

# Package ‘trimcluster’

May 8, 2026

**Title** Cluster Analysis with Trimming

**Version** 0.2-0

**VersionNote** Released 0.1-6 on 2025-06-28 on CRAN

**Depends** R (>= 1.9.0)

**Imports** tclust

**Suggests** fpc

**Description** Trimmed k-means clustering. The method is described in Cuesta-Albertos et al. (1997) <[doi:10.1214/aos/1031833664](https://doi.org/10.1214/aos/1031833664)>.

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**License** GPL (>= 3)

**URL** <https://github.com/valentint/trimcluster>

**BugReports** <https://github.com/valentint/trimcluster/issues>

**Repository** CRAN

**Date/Publication** 2025-07-17 08:40:01 UTC

**NeedsCompilation** no

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trimkmeans	<i>Trimmed k-means clustering</i>
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### Description

The trimmed k-means clustering method by Cuesta-Albertos, Gordaliza and Matran (1997). This optimizes the k-means criterion under trimming a portion of the points.

### Usage

```
trimkmeans(data,k,trim=0.1, scaling=FALSE,
           runs=500, niter1=3, niter2=20, nkeep=5, points=NULL,
           countmode, printcrit, maxit,
           parallel=FALSE, n.cores=-1, trace=0, ...)

## S3 method for class 'tkm'
print(x, ...)
## S3 method for class 'tkm'
plot(x, data, ...)
```

### Arguments

<code>data</code>	matrix or data.frame with raw data
<code>k</code>	integer. Number of clusters.
<code>trim</code>	numeric between 0 and 1. Proportion of points to be trimmed.
<code>scaling</code>	logical. If TRUE, the variables are centered at their means and scaled to unit variance before execution.
<code>runs</code>	The number of random initializations to be performed.
<code>niter1</code>	The number of concentration steps to be performed for the nstart initializations.
<code>niter2</code>	The maximum number of concentration steps to be performed for the nkeep solutions kept for further iteration. The concentration steps are stopped, whenever two consecutive steps lead to the same data partition.
<code>nkeep</code>	The number of iterated initializations (after niter1 concentration steps) with the best values in the target function that are kept for further iterations
<code>points</code>	NULL or a matrix with k vectors used as means to initialize the algorithm. If initial mean vectors are specified, runs should be 1 (otherwise the same initial means are used for all runs).
<code>countmode</code>	(deprecated) optional positive integer. Every countmode algorithm runs trimkmeans shows a message.
<code>printcrit</code>	(deprecated) logical. If TRUE, all criterion values (mean squares) of the algorithm runs are printed.
<code>maxit</code>	(deprecated, use the combination nkeep, niter1 and niter2) The maximum number of concentration steps to be performed. The concentration steps are stopped, whenever two consecutive steps lead to the same data partition.

parallel	A logical value, specifying whether the nstart initializations should be done in parallel.
n.cores	The number of cores to use when paralellizing, only taken into account if parallel=TRUE.
trace	Defines the tracing level, which is set to 0 by default. Tracing level 1 gives additional information on the stage of the iterative process.
x	object of class tkm.
...	further arguments to be transferred to plot or <a href="#">plotcluster</a> .

### Details

The function `trimkmeans()` now calls the function `tkmeans()` from the package `tclust`. This makes the procedure much faster since (a) `tkmeans()` is implemented in C++, (b) a new random initialization is introduced (see the parameters `niter1`, `niter2` and `nkeep` which replace the previous `maxit` and (c) it is possible to run the initialization in parallel (see the argument `parallel` and `ncores`).

`plot.tkm` calls [plotcluster](#) if the dimensionality of the data `p` is 1, shows a scatterplot with non-trimmed regions if `p=2` and discriminant coordinates computed from the clusters (ignoring the trimmed points) if `p>2`.

### Value

An object of class 'tkm' which is a LIST with components

classification	integer vector coding cluster membership with trimmed observations coded as <code>k+1</code> .
means	numerical matrix giving the mean vectors of the <code>k</code> classes.
disttom	vector of squared Euclidean distances of all points to the closest mean.
ropt	maximum value of <code>disttom</code> so that the corresponding point is not trimmed.
k	see above.
trim	see above.
runs	see above.
scaling	see above.

### Author(s)

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### References

Cuesta-Albertos, J. A., Gordaliza, A., and Matran, C. (1997) Trimmed k-Means: An Attempt to Robustify Quantizers, *Annals of Statistics*, 25, 553-576.

### See Also

[plotcluster](#)

**Examples**

```
set.seed(10001)
n1 <-60
n2 <-60
n3 <-70
n0 <-10
nn <- n1+n2+n3+n0
pp <- 2
X <- matrix(rep(0,nn*pp),nrow=nn)
ii <-0
for (i in 1:n1){
  ii <-ii+1
  X[ii,] <- c(5,-5)+rnorm(2)
}
for (i in 1:n2){
  ii <- ii+1
  X[ii,] <- c(5,5)+rnorm(2)*0.75
}
for (i in 1:n3){
  ii <- ii+1
  X[ii,] <- c(-5,-5)+rnorm(2)*0.75
}
for (i in 1:n0){
  ii <- ii+1
  X[ii,] <- rnorm(2)*8
}
tkm1 <- trimkmeans(X, k=3, trim=0.1, runs=5)
## runs=5 is used to save computing time; runs must be >= nkeep

print(tkm1)
plot(tkm1,X)
```

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