

# Package: PUGMM (via r-universe)

September 12, 2024

**Version** 0.1.0

**Title** Parsimonious Ultrametric Gaussian Mixture Models

**Description** Finite Gaussian mixture models with parsimonious extended ultrametric covariance structures estimated via a grouped coordinate ascent algorithm, which is equivalent to the Expectation-Maximization algorithm. The thirteen ultrametric covariance structures implemented allow for the inspection of different hierarchical relationships among variables. The estimation of an ultrametric correlation matrix is included as a function. The methodologies are described in Cavicchia, Vichi, Zaccaria (2024) <[doi:10.1007/s11222-024-10405-9](https://doi.org/10.1007/s11222-024-10405-9)>, Cavicchia, Vichi, Zaccaria (2022) <[doi:10.1007/s11634-021-00488-x](https://doi.org/10.1007/s11634-021-00488-x)> and Cavicchia, Vichi, Zaccaria (2020) <[doi:10.1007/s11634-020-00400-z](https://doi.org/10.1007/s11634-020-00400-z)>.

**Depends** R (>= 4.0)

**Imports** ClusterR, doParallel, foreach, igraph, MASS, Matrix, mclust, mcompanion, ppclust

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**URL** <https://github.com/giorgiazaccaria/PUGMM>

**BugReports** <https://github.com/giorgiazaccaria/PUGMM/issues>

**NeedsCompilation** no

**LazyData** yes

**Encoding** UTF-8

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**Repository** <https://giorgiazaccaria.r-universe.dev>

**RemoteUrl** <https://github.com/giorgiazaccaria/pugmm>

**RemoteRef** HEAD

**RemoteSha** 0b1a5134e8450078423cd7822985e4ce5a43a12e

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penguins	<i>Penguins</i>
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### Description

The data set contains five measurements made on 342 penguins which are classified into three species.

### Usage

```
data(penguins)
```

### Format

A data frame with 342 observations and 5 variables, which are described as follows.

**species** Penguin species (Chinstrap, Adélie, or Gentoo)

**culmen\_length\_mm** Culmen length (mm)

**culmen\_depth\_mm** Culmen depth (mm)

**flipper\_length\_mm** Flipper length (mm)

**body\_mass\_g** Body mass (g)

### Details

Data were collected and made available by Dr. Kristen Gorman and the Palmer Station, Antarctica LTER, a member of the Long Term Ecological Research Network. The categorical variables 'island' and 'sex' have been removed from the original dataset, as well as the incomplete observations on the five variables reported herein.

### Source

Dataset downloaded from Kaggle <https://www.kaggle.com/code/parulpandey/penguin-dataset-the-new-iris>.

**References**

Gorman, K.B., Williams T.D., Fraser W.R. (2014). Ecological sexual dimorphism and environmental variability within a community of Antarctic penguins (genus *Pygoscelis*). *PLoS ONE*, 9(3), e90081.

**Examples**

```
data(penguins)
```

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plot.pugmm	<i>Plotting method for pugmm object</i>
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**Description**

Plots for Parsimonious Ultrametric Gaussian Mixture Models results, such as BIC and path diagrams.

**Usage**

```
## S3 method for class 'pugmm'
plot(x, what = NULL, nrow = NULL, ncol = NULL, cluster_names = NULL, ...)
```

**Arguments**

x	Output from pugmm.
what	A string specifying the type of graph requested. Available choices are:  "\"BIC\" Plot of BIC values for the fitted models. For each $G$ , the best BIC among the ones corresponding to different $m$ is displayed. "\"Path Diagram\" Path diagram representation of the extended ultrametric covariance matrix per component for the best model.
nrow	Number of rows in the graphical window. A new graphical window is opened every 6 plots, i.e., components of pugmm.
ncol	Number of columns in the graphical window. A new graphical window is opened every 6 plots, i.e., components of pugmm.
cluster_names	String of dimension $G$ with the clusters/components' name.
...	Other graphics parameters.

**Value**

No return value since this is a plot method.

**See Also**

[pugmm\(\)](#)

**Examples**

```
data(penguins)
x <- scale(penguins[, 2:5])
pugmm.penguins <- pugmm(x, 3, 1)
plot.pugmm(pugmm.penguins, what = c("BIC", "Path Diagram"))
```

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pugmm

*Parsimonious Ultrametric Gaussian Mixture Models*


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**Description**

Model-based clustering via Parsimonious Ultrametric Gaussian Mixture Models. Hierarchical relationships among variables within and between clusters are inspected. The grouped coordinate ascent algorithm is used for the parameter estimation. The optimal model is selected according to BIC.

**Usage**

```
pugmm(
  X,
  G = NULL,
  m = NULL,
  normalization = NULL,
  model = NULL,
  maxiter = 500,
  tol = 1e-06,
  stop = "aitken",
  rndstart = 1,
  initG = "kmeans",
  initm = "ucms",
  gaussian = "mclust",
  parallel = FALSE
)
```

**Arguments**

<code>X</code>	$(n \times p)$ numeric matrix or data frame, where $n$ and $p$ represent the number of units and variables, respectively. Categorical variables are not allowed.
<code>G</code>	Integer (vector) specifying the number of mixture components (default: $G = 1:5$ ).
<code>m</code>	Integer (vector) specifying the number of variable groups (default: $m = 1:5$ ).
<code>normalization</code>	Character string specifying the data transformation. If <code>NULL</code> , no transformation is applied to the data matrix (default). Other options are: "standard" for the standardization; "center" for centering the data; "range" for the MinMax transformation; "SVD" for the Singular Value Decomposition transformation.

model	Vector of character strings indicating the model names to be fitted. If NULL, all the possible models are fitted (default). See the possible models using <code>available_models()</code> .
maxiter	Integer value specifying the maximum number of iterations of the algorithm (default: <code>maxiter = 500</code> ).
tol	Numeric value specifying the tolerance for the convergence criterion (default: <code>tol = 1e-6</code> ).
stop	Character string specifying the convergence criteria. If "aitken", the Aitken acceleration-based stopping rule is used (default); if "relative", the relative log-likelihood in two sequential iterations is evaluated.
rndstart	Integer value specifying the number of random starts (default: <code>rndstart = 1</code> ).
initG	Character string specifying the method for the initialization of the unit-component membership. If "kmeans", k-means via <code>RcppArmadillo</code> is used (default). Other options are: "random" for random assignment; "kmeansf" for fuzzy c-means (via the function <code>fcm</code> of the package <code>ppclust</code> ).
initm	Character string specifying the method for the initialization of the variable-group membership. If "ucms", the multivariate model to be used for obtaining the variable-group membership estimated is the same <code>model.name</code> used for estimating the Parsimonious Ultrametric Gaussian Mixture Model (default); if "random", a random assignment is performed.
gaussian	Character string specifying the way to compute the log-likelihood. If "mclust", <code>dmvnorm</code> of <code>mclust</code> is used (default); if "canonical", the log-likelihood computation is based upon the canonical representation of an extended ultrametric covariance matrix.
parallel	A logical value, specifying whether the models should be run in parallel.

### Details

The grouped coordinate ascent algorithm used for the estimation of PUGMMs parameters was demonstrated to be equivalent to an Expectation-Maximization algorithm in the GMM framework (Hathaway, 1986).

### Value

An object of class `pugmm` containing the results of the optimal - according to BIC - Parsimonious Ultrametric Gaussian Mixture Model estimation.

`call` Matched call.

`X` Input data matrix.

`G` Number of components of the best model.

`m` Number of variable groups of the best model.

`label` Integer vector of dimension  $n$ , taking values in  $\{1, \dots, G\}$ . It identifies the unit classification according to the maximum a posteriori of the best model.

`pp` Numeric vector of dimension  $G$  containing the prior probabilities for the best model.

`mu` ( $G \times p$ ) numeric matrix containing the component mean vectors (by row) for the best model.

`sigma` List of dimension  $G$  containing the  $(p \times p)$  numeric component extended ultrametric covariance matrices for the best model.

`V` List of dimension  $G$  containing the  $(p \times m)$  binary variable-group membership matrices for the best model.

`Sv` List of dimension  $G$  containing the  $(m \times m)$  numeric diagonal matrices of the group variances for the best model.

`Sw` List of dimension  $G$  containing the  $(m \times m)$  numeric diagonal matrices of the within-group covariances for the best model.

`Sb` List of dimension  $G$  containing the  $(m \times m)$  numeric hallow matrices of the between-group covariances for the best model.

`post`  $(n \times G)$  numeric matrix containing the posterior probabilities for the best model.

`pm` Number of parameters of the best model.

`pm.cov` Number of covariance parameters of the best model.

`pm.free` Number of free parameters of the best model (`pm - (constraints on V + count.constr.SwSb + count.constr.SvSw)`).

`count.constr.SwSb` Number of times the constraint between `Sw` and `Sb` has been turned on for the best model.

`count.constr.SvSw` Number of times the constraint between `Sv` and `Sw` has been turned on for the best model.

`BIC` BIC values for all the fitted models. If BIC is *NA*, the model has not been computed since its structure is equal to another model, while if BIC is *-Inf* the solution has a number of clusters  $< G$ .

`bic` BIC value of the best model.

`loglik` Log-likelihood of the best model.

`loop` Random start corresponding to the selected solution of the best model.

`iter` Number of iterations needed to estimate the best model.

`model.name` Character string denoting the PUGMM model name of the best model among the ones fitted.

`messages` Messages.

## References

- Cavicchia, C., Vichi, M., Zaccaria, G. (2024) Parsimonious ultrametric Gaussian mixture models. *Statistics and Computing*, 34, 108.
- Cavicchia, C., Vichi, M., Zaccaria, G. (2022) Gaussian mixture model with an extended ultrametric covariance structure. *Advances in Data Analysis and Classification*, 16(2), 399-427.
- Hathaway, R. (1986) Another interpretation of the EM algorithm for mixture distributions. *Statistics and Probability Letters*, 4(2), 53-56.

## See Also

[pugmm\\_available\\_models\(\)](#), [plot.pugmm\(\)](#)

## Examples

```
data(penguins)
x <- scale(penguins[, 2:5])
pugmm.penguins <- pugmm(x, 3, 1)
table(penguins$species, pugmm.penguins$label)

pugmm.penguins <- pugmm(x)
pugmm.penguins$G
pugmm.penguins$m
pugmm.penguins$model.name
```

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pugmm\_available\_models

*PUGMM Model Names*

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## Description

Description of the model names used in the *PUGMM* package.

## Usage

```
pugmm_available_models()
```

## Details

The PUGMM model names in the *PUGMM* package are characterized by four letters:

- First letter: it refers to the variable-group membership matrix  $V$ , which can be equal (E) or free to vary (F) across components.
- Second, third, fourth letters: they refer to the matrices of the group variances  $\Sigma_V$ , the within-group covariances  $\Sigma_W$  and the between-group covariances  $\Sigma_B$ , respectively, by indicating if they are unique (U, i.e., equal within and across components), isotropic (I, i.e., equal within components), equal (E, i.e., equal across components) or free to vary across components (F).

## Value

Available models in PUGMM, i.e., the thirteen extended ultrametric covariance structures of PUGMM.

## References

Cavicchia, C., Vichi, M., Zaccaria, G. (2024) Parsimonious ultrametric Gaussian mixture models. *Statistics and Computing*, 34, 108.

## See Also

[pugmm\(\)](#)

**Examples**

```
pugmm_available_models()
```

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rand.member	<i>Random partition of objects into classes</i>
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**Description**

Performs a random partition of objects into classes.

**Usage**

```
rand.member(n.obs, G)
```

**Arguments**

n.obs	Number of objects
G	Number of classes

**Details**

No empty classes can occur.

**Value**

A binary and row-stochastic matrix with *n.obs* rows and *G* columns.

**Examples**

```
rand.member(10, 3)
```

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UCM	<i>Ultrametric Correlation Matrix</i>
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**Description**

Fit an ultrametric correlation matrix on a nonnegative correlation one.

**Usage**

```
UCM(R, m, rndstart, maxiter = 100, eps = 1e-06)
```



**Arguments**

R	$(p \times p)$ nonnegative correlation matrix.
m	Integer specifying the number of variable groups.
rndstart	Integer value specifying the number of random starts.
maxiter	Integer value specifying the maximum number of iterations of the EM algorithm (default: maxiter = 100).
eps	Numeric value specifying the tolerance for the convergence criterion used in the coordinate descent algorithm (default: eps = 1e-6).

**Value**

A list with the following elements:

call Matched call.

V Optimal binary and row-stochastic  $(p \times m)$  variable-group membership matrix.

Rt Optimal  $(p \times p)$  ultrametric correlation matrix.

Rw Optimal  $(m \times m)$  within-concept consistency (diagonal) matrix.

Rb Optimal  $(m \times m)$  between-concept correlation matrix.

of Objective function corresponding to the optimal solution.

loop Random start corresponding to the optimal solution.

i ter Number of iterations needed to obtain the optimal solution.

**References**

Cavicchia, C., Vichi, M., Zaccaria, G. (2020) The ultrametric correlation matrix for modelling hierarchical latent concepts. *Advances in Data Analysis and Classification*, 14(4), 837-853.

**Examples**

```
data(penguins)
R <- cor(penguins[, 2:5])
UCM(R, 4, 1)
```

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