or joke around much with other people
- SE2 If I am feeling depressed, I can usually cheer myself up with humor
- AG3 If someone makes a mistake, I will often tease them about it
- SD4 I let people laugh at me or make fun at my expense more than I should
- AF5 I don't have to work very hard at making other people laugh - I seem to be a naturally humorous person
- SE6 Even when I'm by myself, I'm often amused by the absurdities of life
- AG7 People are never offended or hurt by my sense of humor
- SD8 I will often get carried away in putting myself down if it makes my family or friends laugh
- AF9 I rarely make other people laugh by telling funny stories about myself
- SE10 If I am feeling upset or unhappy I usually try to think of something funny about the situation to make myself feel better
- AG11 When telling jokes or saying funny things, I am usually not very concerned about how other people are taking it
- SD12 I often try to make people like or accept me more by saying something funny about my own weaknesses, blunders, or faults
- AF13 I laugh and joke a lot with my closest friends
- SE14 My humorous outlook on life keeps me from getting overly upset or depressed about things
- AG15 I do not like it when people use humor as a way of criticizing or putting someone down
- SD16 I don't often say funny things to put myself down
- AF17 I usually don't like to tell jokes or amuse people
- SE18 If I'm by myself and I'm feeling unhappy, I make an effort to think of something funny to cheer myself up
- AG19 Sometimes I think of something that is so funny that I can't stop myself from saying it, even if it is not appropriate for the situation
- SD20 I often go overboard in putting myself down when I am making jokes or trying to be funny
- AF21 I enjoy making people laugh
- SE22 If I am feeling sad or upset, I usually lose my sense of humor
- AG23 I never participate in laughing at others even if all my friends are doing it
- SD24 When I am with friends or family, I often seem to be the one that other people make fun of or joke about
- AF25 I don't often joke around with my friends
- SE26 It is my experience that thinking about some amusing aspect of a situation is often a very effective way of coping with problems
- AG27 If I don't like someone, I often use humor or teasing to put them down

- SD28 If I am having problems or feeling unhappy, I often cover it up by joking around, so that even my closest friends don't know how I really feel
- AF29 I usually can't think of witty things to say when I'm with other people
- SE30 I don't need to be with other people to feel amused - I can usually find things to laugh about even when I'm by myself
- AG31 Even if something is really funny to me, I will not laugh or joke about it if someone will be offended
- SD32 Letting others laugh at me is my way of keeping my friends and family in good spirits

## References

Martin, R. A., Puhlik-Doris, P., Larsen, G., Gray, J., & Weir, K. (2003). Individual differences in uses of humor and their relation to psychological well-being: Development of the Humor Styles Questionnaire. *Journal of Research in Personality*, 37(1), 48-75.

## Examples

```
data(hsq)
```

---

local_bootclus	<i>Cluster-wise stability assessment of Joint Dimension Reduction and Clustering methods by bootstrapping.</i>
----------------	--

---

## Description

Assessment of the cluster-wise stability of a joint dimension and clustering method. The data is resampled using bootstrapping and the Jaccard similarities of the original clusters to the most similar clusters in the resampled data are computed. The mean over these similarities is used as an index of the stability of a cluster. The method is similar to the one described in Hennig (2007).

## Usage

```
local_bootclus(data, nclus, ndim = NULL,
method = c("RKM", "FKM", "mixedRKM", "mixedFKM", "clusCA", "MCAk", "iFCB"),
scale = TRUE, center= TRUE, alpha = NULL, nstart=100,
nboot=10, alphak = .5, seed = NULL)
```

## Arguments

data	Continuous, Categorical or Mixed data set
nclus	Number of clusters
ndim	Dimensionality of the solution
method	Specifies the method. Options are RKM for Reduced K-means, FKM for Factorial K-means, mixedRKM for Mixed Reduced K-means, mixedFKM for Mixed Factorial K-means, MCAk for MCA K-means, iFCB for Iterative Factorial Clustering of Binary variables and clusCA for Cluster Correspondence Analysis.

scale	A logical value indicating whether the metric variables should be scaled to have unit variance before the analysis takes place (default = TRUE)
center	A logical value indicating whether the metric variables should be shifted to be zero centered (default = TRUE)
alpha	Adjusts for the relative importance of (mixed) RKM and FKM in the objective function; $\alpha = 1$ reduces to PCA/PCAMIX, $\alpha = 0.5$ to (mixed) reduced K-means, and $\alpha = 0$ to (mixed) factorial K-means
nstart	Number of random starts (default = 100)
nboot	Number of bootstrap pairs of partitions
alphak	Non-negative scalar to adjust for the relative importance of MCA ( $\text{alphak} = 1$ ) and K-means ( $\text{alphak} = 0$ ) in the solution (default = .5). Works only in combination with <code>method = "MCAk"</code>
seed	An integer that is used as argument by <code>set.seed()</code> for offsetting the random number generator when <code>smartStart = NULL</code> . The default value is NULL.

### Details

The algorithm for assessing local cluster stability is similar to that in Hennig (2007) and can be summarized in three steps:

*Step 1. Resampling:* Draw bootstrap samples  $S_i$  and  $T_i$  of size  $n$  from the data and use the original data as evaluation set  $E_i = X$ . Apply a joint dimension reduction and clustering method to  $S_i$  and  $T_i$  and obtain  $C^{\wedge}S_i$  and  $C^{\wedge}T_i$ .

*Step 2. Mapping:* Assign each observation  $x_i$  to the closest centers of  $C^{\wedge}S_i$  and  $C^{\wedge}T_i$  using Euclidean distance, resulting in partitions  $C^{\wedge}XS_i$  and  $C^{\wedge}XT_i$ .

*Step 3. Evaluation:* Obtain the maximum Jaccard agreement between each original cluster  $C_k$  and each one of the two bootstrap clusters,  $C_k^{\wedge}XS_i$  and  $C_k^{\wedge}XT_i$  as measure of agreement and stability, and take the average of each pair.

Inspect the distributions of the maximum Jaccard coefficients to assess the cluster level (local) stability of the solution.

Here are some guidelines for interpretation. Generally, a valid, stable cluster should yield a mean Jaccard similarity value of 0.75 or more. Between 0.6 and 0.75, clusters may be considered as indicating patterns in the data, but which points exactly should belong to these clusters is highly doubtful. Below average Jaccard values of 0.6, clusters should not be trusted. "Highly stable" clusters should yield average Jaccard similarities of 0.85 and above.

While  $B = 100$  is recommended, smaller run numbers could give quite informative results as well, if computation times become too high.

Note that the stability of a cluster is assessed, but stability is not the only important validity criterion - clusters obtained by very inflexible clustering methods may be stable but not valid, as discussed in Hennig (2007).

### Value

nclus	An integer with the number of clusters
clust1	Partitions, $C^{\wedge}XS_i$ of the original data, $X$ , predicted from clustering bootstrap sample $S_i$ (see Details)

clust2	Partitions, $C^{XT}_i$ of the original data, $X$ , predicted from clustering bootstrap sample $T_i$ (see Details)
index1	Indices of the original data rows in bootstrap sample $S_i$
index2	Indices of the original data rows in bootstrap sample $T_i$
Jaccard	Mean Jaccard similarity values

## References

Hennig, C. (2007). Cluster-wise assessment of cluster stability. *Computational Statistics and Data Analysis*, 52, 258-271.

## See Also

[global\\_bootclus](#)

## Examples

```
## 5 bootstrap replicates and nstart = 10 for speed in example,
## use more for real applications
data(iris)
bootres = local_bootclus(iris[,-5], nclus = 3, ndim = 2,
method = "RKM", nboot = 5, nstart = 1, seed = 1234)

boxplot(bootres$Jaccard, xlab = "cluster number", ylab =
"Jaccard similarity")

## 5 bootstrap replicates and nstart = 5 for speed in example,
## use more for real applications
#data(diamond)
#bootres = local_bootclus(diamond[,-7], nclus = 4, ndim = 3,
#method = "mixedRKM", nboot = 5, nstart = 10, seed = 1234)

#boxplot(bootres$Jaccard, xlab = "cluster number", ylab =
#"Jaccard similarity")

## 5 bootstrap replicates and nstart = 1 for speed in example,
## use more for real applications
#data(bribery)
#bootres = local_bootclus(bribery, nclus = 5, ndim = 4,
#method = "clusCA", nboot = 10, nstart = 1, seed = 1234)

#boxplot(bootres$Jaccard, xlab = "cluster number", ylab =
#"Jaccard similarity")
```

---

macro

*Economic Indicators of 20 OECD countries for 1999*

---

### Description

Data on the macroeconomic performance of national economies of 20 countries, members of the OECD (September 1999). The performance of the economies reflects the interaction of six main economic indicators (percentage change from the previous year): gross domestic product (GDP), leading indicator (LI), unemployment rate (UR), interest rate (IR), trade balance (TB), net national savings (NNS).

### Usage

```
data(macro)
```

### Format

A data frame with 20 observations on the following 6 variables.

GDP numeric

LI numeric

UR numeric

IR numeric

TB numeric

NNS numeric

### References

Vichi, M. & Kiers, H. A. (2001). Factorial k-means analysis for two-way data. *Computational Statistics & Data Analysis*, 37(1), 49-64.

---

mybond

*James Bond films*

---

### Description

The data set refers to 26 James Bond films produced up to 2021, based on 10 film characteristics: 7 continuous (year of release, production budget, box office gross in the USA and worldwide, running time, IMDB average rating, Rotten Tomatoes rating) and 3 categorical (Bond actor, native country of the actor playing the villain, native country of the actor playing the Bond girl). All figures in USD are adjusted for inflation. Most of the data was compiled from the Wikipedia page: [https://en.wikipedia.org/wiki/List\\_of\\_James\\_Bond\\_films](https://en.wikipedia.org/wiki/List_of_James_Bond_films).

**Usage**

```
data("mybond")
```

**Format**

A data frame with 26 observations on the following 10 variables.

year Year of release

budget Official production budget (in million USD)

grossusa Box office gross in the USA (in million USD)

grosswrld Box office gross worldwide (in million USD)

rtime Running time in minutes

IMDB IMDB rating

rottentomatoes Rotten Tomatoes rating

actor Bond actor

villaincnt Native country of the actor playing the villain

bondgirlcnt Native country of the actor playing the Bond girl

**Examples**

```
data(mybond)
```

---

plot.clusmca

*Plotting function for clusmca() output.*

---

**Description**

Plotting function that creates a scatterplot of the object scores and/or the attribute scores and the cluster centroids. Optionally, the function returns a series of barplots showing the standardized residuals per attribute for each cluster.

**Usage**

```
## S3 method for class 'clusmca'  
plot(x, dims = c(1,2), what = c(TRUE,TRUE),  
      cludesc = FALSE, topstdres = 20, objlabs = FALSE, attlabs = NULL,  
      subplot = FALSE, max.overlaps=10, ...)
```

**Arguments**

x	Object returned by clusmca()
dims	Numerical vector of length 2 indicating the dimensions to plot on horizontal and vertical axes respectively; default is first dimension horizontal and second dimension vertical
what	Vector of two logical values specifying the contents of the plots. First entry indicates whether a scatterplot of the objects is displayed in principal coordinates. Second entry indicates whether a scatterplot of the attribute categories is displayed in principal coordinates. Cluster centroids are always displayed. The default is c(TRUE, TRUE) and the resultant plot is a biplot of both objects and attribute categories with gamma-based scaling (see van de Velden et al., 2017)
cludesc	A logical value indicating whether a series of barplots is produced showing the largest (in absolute value) standardized residuals per attribute for each cluster (default = FALSE)
topstdres	Number of largest standardized residuals used to describe each cluster (default = 20). Works only in combination with cludesc = TRUE
objlabs	A logical value indicating whether object labels will be plotted; if TRUE row names of the data matrix are used (default = FALSE). Warning: when TRUE, execution time of the plotting function will increase dramatically as the number of objects gets larger
atllabs	Vector of custom attribute labels; if not provided, default labeling is applied
subplot	A logical value indicating whether a subplot with the full distribution of the standardized residuals will appear at the bottom left corner of the corresponding plots. Works only in combination with cludesc = TRUE
max.overlaps	Maximum number of text labels allowed to overlap. Defaults to 10
...	Further arguments to be transferred to clusmca()

**Value**

The function returns a ggplot2 scatterplot of the solution obtained via clusmca() that can be further customized using the **ggplot2** package. When cludesc = TRUE the function also returns a series of ggplot2 barplots showing the largest (or all) standardized residuals per attribute for each cluster.

**References**

- Hwang, H., Dillon, W. R., and Takane, Y. (2006). An extension of multiple correspondence analysis for identifying heterogenous subgroups of respondents. *Psychometrika*, 71, 161-171.
- Iodice D'Enza, A., and Palumbo, F. (2013). Iterative factor clustering of binary data. *Computational Statistics*, 28(2), 789-807.
- van de Velden M., Iodice D'Enza, A., and Palumbo, F. (2017). Cluster correspondence analysis. *Psychometrika*, 82(1), 158-185.

**See Also**

[plot.cluspca](#), [plot.cluspcamix](#)

## Examples

```

data("mybond")
#Cluster Correspondence Analysis with 3 clusters in 2 dimensions after 10 random starts
outclusCA = clusmca(mybond[,8:10], 3, 2, nstart = 100, seed = 234)
#Save the ggplot2 scatterplot
map = plot(outclusCA, max.overlaps = 40)$map
#Customization (adding titles)
map + ggtitle(paste("Cluster CA plot of the James bond categorical data: 3 clusters of sizes ",
                    paste(outclusCA$size, collapse = ", "), sep = "")) +
  xlab("Dim. 1") + ylab("Dim. 2") +
  theme(plot.title = element_text(size = 10, face = "bold", hjust = 0.5))

data("mybond")
#i-FCB with 3 clusters in 2 dimensions after 10 random starts
outclusCA = clusmca(mybond[,8:10], 3, 2, method = "iFCB", nstart= 10)
#Scatterplot with the observations only (dimensions 1 and 2)
#and cluster description plots showing the 20 largest std. residuals
#(with the full distribution showing in subplots)
plot(outclusCA, dim = c(1,2), what = c(TRUE, FALSE), cludesc = TRUE,
      subplot = TRUE)

```

---

plot.cluspca

*Plotting function for cluspca() output.*

---

## Description

Plotting function that creates a scatterplot of the objects, a correlation circle of the variables or a biplot of both objects and variables. Optionally, it returns a parallel coordinate plot showing cluster means.

## Usage

```

## S3 method for class 'cluspca'
plot(x, dims = c(1, 2), cludesc = FALSE,
      what = c(TRUE,TRUE), attlabs, max.overlaps=10, ...)

```

## Arguments

x	Object returned by cluspca()
dims	Numerical vector of length 2 indicating the dimensions to plot on horizontal and vertical axes respectively; default is first dimension horizontal and second dimension vertical
what	Vector of two logical values specifying the contents of the plots. First entry indicates whether a scatterplot of the objects and cluster centroids is displayed and the second entry whether a correlation circle of the variables is displayed. The default is c(TRUE, TRUE) and the resultant plot is a biplot of both objects and variables



cludesc	A logical value indicating if a parallel coordinate plot showing cluster means is produced (default = FALSE)
atllabs	Vector of custom attribute labels; if not provided, default labeling is applied
max.overlaps	Maximum number of text labels allowed to overlap. Defaults to 10
...	Further arguments to be transferred to cluspca()

### Value

The function returns a `ggplot2` scatterplot of the solution obtained via `cluspca()` that can be further customized using the **ggplot2** package. When `cludesc = TRUE` the function also returns a `ggplot2` parallel coordinate plot.

### References

De Soete, G., and Carroll, J. D. (1994). K-means clustering in a low-dimensional Euclidean space. In Diday E. et al. (Eds.), *New Approaches in Classification and Data Analysis*, Heidelberg: Springer, 212-219.

Vichi, M., and Kiers, H.A.L. (2001). Factorial K-means analysis for two-way data. *Computational Statistics and Data Analysis*, 37, 49-64.

### See Also

[plot.clusmca](#), [plot.cluspcamix](#)

### Examples

```
data("macro")
#Factorial K-means (3 clusters in 2 dimensions) after 100 random starts
outFKM = cluspca(macro, 3, 2, method = "FKM", rotation = "varimax")
#Scatterplot (dimensions 1 and 2) and cluster description plot
plot(outFKM, cludesc = TRUE)

data("iris", package = "datasets")
#Compromise solution between PCA and Reduced K-means
#on the iris dataset (3 clusters in 2 dimensions) after 100 random starts
outclusPCA = cluspca(iris[,-5], 3, 2, alpha = 0.3, rotation = "varimax")
table(outclusPCA$cluster, iris[,5])
#Save the ggplot2 scatterplot
map = plot(outclusPCA)$map
#Customization (adding titles)
map + ggtitle(paste("A compromise solution between RKM and FKM on the iris:
3 clusters of sizes ", paste(outclusPCA$size,
collapse = ", ", sep = "")) + xlab("Dimension 1") + ylab("Dimension 2") +
theme(plot.title = element_text(size = 10, face = "bold", hjust = 0.5))
```

---

plot.cluspcamix      *Plotting function for cluspcamix() output.*

---

### Description

Plotting function that creates a scatterplot of the objects, a correlation circle of the variables or a biplot of both objects and variables. Optionally, for metric variables, it returns a parallel coordinate plot showing cluster means and for categorical variables, a series of barplots showing the standardized residuals per attribute for each cluster.

### Usage

```
## S3 method for class 'cluspcamix'
plot(x, dims = c(1, 2), cludesc = FALSE,
     topstdres = 20, objlabs = FALSE, attlabs = NULL, attcatlabs = NULL,
     subplot = FALSE, what = c(TRUE,TRUE), max.overlaps = 10, ...)
```

### Arguments

x	Object returned by cluspcamix()
dims	Numerical vector of length 2 indicating the dimensions to plot on horizontal and vertical axes respectively; default is first dimension horizontal and second dimension vertical
what	Vector of two logical values specifying the contents of the plots. First entry indicates whether a scatterplot of the objects and cluster centroids is displayed and the second entry whether a correlation circle of the variables is displayed. The default is c(TRUE, TRUE) and the resultant plot is a biplot of both objects and variables
cludesc	A logical value indicating if a parallel coordinate plot showing cluster means is produced (default = FALSE)
topstdres	Number of largest standardized residuals used to describe each cluster (default = 20). Works only in combination with cludesc = TRUE
subplot	A logical value indicating whether a subplot with the full distribution of the standardized residuals will appear at the bottom left corner of the corresponding plots. Works only in combination with cludesc = TRUE
objlabs	A logical value indicating whether object labels will be plotted; if TRUE row names of the data matrix are used (default = FALSE). Warning: when TRUE, execution time of the plotting function will increase dramatically as the number of objects gets larger
attlabs	Vector of custom labels of continuous attributes; if not provided, default labeling is applied
attcatlabs	Vector of custom labels of categorical attributes (categories); if not provided, default labeling is applied
max.overlaps	Maximum number of text labels allowed to overlap. Defaults to 10
...	Further arguments to be transferred to cluspcamix()

**Value**

The function returns a `ggplot2` scatterplot of the solution obtained via `cluspcamix()` that can be further customized using the **ggplot2** package. When `cludesc = TRUE`, for metric variables, the function also returns a `ggplot2` parallel coordinate plot and for categorical variables, a series of `ggplot2` barplots showing the largest (or all) standardized residuals per attribute for each cluster.

**References**

van de Velden, M., Iodice D'Enza, A., & Markos, A. (2019). Distance-based clustering of mixed data. *Wiley Interdisciplinary Reviews: Computational Statistics*, e1456.

Vichi, M., Vicari, D., & Kiers, H. A. L. (2019). Clustering and dimension reduction for mixed variables. *Behaviormetrika*. doi:10.1007/s41237-018-0068-6.

**See Also**

[plot.clusmca](#), [plot.cluspca](#)

**Examples**

```
data(diamond)
#Mixed Reduced K-means solution with 3 clusters in 2 dimensions
#after 10 random starts
outmixedRKM = cluspcamix(diamond, 3, 2, method = "mixedRKM", nstart = 10)
#Scatterplot (dimensions 1 and 2)
plot(outmixedRKM, cludesc = TRUE)
```

---

tuneclus

*Cluster quality assessment for a range of clusters and dimensions.*


---

**Description**

This function facilitates the selection of the appropriate number of clusters and dimensions for joint dimension reduction and clustering methods.

**Usage**

```
tuneclus(data, nclusrange = 3:4, ndimrange = 2:3,
method = c("RKM", "FKM", "mixedRKM", "mixedFKM", "clusCA", "iFCB", "MCAK"),
criterion = "asw", dst = "full", alpha = NULL, alphak = NULL,
center = TRUE, scale = TRUE, rotation = "none", nstart = 100,
smartStart = NULL, seed = NULL)

## S3 method for class 'tuneclus'
print(x, ...)

## S3 method for class 'tuneclus'
summary(object, ...)
```

```
## S3 method for class 'tuneclus'
fitted(object, mth = c("centers", "classes"), ...)
```

### Arguments

<code>data</code>	Continuous, Categorical or Mixed data set
<code>nclusrange</code>	An integer vector with the range of numbers of clusters which are to be compared by the cluster validity criteria. Note: the number of clusters should be greater than one
<code>ndimrange</code>	An integer vector with the range of dimensions which are to be compared by the cluster validity criteria
<code>method</code>	Specifies the method. Options are RKM for reduced K-means, FKM for factorial K-means, mixedRKM for mixed reduced K-means, mixedFKM for mixed factorial K-means, MCAk for MCA K-means, iFCB for Iterative Factorial Clustering of Binary variables and clusCA for Cluster Correspondence Analysis
<code>criterion</code>	One of <code>asw</code> , <code>ch</code> or <code>crit</code> . Determines whether average silhouette width, Calinski-Harabasz index or objective value of the selected method is used (default = "asw")
<code>dst</code>	Specifies the data used to compute the distances between objects. Options are <code>full</code> for the original data (after possible scaling) and <code>low</code> for the object scores in the low-dimensional space (default = "full")
<code>alpha</code>	Adjusts for the relative importance of (mixed) RKM and FKM in the objective function; $\alpha = 1$ reduces to PCA, $\alpha = 0.5$ to (mixed) reduced K-means, and $\alpha = 0$ to (mixed) factorial K-means
<code>alphak</code>	Non-negative scalar to adjust for the relative importance of MCA ( $\text{alphak} = 1$ ) and K-means ( $\text{alphak} = 0$ ) in the solution (default = .5). Works only in combination with <code>method = "MCAk"</code>
<code>center</code>	A logical value indicating whether the variables should be shifted to be zero centered (default = TRUE)
<code>scale</code>	A logical value indicating whether the variables should be scaled to have unit variance before the analysis takes place (default = TRUE)
<code>rotation</code>	Specifies the method used to rotate the factors. Options are <code>none</code> for no rotation, <code>varimax</code> for varimax rotation with Kaiser normalization and <code>promax</code> for promax rotation (default = "none")
<code>nstart</code>	Number of starts (default = 100)
<code>smartStart</code>	If NULL then a random cluster membership vector is generated. Alternatively, a cluster membership vector can be provided as a starting solution
<code>seed</code>	An integer that is used as argument by <code>set.seed()</code> for offsetting the random number generator when <code>smartStart = NULL</code> . The default value is NULL.
<code>x</code>	For the print method, a class of <code>clusmca</code>
<code>object</code>	For the summary method, a class of <code>clusmca</code>

meth	For the fitted method, a character string that specifies the type of fitted value to return: "centers" for the observations center vector, or "class" for the observations cluster membership value
...	Not used

### Details

For the K-means part, the algorithm of Hartigan-Wong is used by default.

The hidden print and summary methods print out some key components of an object of class tuneclus.

The hidden fitted method returns cluster fitted values. If method is "classes", this is a vector of cluster membership (the cluster component of the "tuneclus" object). If method is "centers", this is a matrix where each row is the cluster center for the observation. The rownames of the matrix are the cluster membership values.

### Value

clusobjbest	The output of the optimal run of cluspca() or clumca()
nclusbest	The optimal number of clusters
ndimbest	The optimal number of dimensions
critbest	The optimal criterion value for nclusbest clusters and ndimbest dimensions
critgrid	Matrix of size nclusrange x ndimrange with the criterion values for the specified ranges of clusters and dimensions (values are calculated only when the number of clusters is greater than the number of dimensions; otherwise values in the grid are left blank)
criterion	"asw" for average Silhouette width or "ch" for "Calinski-Harabasz"
cluasw	Average Silhouette width values of each cluster, when criterion = "asw"

### References

Calinski, R.B., and Harabasz, J., (1974). A dendrite method for cluster analysis. *Communications in Statistics*, 3, 1-27.

Kaufman, L., and Rousseeuw, P.J., (1990). *Finding Groups in Data: An Introduction to Cluster Analysis*. Wiley, New York.

### See Also

[global\\_bootclus](#), [local\\_bootclus](#)

### Examples

```
# Reduced K-means for a range of clusters and dimensions
data(macro)
# Cluster quality assessment based on the average silhouette width in the low dimensional space
# nstart = 1 for speed in example
# use more for real applications
bestRKM = tuneclus(macro, 3:4, 2:3, method = "RKM",
```

```
criterion = "asw", dst = "low", nstart = 1, seed = 1234)
bestRKM
#plot(bestRKM)

# Cluster Correspondence Analysis for a range of clusters and dimensions
data(bribery)
# Cluster quality assessment based on the Callinski-Harabasz index in the full dimensional space
bestclusCA = tuneclus(bribery, 4:5, 3:4, method = "clusCA",
criterion = "ch", nstart = 20, seed = 1234)
bestclusCA
#plot(bestclusCA, cludesc = TRUE)

# Mixed reduced K-means for a range of clusters and dimensions
data(diamond)
# Cluster quality assessment based on the average silhouette width in the low dimensional space
# nstart = 5 for speed in example
# use more for real applications
bestmixedRKM = tuneclus(diamond[,-7], 3:4, 2:3,
method = "mixedRKM", criterion = "asw", dst = "low",
nstart = 5, seed = 1234)
bestmixedRKM
#plot(bestmixedRKM)
```

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