

# Package: leptokurticMixture (via r-universe)

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**Title** Implements Parsimonious Finite Mixtures of Multivariate  
Elliptical Leptokurtic-Normals

**Version** 1.1

**Description** A way to fit Parsimonious Finite Mixtures of Multivariate  
Elliptical Leptokurtic-Normals. Two methods of estimation are  
implemented.

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compareEstimation	<i>Compare the two methods of estimation</i>
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### Description

Compare the two methods of estimation for fitting a finite mixture of multivariate elliptical leptokurtic-normal distributions; fixed point iterations and MM algorithm.

### Usage

```
compareEstimation(  
  mod = NULL,  
  data = NULL,  
  G = NULL,  
  n = 10^4,  
  tol = 1e-06,  
  wt = NULL,  
  n0 = 25,  
  lab = NULL  
)
```

### Arguments

mod	A character of length 4 such as "VVVV", indicating the model; the covariance and beta parameters.
data	A n x p matrix of observations.
G	The number of components to fit.
n	The maximum number of EM iterations.
tol	The tolerance for the stopping rule; lack of progress. The default is 1e-6 but it depends on the dataset.
wt	a (n x d) matrix of weights for initialization if NULL, then a random weight matrix is generated.
n0	Given wt, the number of iterations used to obtain the initial parameters
lab	Using given labels (lab) as starting values.

### Value

A vector of times, number of iterations and log-likelihood values.

**Description**

Performs a number of iterations of the EM for the multivariate elliptical leptokurtic-normal (MLN) distribution until the tolerance for the lack progress or the maximum number of iterations is reached. An implementation of parsimonious clustering models via the eigen-decomposition of the scatter matrix and allowing the concentration parameter to be varying, equal or fixed across components.

**Usage**

```
EM(
  data = NULL,
  G = 2,
  model = NULL,
  km1 = c(1, 0, 1),
  n = 10,
  epsilon = 0.01,
  gpar0 = NULL,
  estimation = 1,
  label = NULL
)
```

**Arguments**

<code>data</code>	A $n \times p$ matrix of observations.
<code>G</code>	A integer determine the number of components of the mixture model.
<code>model</code>	a character of length 4 such as "VVVV", indicating the model; the covariance and beta parameters. The 1st position controls, lambda, the volume; "V" varying across components or "E" equal across components. The 2nd position controls the eigenvalues; "V" varying across components, "E" equal across components or "I" the identity matrix. The 3rd position controls the orientation; "V" varying across components, "E" equal across components or "I" the identity matrix. The 4th position controls the concentration, beta; "V" varying across components, "E" equal across components, "F" fixed at the maximum value.
<code>km1</code>	a vector of length 3 indicating, the number of k-means starts, number of random starts and the number of EM iterations used for each start
<code>n</code>	The maximum number of EM iterations.
<code>epsilon</code>	The tolerance for the stopping rule; lack of progress. The default is $1e-6$ but it depends on the dataset.
<code>gpar0</code>	A list of model parameters .
<code>estimation</code>	If 1 (default) use the fixed point iterations and if 2 the MM algorithm.
<code>label</code>	If NULL then the data has no known groups. If is integer then some of the observations have known groups. If <code>label[i]=k</code> then observation belongs to group k. If <code>label[i]=0</code> then observation has no known group.

**Value**

A list with following items

- loglik - A vector of the loglikelihood values
- gpar - A list containing the parameters values
- z - A  $n \times G$  matrix of the posterior probabilities
- map - A vector the maximum a posteriori derived from z
- label - The input provided.
- numpar - The number of free parameters in the fitted model.
- maxLoglik - The largest value from loglik.

**Examples**

```
x1 = rmln(n=100, d=4, mu=rep(5,4), diag(4), beta=2)
x2 = rmln(n=100, d=4, mu=rep(-5,4), diag(4), beta=2)
x = rbind( x1,x2)
mlnFit = EM(data=x, G=2, model="VVVF")
```

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pmln

*Parsimonious model-based clustering with the multivariate elliptical leptokurtic-normal*

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**Description**

Performs parsimonious clustering with the multivariate elliptical leptokurtic-normal (MLN). There are 14 possible scale matrix structure and 2 for the kurtosis parameter for a total of 28 models.

**Usage**

```
pmln(
  data = NULL,
  G = 1:3,
  covModels = NULL,
  betaModels = "B",
  kml = c(1, 0, 1),
  label = NULL,
  scale.data = TRUE,
  veo = FALSE,
  iterMax = 1000,
  tol = 1e-08,
  pprogress = FALSE,
  method = "FP"
)
```

**Arguments**

<code>data</code>	A $n \times p$ matrix of observations.
<code>G</code>	A integer determine the number of components of the mixture model.
<code>covModels</code>	if NULL fit 14 possible scale matrix structures. Otherwise a character vector where each element has length 3. e.g. <code>c("VVV", "EEE")</code> A character of length 4 such as "VVVV", indicating the model; the covariance and beta parameters. The 1st position controls, lambda, the volume; "V" varying across components or "E" equal across components. The 2nd position controls the eigenvalues; V" varying across components, "E" equal across components or "I" the identity matrix. The 3rd position controls the orientation; "V" varying across components, "E" equal across components or "I" the identity matrix.
<code>betaModels</code>	set to "V", "E", "B", "F". "V" varying across components, "E" equal across components, "B" consider both "V" & "E", "F" fixed at the maximum value.
<code>km1</code>	a vector of length 3 indicating, the number of k-means starts, number of random starts and the number of EM iterations used for each start
<code>label</code>	If NULL then the data has no known groups. If is integer then some of the observations have known groups. If <code>label[i]=k</code> then observation belongs to group k. If <code>label[i]=0</code> then observation has no known group.
<code>scale.data</code>	Should the data be scaled before clustering. The default is TRUE.
<code>veo</code>	"Variables exceed observations". If TRUE, fit the model even though the number variables in the model exceeds the number of observations.
<code>iterMax</code>	The maximum number of EM iterations for each model fitted.
<code>tol</code>	The tol for the stopping rule; lack of progress. The default is 1e-6 but it depends on the data set.
<code>pprogress</code>	If TRUE print the progress of the function.
<code>method</code>	If FP use the fixed point iteration method otherwise if MM use the MM method.

**Value**

A list of

- `startobject` - A statement on how the models were initialized
- `gpar` - A list of parameter values for the model chosen by the BIC
- `loglik` - A vector of the log-likelihoods values
- `z` - A  $n \times G$  matrix of the posterior probabilities from the model chosen by the BIC
- `map` - A vector the maximum a posteriori derived from `z`
- `BIC` - An array with dimensions (G, number of fitted models, 3). The last dimension indices the loglik, number of free parameters and BIC for each fitted model.
- `bicModel` - Information as list on the model chosen by the BIC.

**Examples**

```
x1 = rmln(n=100, d=4, mu=rep(5,4), diag(4), beta=2)
x2 = rmln(n=100, d=4, mu=rep(-5,4), diag(4), beta=2)
x = rbind(x1,x2)
m1nFit = pmln(data=x, G=2, covModels=c("VVV", "EEE"), betaModels="B")
```

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rmln	<i>Generate realizations from the multivariate elliptical leptokurtic-normal distribution</i>
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**Description**

This function calculates the log cumulative density function for the multivariate-t with scale matrix equal to the identity matrix. It finds the mode and then uses Gaussian quadrature to estimate the integral.

**Usage**

```
rmln(n = NULL, d = NULL, mu = NULL, Sigma = NULL, beta = NULL)
```

**Arguments**

n	number of observations
d	the dimension of the observations
mu	location parameter of length d
Sigma	(d x d) scatter matrix
beta	the concentration parameter

**Value**

A (n x d) matrix of realizations

**Examples**

```
x = rmln(n=10, d=4, mu=rep(0,4), diag(4), beta=2)
```

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