

# Package ‘EMCluster’

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**Title** EM Algorithm for Model-Based Clustering of Finite Mixture  
Gaussian Distribution

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**LazyData** yes

**Description** EM algorithms and several efficient  
initialization methods for model-based clustering of finite  
mixture Gaussian distribution with unstructured dispersion  
in both of unsupervised and semi-supervised learning.

**License** Mozilla Public License 2.0

**BugReports** <https://github.com/snoweye/EMCluster/issues>

**URL** <https://github.com/snoweye/EMCluster>

**NeedsCompilation** yes

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EMCluster-package	<i>EM Algorithm for Model-Based Clustering of Finite Mixture Gaussian Distribution</i>
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## Description

EMCluster provides EM algorithms and several efficient initialization methods for model-based clustering of finite mixture Gaussian distribution with unstructured dispersion in both of unsupervised and semi-supervised clustering.

## Details

The install command is simply as

```
> R CMD INSTALL EMCluster_0.2-0.tar.gz
```

from a command mode or

```
R> install.packages("EMCluster")
```

inside an R session.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [emcluster](#).

**Examples**

```
## Not run:
demo(allinit, 'EMCluster', ask = F, echo = F)
demo(allinit_ss, 'EMCluster', ask = F, echo = F)

## End(Not run)
```

---

Assign Class

*Assign Class Id*

---

**Description**

This function assigns cluster id to each observation in  $x$  according to the desired model `emobj` or specified parameters `pi`, `Mu`, and `LTSigma`.

**Usage**

```
assign.class(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL,
            lab = NULL, return.all = TRUE)
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>emobj</code>	the desired model which is a list mainly contains <code>pi</code> , <code>Mu</code> , and <code>LTSigma</code> , usually a returned object from <code>init.EM</code> .
<code>pi</code>	the mixing proportion, length $K$ .
<code>Mu</code>	the centers of clusters, dimension $K \times p$ .
<code>LTSigma</code>	the lower triangular matrices of dispersion, dimension $K \times p(p + 1)/2$ .
<code>lab</code>	labeled data for semi-supervised clustering, length $n$ .
<code>return.all</code>	if returning with a whole <code>emobj</code> object.

**Details**

This function are based either an input emobj or inputs pi, Mu, and LTSigma to assign class id to each observation of  $x$ .

If lab is submitted, then the observation with label id greater 0 will not be assigned new class.

**Value**

This function returns a list containing mainly two new variables: nc (length  $K$  numbers of observations in each class) and class (length  $n$  class id).

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [emcluster](#).

**Examples**

```
library(EMCluster, quietly = TRUE)
set.seed(1234)
x2 <- da2$da

ret <- init.EM(x2, nclass = 2)
ret.new <- assign.class(x2, ret, return.all = FALSE)
str(ret.new)
```

---

Conversion

*Convert Matrices in Different Format*

---

**Description**

These utility functions are to convert matrices in different formats.

**Usage**

```
LTSigma2variance(x)
variance2LTSigma(x)
LTSigma2var(x1, p = NULL)
var2LTSigma(x1)
class2Gamma(class)
Gamma2class(Gamma)
```

**Arguments**

x	a matrix/array to be converted, the dimension could be $K \times p(p + 1)/2$ or $p \times p \times K$ .
x1	a vector/matrix to be converted, the length and dimension could be $p(p + 1)/2$ and $p \times p$ .
p	dimension of matrix.
class	id of clusters for each observation, length $n$ .
Gamma	containing posterior probabilities if normalized, otherwise containing component densities weighted by mixing proportion, dimension $n \times K$ .

**Details**

LTSigma2variance converts LTSigma format to 3D array, and variance2LTSigma is the inversion function.

LTsigma2var converts LTsigma format to a matrix, and var2LTsigma is the inversion function. Note that LTsigma is one component of LTSigma.

class2Gamma converts id to a Gamma matrix where with probability 1 for the cluster where the observation belongs to, and Gamma2class converts posterior to cluster id where largest posterior is picked for each observation.

**Value**

A vector/matrix/array is returned.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [emcluster](#).

**Examples**

```
## Not run:
library(EMcluster, quietly = TRUE)
x <- da2$LTSigma
class <- da2$class

y <- LTSigma2variance(x)
str(y)
y <- variance2LTSigma(y)
str(y)
sum(x != y)
```

```
Gamma <- class2Gamma(class)
class.new <- Gamma2class(Gamma)
sum(class != class.new)

## End(Not run)
```

---

Dataset

*Dataset for demonstrations*

---

## Description

There are four small datasets to test and demonstrate **EMCluster**.

## Usage

```
da1
da2
da3
```

## Format

da1, da2, da3 are in list.

## Details

da1 has 500 observations in two dimensions da1\$da\$x and da1\$da\$y, and they are in 10 clusters given in da1\$class.

da2 has 2,500 observations in two dimensions, too. The true parameters are given in da1\$pi, da1\$Mu, and da1\$LTSigma. There are 40 clusters given in da1\$class for this dataset.

da3 is similar to da2, but with lower overlaps between clusters.

## Author(s)

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

## References

<https://www.stat.iastate.edu/people/ranjan-maitra>

**Description**

These are core functions of **EMCluster** performing EM algorithm for model-based clustering of finite mixture multivariate Gaussian distribution with unstructured dispersion.

**Usage**

```
emcluster(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL,
          lab = NULL, EMC = .EMC, assign.class = FALSE)
shortemcluster(x, emobj = NULL, pi = NULL, Mu = NULL,
               LTSigma = NULL, maxiter = 100, eps = 1e-2)
simple.init(x, nclass = 1)
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>emobj</code>	the desired model which is a list mainly contains <code>pi</code> , <code>Mu</code> , and <code>LTSigma</code> , usually a returned object from <code>init.EM</code> .
<code>pi</code>	the mixing proportion, length $K$ .
<code>Mu</code>	the centers of clusters, dimension $K \times p$ .
<code>LTSigma</code>	the lower triangular matrices of dispersion, $K \times p(p + 1)/2$ .
<code>lab</code>	labeled data for semi-supervised clustering, length $n$ .
<code>EMC</code>	the control for the EM iterations.
<code>assign.class</code>	if assigning class id.
<code>maxiter</code>	maximum number of iterations.
<code>eps</code>	convergent tolerance.
<code>nclass</code>	the desired number of clusters, $K$ .

**Details**

The `emcluster` mainly performs EM iterations starting from the given parameters `emobj` without other initializations.

The `shortemcluster` performs short-EM iterations as described in `init.EM`.

**Value**

The `emcluster` returns an object `emobj` with class `emret` which can be used in post-process or other functions such as `e.step`, `m.step`, `assign.class`, `em.ic`, and `dmixmvn`.

The `shortemcluster` also returns an object `emobj` with class `emret` which is the best of several random initializations.

The `simple.init` utilizes `rand.EM` to obtain a simple initial.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [e.step](#), [m.step](#), [.EMControl](#).

**Examples**

```
library(EMcluster, quietly = TRUE)
set.seed(1234)
x1 <- da1$da

emobj <- simple.init(x1, nclass = 10)
emobj <- shortemcluster(x1, emobj)
summary(emobj)

ret <- emcluster(x1, emobj, assign.class = TRUE)
summary(ret)
```

---

EM Control

*EM Control Generator and Controller*

---

**Description**

The `.EMControl` generates an EM control (`.EMC`) controlling the options and conditions of EM algorithms, i.e. this function generate a default template. One can either modify `.EMC` or employ this function to control EM algorithms. By default, `.EMC`, `.EMC.Rnd`, and `.EC.Rndp` are three native controllers as the **EMCluster** is loaded.

**Usage**

```
.EMControl(alpha = 0.99, short.iter = 200, short.eps = 1e-2,
  fixed.iter = 1, n.candidate = 3,
  em.iter = 1000, em.eps = 1e-6, exhaust.iter = 5)
.EMC
.EMC.Rnd
.EMC.Rndp
```



**Arguments**

<code>alpha</code>	only used in emgroup for "SVD" initialization.
<code>short.iter</code>	number of short-EM steps, default = 200.
<code>short.eps</code>	tolerance of short-EM steps, default = 1e-2.
<code>fixed.iter</code>	fixed iterations of EM for "RndEM" initialization, default = 1.
<code>n.candidate</code>	reserved for other initialization methods (unimplemented).
<code>em.iter</code>	maximum number of long-EM steps, default = 1000.
<code>em.eps</code>	tolerance of long-EM steps, default = 1e-6.
<code>exhaust.iter</code>	number of iterations for "exhaustEM" initialization, default = 5.

**Details**

`exhaust.iter` and `fixed.iter` are used to control the iterations of initialization procedures.

`short.iter` and `short.eps` are used to control the short-EM iterations.

`em.iter` and `em.eps` are used to control the long-EM iterations.

Moreover, `short.eps` and `em.eps` are for checking convergence of the iterations.

**Value**

This function returns a list as `.EMC` by default.

The `.EMC.Rnd` is equal to `.EMControl(short.eps = Inf)` and usually used by the `rand.EM` method.

The `.EMC.Rndp` is equal to `.EMControl(fixed.iter = 5)` where each random initials run 5 EM iterations in the `rand.EM` method.

**Author(s)**

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [emcluster](#).

**Examples**

```
library(EMcluster, quietly = TRUE)

.EMC <- .EMControl()
.EMC.Rnd <- .EMControl(short.eps = Inf)
.EMC.Rndp <- .EMControl(fixed.iter = 5)
```

**Description**

These functions are tools for compute information criteria for the fitted models.

**Usage**

```
em.ic(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL,
      llhdval = NULL)
em.aic(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL)
em.bic(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL)
em.clc(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL)
em.icl(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL)
em.icl.bic(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL)
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>emobj</code>	the desired model which is a list mainly contains <code>pi</code> , <code>Mu</code> , and <code>LTSigma</code> , usually a returned object from <code>init.EM</code> .
<code>pi</code>	the mixing proportion, length $K$ .
<code>Mu</code>	the centers of clusters, dimension $K \times p$ .
<code>LTSigma</code>	the lower triangular matrices of dispersion, $K \times p(p + 1)/2$ .
<code>llhdval</code>	the total log likelihood value of <code>x</code> given <code>emobj</code> .

**Details**

The `em.ic` calls all other functions to compute AIC (`em.aic`), BIC (`em.bic`), CLC (`em.clc`), ICL (`em.icl`), and ICL.BIC (`em.icl.bic`). All are useful information criteria for model selections, mainly choosing number of cluster.

**Value**

`em.ic` returns a list containing all other information criteria for given the data `x` and the desired model `emobj`.

**Author(s)**

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#).

**Examples**

```
library(EMcluster, quietly = TRUE)
x2 <- da2$da

emobj <- list(pi = da2$pi, Mu = da2$Mu, LTSigma = da2$LTSigma)
em.ic(x2, emobj = emobj)
```

---

Initialization and EM *Initialization and EM Algorithm*

---

**Description**

These functions perform initializations (including `em.EM` and `RndEM`) followed by the EM iterations for model-based clustering of finite mixture multivariate Gaussian distribution with unstructured dispersion in both of unsupervised and semi-supervised clusterings.

**Usage**

```
init.EM(x, nclass = 1, lab = NULL, EMC = .EMC,
        stable.solution = TRUE, min.n = NULL, min.n.iter = 10,
        method = c("em.EM", "Rnd.EM"))
em.EM(x, nclass = 1, lab = NULL, EMC = .EMC,
       stable.solution = TRUE, min.n = NULL, min.n.iter = 10)
rand.EM(x, nclass = 1, lab = NULL, EMC = .EMC.Rnd,
        stable.solution = TRUE, min.n = NULL, min.n.iter = 10)
exhaust.EM(x, nclass = 1, lab = NULL,
           EMC = .EMControl(short.iter = 1, short.eps = Inf),
           method = c("em.EM", "Rnd.EM"),
           stable.solution = TRUE, min.n = NULL, min.n.iter = 10);
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>nclass</code>	the desired number of clusters, $K$ .
<code>lab</code>	labeled data for semi-supervised clustering, length $n$ .
<code>EMC</code>	the control for the EM iterations.
<code>stable.solution</code>	if returning a stable solution.
<code>min.n</code>	restriction for a stable solution, the minimum number of observations for every final clusters.
<code>min.n.iter</code>	restriction for a stable solution, the minimum number of iterations for trying a stable solution.
<code>method</code>	an initialization method.

## Details

The `init.EM` calls either `em.EM` if `method="em.EM"` or `rand.EM` if `method="Rnd.EM"`.

The `em.EM` has two steps: short-EM has loose convergent tolerance controlled by `.EMC$short.eps` and try several random initializations controlled by `.EMC$short.iter`, while long-EM starts from the best short-EM result (in terms of log likelihood) and run to convergence with a tight tolerance controlled by `.EMC$em.eps`.

The `rand.EM` also has two steps: first randomly pick several random initializations controlled by `.EMC$short.iter`, and second starts from the best of the random result (in terms of log likelihood) and run to convergence.

The `lab` is only for the semi-supervised clustering, and it contains pre-labeled indices between 1 and  $K$  for labeled observations. Observations with index 0 is non-labeled and has to be clustered by the EM algorithm. Indices will be assigned by the results of the EM algorithm. See `demo(allinit_ss, 'EMCluster')` for details.

The `exhaust.EM` also calls the `init.EM` with different EMC and perform `exhaust.iter` times of EM algorithm with different initials. The best result is returned.

## Value

These functions return an object `emobj` with class `emret` which can be used in post-process or other functions such as `e.step`, `m.step`, `assign.class`, `em.ic`, and `dmixmvm`.

## Author(s)

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

## References

<https://www.stat.iastate.edu/people/ranjan-maitra>

## See Also

[emcluster](#), [.EMControl](#).

## Examples

```
## Not run:
library(EMCluster, quietly = TRUE)
set.seed(1234)
x <- da1$da

ret.em <- init.EM(x, nclass = 10, method = "em.EM")
ret.Rnd <- init.EM(x, nclass = 10, method = "Rnd.EM", EMC = .EMC.Rnd)

emobj <- simple.init(x, nclass = 10)
ret.init <- emcluster(x, emobj, assign.class = TRUE)

par(mfrow = c(2, 2))
plotem(ret.em, x)
plotem(ret.Rnd, x)
```

```
plotem(ret.init, x)

## End(Not run)
```

---

Jaccard Index

*Jaccard Index*

---

### Description

This function returns the Jaccard index for binary ids.

### Usage

```
Jaccard.Index(x, y)
```

### Arguments

x	true binary ids, 0 or 1.
y	predicted binary ids, 0 or 1.

### Details

All ids, x and y, should be either 0 (not active) or 1 (active). Any value other than 1 will be converted to 0.

### Value

Return the value of Jaccard index.

### Author(s)

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

### References

<https://www.stat.iastate.edu/people/ranjan-maitra>

### Examples

```
library(EMcluster, quietly = TRUE)

x.id <- c(1, 1, 1, 0, 0, 0, 3, 3, 3)
y.id <- c(0, 1, 0, 1, 1, 1, 0, 1, 1)

Jaccard.Index(x.id, y.id)
```

---

**Likelihood Mixture Tests***Likelihood Mixture Tests*

---

**Description**

This function test two mixture Gaussian models with unstructured covariance matrix and different numbers of clusters.

**Usage**

```
lmt(emobj.0, emobj.a, x, tau = 0.5, n.mc.E.delta = 1000,  
    n.mc.E.chi2, verbose = FALSE)
```

**Arguments**

<code>emobj.0</code>	a emret object for the null hypothesis.
<code>emobj.a</code>	a emret object for the alternative hypothesis.
<code>x</code>	the data matrix, dimension $n \times p$ .
<code>tau</code>	proportion of null and alternative hypotheses.
<code>n.mc.E.delta</code>	number of Monte Carlo simulations for expectation of delta (difference of logL).
<code>n.mc.E.chi2</code>	number of Monte Carlo simulations for expectation of chisquare statistics.
<code>verbose</code>	if verbose.

**Details**

This function calls several subroutines to compute information, likelihood ratio statistics, degrees of freedom, non-centrality of chi-squared distributions ...etc. Based on Monte Carlo methods to estimate parameters of likelihood mixture tests, this function return a p-value for testing  $H_0$ : `emobj.0` v.s.  $H_a$ : `emobj.a`.

**Value**

A list of class `lmt` are returned.

**Author(s)**

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#).

**Examples**

```
## Not run:
library(EMCluster, quietly = TRUE)
set.seed(1234)

x <- as.matrix(iris[, 1:4])
p <- ncol(x)
min.n <- p * (p + 1) / 2
.EMC$short.iter <- 200

ret.2 <- init.EM(x, nclass = 2, min.n = min.n, method = "Rnd.EM")
ret.3 <- init.EM(x, nclass = 3, min.n = min.n, method = "Rnd.EM")
ret.4 <- init.EM(x, nclass = 4, min.n = min.n, method = "Rnd.EM")

(lmt.23 <- lmt(ret.2, ret.3, x))
(lmt.34 <- lmt(ret.3, ret.4, x))
(lmt.24 <- lmt(ret.2, ret.4, x))

## End(Not run)
```

---

LMT Functions

*Likelihood Mixture Test (LMT) Functions of EMCluster*

---

**Description**

All likelihood mixture test (LMT) functions are for testing and can be utilized by advanced developers with caution.

Currently, these are only for workflows.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

---

MVN

*Density of (Mixture) Multivariate Normal Distribution*

---

**Description**

These functions are tools for compute density of (mixture) multivariate Gaussian distribution with unstructured dispersion.

**Usage**

```
dmvn(x, mu, LTsigma, log = FALSE)
dlmvn(x, mu, LTsigma, log = TRUE)
dmixmvn(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL, log = FALSE)
logL(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL)
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>mu</code>	the centers of clusters, length $p$ .
<code>LTsigma</code>	the lower triangular matrices of dispersion, length $p(p + 1)/2$ .
<code>log</code>	if logarithm returned.
<code>emobj</code>	the desired model which is a list mainly contains <code>pi</code> , <code>Mu</code> , and <code>LTSigma</code> , usually a returned object from <code>init.EM</code> .
<code>pi</code>	the mixing proportion, length $K$ .
<code>Mu</code>	the centers of clusters, dimension $K \times p$ .
<code>LTSigma</code>	the lower triangular matrices of dispersion, $K \times p(p + 1)/2$ .

**Details**

The `dmvn` and `dlmvn` compute density and log density of multivariate distribution.

The `dmixmvn` computes density of mixture multivariate distribution and is based either an input `emobj` or inputs `pi`, `Mu`, and `LTSigma` to assign class id to each observation of `x`.

The `logL` returns the value of the observed log likelihood function of the parameters at the current values of the parameters `pi`, `Mu`, and `LTSigma`, with the supplied data matrix `x`.

**Value**

A density value is returned.

**Author(s)**

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [emcluster](#).



**Examples**

```

library(EMcluster, quietly = TRUE)
x2 <- da2$da
x3 <- da3$da

emobj2 <- list(pi = da2$pi, Mu = da2$Mu, LTSigma = da2$LTSigma)
emobj3 <- list(pi = da3$pi, Mu = da3$Mu, LTSigma = da3$LTSigma)

logL(x2, emobj = emobj2)
logL(x3, emobj = emobj3)

dmixmvn2 <- dmixmvn(x2, emobj2)
dmixmvn3 <- dmixmvn(x3, emobj3)

dlmvn(da2$da[1,], da2$Mu[1,], da2$LTSigma[1,])
log(dmvn(da2$da[1,], da2$Mu[1,], da2$LTSigma[1,]))

```

---

Other Initializations *Other Initializations*

---

**Description**

Two more functions with different initialization method.

**Usage**

```

starts.via.svd(x, nclass = 1, method = c("em", "kmeans"),
              EMC = .EMC)
emgroup(x, nclass = 1, EMC = .EMC)

```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>nclass</code>	the desired number of clusters, $K$ .
<code>method</code>	method with the svd initializations.
<code>EMC</code>	the control for the EM iterations.

**Details**

The `starts.via.svd` utilizes SVD to initial parameters, and the `emgroup` runs the EM algorithm starting from the initial.

**Value**

The `starts.via.svd` returns an object with class `svd`, and the `emgroup` returns and object `emobj` with class `emret`.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [.EMControl](#).

**Examples**

```
library(EMcluster, quietly = TRUE)
set.seed(1234)
x1 <- da1$da

emobj <- emgroup(x1, nclass = 10)
summary(emobj)

ret.0 <- starts.via.svd(x1, nclass = 10, method = "kmeans")
summary(ret.0)
```

---

Plot EM Results

*Plot Two Dimensional Data with clusters*

---

**Description**

The functions plot two dimensional data for clusters.

**Usage**

```
plotem(emobj, x, main = NULL, xlab = NULL, ylab = NULL,
      ...)
plot2d(x, emobj = NULL, k = NULL, color.pch = 1,
      append.BN = TRUE, ...)
```

**Arguments**

<code>emobj</code>	the desired model which is a list mainly contains <code>pi</code> , <code>Mu</code> , and <code>LTSigma</code> , usually a returned object from <code>init.EM</code> .
<code>x</code>	the data matrix, dimension $n \times p$ .
<code>main</code>	title of plot.
<code>xlab</code>	label of x-axis.
<code>ylab</code>	label of y-axis.
<code>...</code>	other parameters to the plot.

k	index for symbols.
color.pch	color and style for symbols.
append.BN	if appending bivariate normal ellipsoid.

**Details**

This a simple x-y lot.

**Value**

A plot is returned.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [emcluster](#).

**Examples**

```
## Not run:  
library(EMcluster, quietly = TRUE)  
x1 <- da1$da  
  
ret.1 <- starts.via.svd(x1, nclass = 10, method = "em")  
summary(ret.1)  
  
plotem(ret.1, x1)  
  
## End(Not run)
```

---

Plot Multivariate Data

*Plot Multivariate Data*

---

**Description**

The function plots multivariate data for clusters as the parallel coordinates plot.

**Usage**

```
plotmd(x, class = NULL, xlab = "Variables", ylab = "Data", ...)
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>class</code>	class id for all observations.
<code>xlab</code>	label of x-axis.
<code>ylab</code>	label of y-axis.
<code>...</code>	other parameters to the plot.

**Details**

This a simplified parallel coordinate plot.

**Value**

A plot is returned.

**Author(s)**

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#), [emcluster](#).

**Examples**

```
## Not run:
library(EMCluster, quietly = TRUE)
set.seed(1234)

x <- as.matrix(iris[, 1:4], ncol = 4)
ret <- em.EM(x, nclass = 5)
plotmd(x, ret$class)

## End(Not run)
```

---

Plot Projection and Contour

*Plot Contour*

---

## Description

The function plots multivariate data on 2D plane with contour. Typically, the contour is built via projection pursuit or SVD algorithms, such as [project.on.2d\(\)](#).

## Usage

```
plotppcontour(da, Pi, Mu, S, class, class.true = NULL, n.grid = 128,  
              angle = 0, xlab = "", ylab = "", main = "")
```

## Arguments

da	a projected data matrix, dimension $n \times 2$ .
Pi	proportion, length $K$ .
Mu	the projected centers of cluster, dimension $K \times 2$ .
S	projected matrices of dispersion, dimension $p \times p \times K$ .
class	id of classifications, length $n$ .
class.true	ture id of classifications if available, length $n$ .
n.grid	number of grid points.
angle	a rotation angle (0 to $2\pi$ ).
xlab	an option for <code>plot()</code> function.
ylab	an option for <code>plot()</code> function.
main	an option for <code>plot()</code> function.

## Details

This function plots projection output of `project.on.2d()`.

da, Mu, and S are projected by some projection matrices obtained via SVD or projection pursuit algorithms. The projection is made on a 2D plane in the direction in which clusters of data x are most distinguishable to visualize.

## Value

A 2D projection plot is returned.

## Note

Only distinguishable for up to 7 clusters due to the limited color schemes.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[project.on.2d\(\)](#).

**Examples**

```
## Not run:
library(EMCluster, quietly = TRUE)
library(MASS, quietly = TRUE)
set.seed(1234)

### Crabs.
x <- as.matrix(crabs[, 4:8])
ret <- init.EM(x, nclass = 4, min.n = 20)
ret.proj <- project.on.2d(x, ret)

### Plot.
pdf("crabs_ppcontour.pdf", height = 5, width = 5)
plotppcontour(ret.proj$da, ret.proj$Pi, ret.proj$Mu, ret.proj$S,
              ret.proj$class, angle = pi/6, main = "Crabs K = 4")
dev.off()

## End(Not run)
```

---

Post I Information Functions

*Post I Information Functions of EMCluster*

---

**Description**

All post I information functions are for computing relative quantities and can be utilized by advanced developers with caution.

Currently, these are only for workflows.

**Author(s)**

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

## Description

Several classes are declared in **EMCluster**, and these are functions to print and summary objects.

## Usage

```
## S3 method for class 'emret'  
print(x, digits = max(4, getOption("digits") - 3), ...)  
## S3 method for class 'emret'  
summary(object, ...)  
## S3 method for class 'svd'  
summary(object, ...)
```

## Arguments

x	an object with the class attributes.
digits	for printing out numbers.
object	an object with the class attributes.
...	other possible options.

## Details

These are useful functions for summarizing and debugging.

## Value

The results will cat or print on the STDOUT by default.

## Author(s)

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

## References

<https://www.stat.iastate.edu/people/ranjan-maitra>

## See Also

[init.EM](#), [emcluster](#), [starts.via.svd](#).

**Examples**

```
## Not run:
library(EMCluster, quietly = TRUE)
x2 <- da2$da

emobj <- list(pi = da2$pi, Mu = da2$Mu, LTSigma = da2$LTSigma)
eobj <- e.step(x2, emobj = emobj)
emobj <- m.step(x2, emobj = eobj)
summary(emobj)

ret <- starts.via.svd(x2, nclass = 10, method = "kmeans")
summary(ret)

## End(Not run)
```

---

 Projection On 2D

*Produce Projection on 2D*


---

**Description**

The function projects multivariate data on 2D plane which can be displayed by [plotppcontour\(\)](#) later.

**Usage**

```
project.on.2d(x, emobj = NULL, pi = NULL, Mu = NULL,
             LTSigma = NULL, class = NULL, method = c("PP", "SVD"))
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>emobj</code>	the desired model which is a list mainly contains <code>pi</code> , <code>Mu</code> , and <code>LTSigma</code> , usually a returned object from <code>init.EM</code> .
<code>pi</code>	the mixing proportion, length $K$ .
<code>Mu</code>	the centers of clusters, dimension $K \times p$ .
<code>LTSigma</code>	the lower triangular matrices of dispersion, $K \times p(p + 1)/2$ .
<code>class</code>	id of classifications, length $n$ .
<code>method</code>	either projection pursuit or singular value decomposition.

**Details**

This function produces projection outputs of `x` and `emobj`.



**Value**

A projection is returned which is a list contains

- `da` is a  $n \times 2$  projected matrix of `x`.
- `Pi` is the original proportion `emobj$pi` of length  $K$ .
- `Mu` is a  $K \times 2$  projected matrix of `emobj$Mu`.
- `S` is a  $2 \times 2 \times K$  projected array of `emobj$Sigma`.
- `class` is the original class id `emobj$class`.
- `proj.mat` is the projection matrix of dimension  $p \times 2$ .

**Author(s)**

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[project.on.2d\(\)](#).

**Examples**

```
## Not run:
library(EMcluster, quietly = TRUE)
set.seed(1234)

### Iris.
x <- as.matrix(iris[, 1:4])
ret <- init.EM(x, nclass = 3, min.n = 30)
ret.proj <- project.on.2d(x, ret)

### Plot.
pdf("iris_ppcontour.pdf", height = 5, width = 5)
plotppcontour(ret.proj$da, ret.proj$Pi, ret.proj$Mu, ret.proj$S,
              ret.proj$class, main = "Iris K = 3")
dev.off()

## End(Not run)
```

---

Rand Index

*Rand Index and Adjusted Rand Index*

---

### Description

This function returns the Rand index and the adjusted Rand index for given true class ids and predicted class ids.

### Usage

```
RRand(trc1, prc1, lab = NULL)
```

### Arguments

trc1	true class ids.
prc1	predicted class ids.
lab	known ids for semi-supervised clustering.

### Details

All ids, trc1 and prc1, should be positive integers and started from 1 to K, and the maximums are allowed to be different.

lab used in semi-supervised clustering contains the labels which are known before clustering. It should be positive integer and started from 1 for labeled data and 0 for unlabeled data.

### Value

Return a Class RRand contains Rand index and adjusted Rand index.

### Author(s)

Wei-Chen Chen <wccsnow@gmail.com> and Ranjan Maitra.

### References

<https://www.stat.iastate.edu/people/ranjan-maitra>

### Examples

```
library(EMCluster, quietly = TRUE)

true.id <- c(1, 1, 1, 2, 2, 2, 3, 3, 3)
pred.id <- c(2, 1, 2, 1, 1, 1, 2, 1, 1)
label   <- c(0, 0, 0, 0, 1, 0, 2, 0, 0)

RRand(true.id, pred.id)
RRand(true.id, pred.id, lab = label)
```

---

Recolor Classification IDs  
*Recolor Classification IDs*

---

**Description**

These functions return new classification IDs.

**Usage**

```
recolor(id.target, id.class, scatter.class = NULL, scatter.target = NULL)
rematch(tg.id, cl.id)
recode(id)
```

**Arguments**

<code>id.target</code>	target class ids.
<code>id.class</code>	original class ids.
<code>scatter.class</code>	scatter class ids.
<code>scatter.target</code>	scatter target class ids.
<code>id</code>	class ids.
<code>tg.id</code>	target class ids.
<code>cl.id</code>	class ids.

**Details**

The function `recolor` colors `id.target` in accordance with the most likely candidate in `id.class`. Note that if `scatter` is present, then the class given by 0 is represented as `scatter` and it is assumed to be the same for both classifications.

The function `rematch` returns a list as `id.trcl` and `id.prcl`. It is the heart of the `recolor` function and is usually called from `recolor`.

The function `recode` reorders classes to eliminate group ids without any members. It is assumed that the group ids are integers.

**Value**

See Details.

**Author(s)**

Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**Examples**

```
## Not run:
library(EMCluster, quietly = TRUE)

true.id <- c(1, 1, 1, 2, 2, 2, 3, 3, 3)
pred.id <- c(2, 1, 2, 1, 1, 1, 2, 1, 1)

recolor(pred.id, true.id)

## End(Not run)
```

---

 Single Step

*Single E- and M-step*


---

**Description**

These functions are single E- and M-step of EM algorithm for model-based clustering of finite mixture multivariate Gaussian distribution with unstructured dispersion.

**Usage**

```
e.step(x, emobj = NULL, pi = NULL, Mu = NULL, LTSigma = NULL,
       norm = TRUE)
m.step(x, emobj = NULL, Gamma = NULL, assign.class = FALSE)
```

**Arguments**

<code>x</code>	the data matrix, dimension $n \times p$ .
<code>emobj</code>	the desired model which is a list mainly contains <code>pi</code> , <code>Mu</code> , and <code>LTSigma</code> , usually a returned object from <code>init.EM</code> .
<code>pi</code>	the mixing proportion, length $K$ .
<code>Mu</code>	the centers of clusters, dimension $K \times p$ .
<code>LTSigma</code>	the lower triangular matrices of dispersion, $K \times p(p+1)/2$ .
<code>norm</code>	if returning normalized Gamma.
<code>Gamma</code>	containing posterior probabilities if normalized, otherwise containing component densities weighted by mixing proportion, dimension $n \times K$ .
<code>assign.class</code>	if assigning class id.

**Details**

These two functions are mainly used in debugging for development and post process after model fitting.

**Value**

The `e.step` returns a list contains Gamma, the posterior probabilities if `norm=TRUE`, otherwise it contains component densities. This is one E-step and Gamma is used to update `emobj` in the M-step next.

The `m.step` returns a new `emobj` according to the Gamma from the E-step above.

**Author(s)**

Wei-Chen Chen <[wccsnow@gmail.com](mailto:wccsnow@gmail.com)> and Ranjan Maitra.

**References**

<https://www.stat.iastate.edu/people/ranjan-maitra>

**See Also**

[init.EM](#).

**Examples**

```
library(EMcluster, quietly = TRUE)
x2 <- da2$da

emobj <- list(pi = da2$pi, Mu = da2$Mu, LTSigma = da2$LTSigma)
eobj <- e.step(x2, emobj = emobj)
emobj <- m.step(x2, emobj = eobj)
emobj
```

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