

# Package ‘fmx’

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**Type** Package

**Title** Finite Mixture Parametrization

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**Description** A parametrization framework for finite mixture distribution using S4 objects. Density, cumulative density, quantile and simulation functions are defined. Currently normal, Tukey g-&-h, skew-normal and skew-t distributions are well tested. The gamma, negative binomial distributions are being tested.

**License** GPL-2

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**Description**

A parametrization framework for finite mixture distribution using S4 objects.

Density, cumulative density, quantile and simulation functions are defined.

Currently normal, Tukey *g*-&-*h*, skew-normal and skew-*t* distributions are well tested. The gamma, negative binomial distributions are being tested.

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---

approxdens

*Empirical Density Function*

---

**Description**

..

**Usage**

```
approxdens(x, ...)
```

**Arguments**

`x`                    [numeric vector](#), observations  
`...`                additional parameters of [density.default](#)

**Details**

[approx](#) inside [density.default](#)

another 'layer' of [approxfun](#)

**Value**

Function [approxdens](#) returns a [function](#).

**Examples**

```
x = rnorm(1e3L)
f = approxdens(x)
f(x[1:3])
```

---

|        |  |
|--------|--|
| as.fmx | <i>Turn Various Objects to <a href="#">fmx</a> Class</i> |
|--------|--|

---

**Description**

Turn various objects created in other R packages to [fmx](#) class.

**Usage**

```
as.fmx(x, ...)
```

**Arguments**

|     |   |
|-----|---|
| x   | an R object   |
| ... | additional parameters, see <b>Arguments</b> in individual S3 dispatches |

**Details**

Various mixture distribution estimates obtained from other R packages are converted to [fmx](#) class, so that we could take advantage of all methods defined for [fmx](#) objects.

**Value**

S3 generic function [as.fmx](#) returns an [fmx](#) object.

---

|                |   |
|----------------|---|
| as.fmx.fitdist | <i>Convert <a href="#">fitdist</a> Objects to <a href="#">fmx</a> Class</i> |
|----------------|---|

---

**Description**

To convert [fitdist](#) objects (from package [fitdistrplus](#)) to [fmx](#) class.

**Usage**

```
## S3 method for class 'fitdist'
as.fmx(x, ...)
```

**Arguments**

|     |                                |
|-----|--------------------------------|
| x   | <a href="#">fitdist</a> object |
| ... | ..                             |

**Value**

Function [as.fmx.fitdist](#) returns an [fmx](#) object.

**Examples**

```
library(fitdistrplus)
# ?fitdist
data(endosulfan, package = 'fitdistrplus')
ATV <- subset(endosulfan, group == 'NonArthroInvert')$ATV
log10ATV <- log10(ATV)
fln <- fitdist(log10ATV, distr = 'norm')
(fln2 <- as.fmx(fln))
hist.default(log10ATV, freq = FALSE)
curve(dfmx(x, dist = fln2), xlim = range(log10ATV), add = TRUE)
```

as.fmx.mixEM

*Convert mixEM Objects to [fmx](#) Class***Description**

To convert mixEM objects (from package [mixtools](#)) to [fmx](#) class.

Currently only the returned value of [normalmixEM](#) and [gammamixEM](#) are supported

**Usage**

```
## S3 method for class 'mixEM'
as.fmx(x, data = x[["x"]], ...)
```

**Arguments**

|      |                                |
|------|--------------------------------|
| x    | mixEM object                   |
| data | <a href="#">numeric vector</a> |
| ...  | ..                             |

**Value**

Function [as.fmx.mixEM](#) returns an [fmx](#) object.

**Note**

[plot.mixEM](#) not plot [gammamixEM](#) returns, as of 2022-09-19.

**Examples**

```
library(mixtools)
(wait = as.fmx(normalmixEM(faithful$waiting, k = 2)))
hist.default(faithful$waiting, freq = FALSE)
curve(dfmx(x, dist = wait), xlim = range(faithful$waiting), add = TRUE)
```

---

|               |   |
|---------------|---|
| as.fmx.Normal | <i>Convert</i> Normal <i>fit</i> <i>from</i> <a href="https://CRAN.R-project.org/package=mixsmsn">Rhrefhttps://CRAN.R-project.org/package=mixsmsn</a> to <i>fmx</i> |
|---------------|---|

---

## Description

To convert Normal object (from package **mixsmsn**) to **fmx** class.

## Usage

```
## S3 method for class 'Normal'
as.fmx(x, data, ...)
```

## Arguments

|      |  |
|------|--|
| x    | 'Normal' object, returned from <a href="#">smsn.mix</a> with parameter family = 'Normal' |
| data | <a href="#">numeric vector</a>   |
| ...  | additional parameters, currently not in use  |

## Value

Function [as.fmx.Normal](#) returns an **fmx** object.

## Note

[smsn.mix](#) does not offer a parameter to keep the input data, as of 2021-10-06.

## Examples

```
library(mixsmsn)
# ?smsn.mix
arg1 = c(mu = 5, sigma2 = 9, lambda = 5, nu = 5)
arg2 = c(mu = 20, sigma2 = 16, lambda = -3, nu = 5)
arg3 = c(mu = 35, sigma2 = 9, lambda = -6, nu = 5)
set.seed(120); x = rmix(n = 1e3L, p=c(.5, .2, .3), family = 'Skew.t',
  arg = list(unnname(arg1), unnname(arg2), unnname(arg3)))

# Normal
class(m2 <- smsn.mix(x, nu = 3, g = 3, family = 'Normal', calc.im = FALSE))
mix.hist(y = x, model = m2)
m2a = as.fmx(m2, data = x)
hist(x, freq = FALSE)
curve(dfmx(x, dist = m2a), xlim = range(x), add = TRUE)
```

---

as.fmx.Skew.normal      *Convert Skew.normal Object to fmx*

---

## Description

To convert Skew.normal object (from package **mixsmsn**) to **fmx** class.

## Usage

```
## S3 method for class 'Skew.normal'  
as.fmx(x, data, ...)
```

## Arguments

|      |   |
|------|---|
| x    | 'Skew.normal' object, returned from <a href="#">smsn.mix</a> with parameter family = 'Skew.normal'. |
| data | <a href="#">numeric vector</a>  |
| ...  | additional parameters, currently not in use   |

## Value

Function [as.fmx.Skew.normal](#) returns an **fmx** object.

## Note

[smsn.mix](#) does not offer a parameter to keep the input data, as of 2021-10-06.

## Examples

```
library(mixsmsn)  
# ?smsn.mix  
arg1 = c(mu = 5, sigma2 = 9, lambda = 5, nu = 5)  
arg2 = c(mu = 20, sigma2 = 16, lambda = -3, nu = 5)  
arg3 = c(mu = 35, sigma2 = 9, lambda = -6, nu = 5)  
set.seed(120); x = rmix(n = 1e3L, p=c(.5, .2, .3), family = 'Skew.t',  
  arg = list(unnamed(arg1), unnamed(arg2), unnamed(arg3)))  
  
# Skew Normal  
class(m1 <- smsn.mix(x, nu = 3, g = 3, family = 'Skew.normal', calc.im = FALSE))  
mix.hist(y = x, model = m1)  
m1a = as.fmx(m1, data = x)  
(l1a = logLik(m1a))  
hist(x, freq = FALSE)  
curve(dfmx(x, dist = m1a), xlim = range(x), add = TRUE)
```

---

|               |  |
|---------------|--|
| as.fmx.Skew.t | <i>Convert</i> Skew.t <i>fit</i> <i>from</i> R<br><a href="https://CRAN.R-project.org/package=mixsmsn">https://CRAN.R-project.org/package=mixsmsn</a> <i>to</i> <i>fmX</i> |
|---------------|--|

---

## Description

To convert Skew.t object (from package **mixsmsn**) to **fmX** class.

## Usage

```
## S3 method for class 'Skew.t'
as.fmx(x, data, ...)
```

## Arguments

|      |  |
|------|--|
| x    | 'Skew.t' object, returned from <a href="#">smsn.mix</a> with parameter family = 'Skew.t' |
| data | <a href="#">numeric vector</a>   |
| ...  | additional parameters, currently not in use  |

## Value

Function [as.fmx.Skew.t](#) returns an **fmX** object.

## Note

[smsn.mix](#) does not offer a parameter to keep the input data, as of 2021-10-06.

## Examples

```
# mixsmsn::smsn.mix with option `family = 'Skew.t'` is slow

library(mixsmsn)
# ?smsn.mix
arg1 = c(mu = 5, sigma2 = 9, lambda = 5, nu = 5)
arg2 = c(mu = 20, sigma2 = 16, lambda = -3, nu = 5)
arg3 = c(mu = 35, sigma2 = 9, lambda = -6, nu = 5)
set.seed(120); x = rmix(n = 1e3L, p=c(.5, .2, .3), family = 'Skew.t',
  arg = list(unnamed(arg1), unnamed(arg2), unnamed(arg3)))

# Skew t
class(m3 <- smsn.mix(x, nu = 3, g = 3, family = 'Skew.t', calc.im = FALSE))
mix.hist(y = x, model = m3)
m3a = as.fmx(m3, data = x)
hist(x, freq = FALSE)
curve(dfmx(x, dist = m3a), xlim = range(x), add = TRUE)
(l3a = logLik(m3a))
stopifnot(all.equal.numeric(AIC(l3a), m3$aaic), all.equal.numeric(BIC(l3a), m3$bic))
```



---

|          |                           |  |
|----------|---------------------------|--|
| as.fmx.t | <i>Convert t fit from</i> | <a href="https://CRAN.R-project.org/package=mixsmsn">Rhrefhttps://CRAN.R-project.org/package=mixsmsn</a> to <i>fmx</i> |
|----------|---------------------------|--|

---

## Description

To convert t object (from package **mixsmsn**) to **fmx** class.

## Usage

```
## S3 method for class 't'
as.fmx(x, data, ...)
```

## Arguments

|      |  |
|------|--|
| x    | 't' object, returned from <a href="#">smsn.mix</a> with parameter family = 't' |
| data | <a href="#">numeric vector</a>   |
| ...  | additional parameters, currently not in use                                    |

## Value

Function [as.fmx.t](#) has not been completed yet

## Note

[smsn.mix](#) does not offer a parameter to keep the input data, as of 2021-10-06.

## Examples

```
library(mixsmsn)
# ?smsn.mix
arg1 = c(mu = 5, sigma2 = 9, lambda = 5, nu = 5)
arg2 = c(mu = 20, sigma2 = 16, lambda = -3, nu = 5)
arg3 = c(mu = 35, sigma2 = 9, lambda = -6, nu = 5)
set.seed(120); x = rmix(n = 1e3L, p=c(.5, .2, .3), family = 'Skew.t',
  arg = list(unnamed(arg1), unnamed(arg2), unnamed(arg3)))

# t
class(m4 <- smsn.mix(x, nu = 3, g = 3, family = 't', calc.im = FALSE))
mix.hist(y = x, model = m4)
# as.fmx(m4, data = x) # not ready yet!!
```

---

 coef.fmx

*Parameter Estimates of [fmx](#) object*


---

**Description**

..

**Usage**

```
## S3 method for class 'fmx'
coef(object, internal = FALSE, ...)
```

**Arguments**

|          |   |
|----------|---|
| object   | <a href="#">fmx</a> object  |
| internal | <b>logical</b> scalar, either for the user-friendly parameters (FALSE, default) (e.g., mean, sd for normal mixture, and A, B, g, h for Tukey <i>g</i> -and- <i>h</i> mixture), or for the internal/unconstrained parameters (TRUE). |
| ...      | place holder for S3 naming convention   |

**Details**

Function [coef.fmx](#) returns the estimates of the user-friendly parameters (parm = 'user'), or the internal/unconstrained parameters (parm = 'internal'). When the distribution has constraints on one or more parameters, function [coef.fmx](#) does not return the estimates (which is constant 0) of the constrained parameters.

**Value**

Function [coef.fmx](#) returns a **numeric vector**.

---

 confint.fmx

*Confidence Interval of [fmx](#) Object*


---

**Description**

...

**Usage**

```
## S3 method for class 'fmx'
confint(object, ..., level = 0.95)
```

**Arguments**

object            [fmx](#) object  
 ...                place holder for S3 naming convention  
 level             confidence level, default 95%.

**Details**

[confint.fmx](#) returns the Wald-type confidence intervals based on the user-friendly parameters (`parm = 'user'`), or the internal/unconstrained parameters (`parm = 'internal'`). When the distribution has constraints on one or more parameters, function [confint.fmx](#) does not return the confident intervals of for the constrained parameters.

**Value**

[confint.fmx](#) returns a [matrix](#)

---

|         |   |
|---------|---|
| dbl2fmx | <i>Inverse of <a href="#">fmx2dbl</a>, for internal use</i> |
|---------|---|

---

**Description**

..

**Usage**

```
dbl2fmx(x, K, distname, ...)
```

**Arguments**

x                    [numeric vector](#), unrestricted parameters  
 K                    [integer](#) scalar  
 distname            [character](#) scalar  
 ...                  additional parameters, not currently used

**Details**

Only used in downstream function `QuantileGH::QLMDe` and unexported function `QuantileGH:::qfmx_gr`, not compute intensive.

**Value**

Function [dbl2fmx](#) returns a [list](#) with two elements `$pars` and `$w`

---

`dfmx`*Density, Distribution and Quantile of Finite Mixture Distribution*

---

**Description**

Density function, distribution function, quantile function and random generation for a finite mixture distribution with normal or Tukey  $g$ -&- $h$  components.

**Usage**

```
dfmx(  
  x,  
  dist,  
  distname = dist@distname,  
  K = dim(pars)[1L],  
  pars = dist@pars,  
  w = dist@w,  
  ...,  
  log = FALSE  
)  
  
pfmx(  
  q,  
  dist,  
  distname = dist@distname,  
  K = dim(pars)[1L],  
  pars = dist@pars,  
  w = dist@w,  
  ...,  
  lower.tail = TRUE,  
  log.p = FALSE  
)  
  
qfmx(  
  p,  
  dist,  
  distname = dist@distname,  
  K = dim(pars)[1L],  
  pars = dist@pars,  
  w = dist@w,  
  interval = qfmx_interval(dist = dist),  
  ...,  
  lower.tail = TRUE,  
  log.p = FALSE  
)  
  
rfmx(  
  n,  
  dist,  
  distname = dist@distname,  
  K = dim(pars)[1L],  
  pars = dist@pars,  
  w = dist@w,  
  interval = qfmx_interval(dist = dist),  
  ...,  
  lower.tail = TRUE,  
  log.p = FALSE  
)
```

```

n,
dist,
distname = dist@distname,
K = dim(pars)[1L],
pars = dist@pars,
w = dist@w
)

```

### Arguments

|                                   |   |
|-----------------------------------|---|
| <code>x, q</code>                 | <b>numeric vector</b> , quantiles, NA_real_ value(s) allowed.   |
| <code>dist</code>                 | <b>fmx</b> object, a finite mixture distribution  |
| <code>distname, K, pars, w</code> | auxiliary parameters, whose default values are determined by argument <code>dist</code> . The user-specified <b>vector</b> of <code>w</code> does not need to sum up to 1; $w/\text{sum}(w)$ will be used internally. |
| <code>...</code>                  | additional parameters   |
| <code>log, log.p</code>           | <b>logical</b> scalar. If TRUE, probabilities are given as $\log(p)$ .  |
| <code>lower.tail</code>           | <b>logical</b> scalar. If TRUE (default), probabilities are $Pr(X \leq x)$ , otherwise, $Pr(X > x)$ .   |
| <code>p</code>                    | <b>numeric vector</b> , probabilities.  |
| <code>interval</code>             | length two <b>numeric vector</b> , interval for root finding, see <a href="#">vuniroot2</a> and <a href="#">vuniroot</a>  |
| <code>n</code>                    | <b>integer</b> scalar, number of observations.  |

### Details

A computational challenge in function `dfmx` is when mixture density is very close to 0, which happens when the per-component log densities are negative with big absolute values. In such case, we cannot compute the log mixture densities (i.e., `-Inf`), for the log-likelihood using function `logLik.fmx`. Our solution is to replace these `-Inf` log mixture densities by the weighted average (using the mixing proportions of `dist`) of the per-component log densities.

Function `qfmx` gives the quantile function, by numerically solving `pfmx`. One major challenge when dealing with the finite mixture of Tukey *g*-&-*h* family distribution is that Brent–Dekker’s method needs to be performed in both `pGH` and `qfmx` functions, i.e. *two layers* of root-finding algorithm.

### Value

Function `dfmx` returns a **numeric vector** of probability density values of an `fmx` object at specified quantiles `x`.

Function `pfmx` returns a **numeric vector** of cumulative probability values of an `fmx` object at specified quantiles `q`.

Function `qfmx` returns an unnamed **numeric vector** of quantiles of an `fmx` object, based on specified cumulative probabilities `p`.

Function `rfmx` generates random deviates of an `fmx` object.

**Note**

Function `qnorm` returns an unnamed **vector** of quantiles, although `quantile` returns a named **vector** of quantiles.

**Examples**

```
library(ggplot2)

(e1 = fmx('norm', mean = c(0,3), sd = c(1,1.3), w = c(1, 1)))
curve(dfmx(x, dist = e1), xlim = c(-3,7))
ggplot() + geom_function(fun = dfmx, args = list(dist = e1)) + xlim(-3,7)
ggplot() + geom_function(fun = pfmx, args = list(dist = e1)) + xlim(-3,7)
hist(rfmx(n = 1e3L, dist = e1), main = '1000 obs from e1')

x = (-3):7
round(dfmx(x, dist = e1), digits = 3L)
round(p1 <- pfmx(x, dist = e1), digits = 3L)
stopifnot(all.equal.numeric(qfmx(p1, dist = e1), x, tol = 1e-4))

(e2 = fmx('GH', A = c(0,3), g = c(.2, .3), h = c(.2, .1), w = c(2, 3)))
ggplot() + geom_function(fun = dfmx, args = list(dist = e2)) + xlim(-3,7)

round(dfmx(x, dist = e2), digits = 3L)
round(p2 <- pfmx(x, dist = e2), digits = 3L)
stopifnot(all.equal.numeric(qfmx(p2, dist = e2), x, tol = 1e-4))

(e3 = fmx('GH', g = .2, h = .01)) # one-component Tukey
ggplot() + geom_function(fun = dfmx, args = list(dist = e3)) + xlim(-3,5)
set.seed(124); r1 = rfmx(1e3L, dist = e3);
set.seed(124); r2 = TukeyGH77::rGH(n = 1e3L, g = .2, h = .01)
stopifnot(identical(r1, r2)) # but ?rfmx has much cleaner code
round(dfmx(x, dist = e3), digits = 3L)
round(p3 <- pfmx(x, dist = e3), digits = 3L)
stopifnot(all.equal.numeric(qfmx(p3, dist = e3), x, tol = 1e-4))

a1 = fmx('GH', A = c(7,9), B = c(.8, 1.2), g = c(.3, 0), h = c(0, .1), w = c(1, 1))
a2 = fmx('GH', A = c(6,9), B = c(.8, 1.2), g = c(-.3, 0), h = c(.2, .1), w = c(4, 6))
library(ggplot2)
(p = ggplot() +
  geom_function(fun = pfmx, args = list(dist = a1), mapping = aes(color = 'g2=h1=0')) +
  geom_function(fun = pfmx, args = list(dist = a2), mapping = aes(color = 'g2=0')) +
  xlim(3,15) +
  scale_y_continuous(labels = scales::percent) +
  labs(y = NULL, color = 'models') +
  coord_flip())
p + theme(legend.position = 'none')

# to use [rfmx] without \pkg{fmx}
(d = fmx(distname = 'GH', A = c(-1,1), B = c(.9,1.1), g = c(.3,-.2), h = c(.1,.05), w = c(2,3)))
d@pars
```

```

set.seed(14123); x = rfm(x = 1e3L, dist = d)
set.seed(14123); x_raw = rfm(x = 1e3L,
  distname = 'GH', K = 2L,
  pars = rbind(
    c(A = -1, B = .9, g = .3, h = .1),
    c(A = 1, B = 1.1, g = -.2, h = .05)
  ),
  w = c(.4, .6)
)
stopifnot(identical(x, x_raw))

```

---

|          |  |
|----------|--|
| distArgs | <i>Name(s) of Formal Argument(s) of Distribution</i> |
|----------|--|

---

### Description

To obtain the name(s) of distribution parameter(s).

### Usage

```
distArgs(distname)
```

### Arguments

distname      [character](#) scalar, name of distribution

### Value

Function [distArgs](#) returns a [character vector](#).

### See Also

[formalArgs](#)

---

|          |                          |
|----------|--------------------------|
| distType | <i>Distribution Type</i> |
|----------|--------------------------|

---

### Description

..

### Usage

```
distType(type = c("discrete", "nonNegContinuous", "continuous"))
```

**Arguments**

type            [character](#) scalar

**Value**

Function [distType](#) returns a [character vector](#).

---

|               |  |
|---------------|--|
| dist_logtrans | <i>Distribution Parameters that needs to have a log-transformation</i> |
|---------------|--|

---

**Description**

..

**Usage**

```
dist_logtrans(distname)
```

**Arguments**

distname        [character](#) scalar, name of distribution

**Value**

Function [dist\\_logtrans](#) returns an [integer](#) scalar

---

|     |  |
|-----|--|
| fmx | <i>Create fmx Object for Finite Mixture Distribution</i> |
|-----|--|

---

**Description**

To create [fmx](#) object for finite mixture distribution.

**Usage**

```
fmx(distname, w = 1, ...)
```

**Arguments**

distname        [character](#) scalar

w                (optional) [numeric vector](#). Does not need to sum up to 1; w/sum(w) will be used internally.

...              mixture distribution parameters. See function [dGH](#) for the names and default values of Tukey *g-&-h* distribution parameters, or [dnorm](#) for the names and default values of normal distribution parameters.



**Value**

Function `fmx` returns an `fmx` object.

**Examples**

```
(e1 = fmx('norm', mean = c(0,3), sd = c(1,1.3), w = c(1, 1)))
isS4(e1) # TRUE
slotNames(e1)

(e2 = fmx('GH', A = c(0,3), g = c(.2, .3), h = c(.2, .1), w = c(2, 3)))

(e3 = fmx('GH', A = 0, g = .2, h = .2)) # one-component Tukey
```

---

fmx-class

*fmx Class: Finite Mixture Parametrization*


---

**Description**

An S4 object to specify the parameters and type of distribution of a one-dimensional finite mixture distribution.

**Slots**

`distname` **character** scalar, name of parametric distribution of the mixture components. Currently, normal ('norm') and Tukey *g*-&-*h* ('GH') distributions are supported.

`pars` **double matrix**, all distribution parameters in the mixture. Each row corresponds to one component. Each column includes the same parameters of all components. The order of rows corresponds to the (non-strictly) increasing order of the component location parameters. The columns match the formal arguments of the corresponding distribution, e.g., 'mean' and 'sd' for **normal** mixture, or 'A', 'B', 'g' and 'h' for Tukey *g*-&-*h* mixture.

`w` **numeric vector** of mixing proportions that must sum to 1

`data` (optional) **numeric vector**, the one-dimensional observations

`data.name` (optional) **character** scalar, a human-friendly name of the observations

`vcov_internal` (optional) variance-covariance **matrix** of the internal (i.e., unconstrained) estimates

`vcov` (optional) variance-covariance **matrix** of the mixture distribution (i.e., constrained) estimates

`Kolmogorov, CramerVonMises, KullbackLeibler` (optional) **numeric** scalars

fmx2dbl

Reparameterization of *fmx* Object**Description**

To convert the parameters of *fmx* object into unrestricted parameters.

**Usage**

```
fmx2dbl(
  x,
  distname = x@distname,
  pars = x@pars,
  K = dim(pars)[1L],
  w = x@w,
  ...
)
```

**Arguments**

|          |   |
|----------|---|
| x        | <i>fmx</i> object                         |
| distname | character scalar, default x@distname      |
| pars     | numeric matrix, default x@pars            |
| K        | integer scalar, default value from x      |
| w        | numeric vector, default x@w               |
| ...      | additional parameters, not currently used |

**Details**

For the first parameter

- $A_1 \rightarrow A_1$
- $A_2 \rightarrow A_1 + \exp(\log(d_1))$
- $A_k \rightarrow A_1 + \exp(\log(d_1)) + \dots + \exp(\log(d_{k-1}))$

For mixing proportions to multinomial logits.

For 'norm': sd -> log(sd) for 'GH': B -> log(B), h -> log(h)

**Value**

Function *fmx2dbl* returns a numeric vector.

**See Also**

[dbl2fmx](#)

---

|                |  |
|----------------|--|
| fmx_constraint | <i>Parameter Constraint(s) of Mixture Distribution</i> |
|----------------|--|

---

### Description

Determine the parameter constraint(s) of a finite mixture distribution `fmx`, either by the value of parameters of such mixture distribution, or by a user-specified string.

### Usage

```
fmx_constraint(
  dist,
  distname = dist@distname,
  K = dim(dist@pars)[1L],
  pars = dist@pars
)
```

### Arguments

|                       |  |
|-----------------------|--|
| <code>dist</code>     | (optional) <code>fmx</code> object   |
| <code>distname</code> | <code>character</code> scalar, name of distribution (see <code>fmx</code> ), default value determined by <code>dist</code>                                   |
| <code>K</code>        | <code>integer</code> scalar, number of components, default value determined by <code>dist</code>   |
| <code>pars</code>     | <code>double matrix</code> , distribution parameters of a finite mixture distribution (see <code>fmx</code> ), default value determined by <code>dist</code> |

### Value

Function `fmx_constraint` returns the indices of internal parameters (only applicable to Tukey *g*-&-*h* mixture distribution, yet) to be constrained, based on the input `fmx` object `dist`.

### Examples

```
(d0 = fmx('GH', A = c(1,4), g = c(.2,.1), h = c(.05,.1), w = c(1,1)))
(c0 = fmx_constraint(d0))
user_constraint(character(), distname = 'GH', K = 2L) # equivalent

(d1 = fmx('GH', A = c(1,4), g = c(.2,0), h = c(0,.1), w = c(1,1)))
(c1 = fmx_constraint(d1))
user_constraint(c('g2', 'h1'), distname = 'GH', K = 2L) # equivalent

(d2 = fmx('GH', A = c(1,4), g = c(.2,0), h = c(.15,.1), w = c(1,1)))
(c2 = fmx_constraint(d2))
user_constraint('g2', distname = 'GH', K = 2L) # equivalent
```

---

|               |                                    |
|---------------|------------------------------------|
| fmx_diagnosis | <i>Diagnoses for fmx Estimates</i> |
|---------------|------------------------------------|

---

### Description

Diagnoses for [fmx](#) estimates.

### Usage

```
Kolmogorov_fmx(object, data = object@data, ...)
```

```
KullbackLeibler_fmx(object, data = object@data, ...)
```

```
CramerVonMises_fmx(object, data = object@data, ...)
```

### Arguments

|        |  |
|--------|--|
| object | <a href="#">fmx</a> object, or an R object convertible to an <a href="#">fmx</a> object                            |
| data   | <a href="#">double vector</a> , observed data. Default is <code>object@data</code> , the data used for estimation. |
| ...    | additional parameters, currently not in use  |

### Details

Function [Kolmogorov\\_fmx](#) calculates Kolmogorov distance.

Function [KullbackLeibler\\_fmx](#) calculates Kullback-Leibler divergence. The R code is adapted from `LaplacesDemon::KLD`.

Function [CramerVonMises\\_fmx](#) calculates Cramer-von Mises quadratic distance (via [cvm.test](#)).

### Value

Functions [Kolmogorov\\_fmx](#), [KullbackLeibler\\_fmx](#), [CramerVonMises\\_fmx](#) all return [numeric](#) scalars.

### See Also

`dgof::cvmf.test`

---

|        |   |
|--------|---|
| getTeX | <i>TeX Label (of Parameter Constraint(s)) of fmx Object</i> |
|--------|---|

---

**Description**

Create TeX label of (parameter constraint(s)) of [fmx](#) object

**Usage**

```
getTeX(dist, print_K = FALSE)
```

**Arguments**

`dist`            [fmx](#) object  
`print_K`        [logical](#) scalar, whether to print the number of components  $K$ . Default FALSE.

**Value**

Function [getTeX](#) returns a [character](#) scalar (of TeX expression) of the constraint, primarily intended for end-users in plots.

**Examples**

```
(d0 = fmx('GH', A = c(1,4), g = c(.2,.1), h = c(.05,.1), w = c(1,1)))
getTeX(d0)
```

```
(d1 = fmx('GH', A = c(1,4), g = c(.2,0), h = c(0,.1), w = c(1,1)))
getTeX(d1)
```

```
(d2 = fmx('GH', A = c(1,4), g = c(.2,0), h = c(.15,.1), w = c(1,1)))
getTeX(d2)
```

---

|                 |                                       |
|-----------------|---------------------------------------|
| Kolmogorov_dist | <i>One-Sample Kolmogorov Distance</i> |
|-----------------|---------------------------------------|

---

**Description**

To calculate the one-sample Kolmogorov distance between observations and a distribution.

**Usage**

```
Kolmogorov_dist(x, null, alternative = c("two.sided", "less", "greater"), ...)
```

**Arguments**

|             |  |
|-------------|--|
| x           | numeric vector, observations $x$   |
| null        | cumulative distribution function   |
| alternative | character scalar, alternative hypothesis, either 'two.sided' (default), 'less', or 'greater' |
| ...         | additional arguments of null   |

**Details**

Function `Kolmogorov_dist` is different from `ks.test` in the following aspects

- Ties in observations are supported. The step function of empirical distribution is inspired by `ecdf`. This is superior than  $(0 : (n - 1)) / n$  in `ks.test`.
- Discrete distribution (with discrete observation) is supported.
- Discrete distribution (with continuous observation) is not supported yet. This will be an easy modification in future.
- Only the one-sample Kolmogorov distance, not the one-sample Kolmogorov test, is returned, to speed up the calculation.

**Value**

Function `Kolmogorov_dist` returns a numeric scalar.

**Examples**

```
# from ?stats::ks.test
x1 = rnorm(50)
ks.test(x1+2, y = pgamma, shape = 3, rate = 2)
Kolmogorov_dist(x1+2, null = pgamma, shape = 3, rate = 2) # exactly the same

# discrete distribution
x2 <- rbinom(n = 1e2L, size = 500, prob = .4)
suppressWarnings(ks.test(x2, y = pnbinom, size = 500, prob = .4)) # warning on ties
Kolmogorov_dist(x2, null = pnbinom, size = 500, prob = .4) # wont be the same
```

---

logLik.fitdist

*Log-Likelihood of fitdist Object*


---

**Description**

..

**Usage**

```
## S3 method for class 'fitdist'
logLik(object, ...)
```

**Arguments**

object            [fitdist](#) object  
 ...                additional parameters, currently not in use

**Details**

Output of [fitdist](#) has elements \$loglik, \$aic and \$bic, but they are simply [numeric](#) scalars. `fitdistrplus:::logLik.fitdist` simply returns these elements.

**Value**

Function [logLik.fitdist](#) returns a [logLik](#) object, which could be further used by [AIC](#) and [BIC](#).  
 (I have written to the authors)

---

|            |   |
|------------|---|
| logLik.fmx | <i>Log-Likelihood of <a href="#">fmx</a> Object</i> |
|------------|---|

---

**Description**

..

**Usage**

```
## S3 method for class 'fmx'
logLik(object, data = object@data, ...)
```

**Arguments**

object            [fmx](#) object  
 data              [double vector](#), actual observations  
 ...                place holder for S3 naming convention

**Details**

Function [logLik.fmx](#) returns a [logLik](#) object indicating the log-likelihood. An additional attribute `attr(, 'logl')` indicates the point-wise log-likelihood, to be use in Vuong's closeness test.

**Value**

Function [logLik.fmx](#) returns a [logLik](#) object with an additional attribute `attr(, 'logl')`.

---

|              |   |
|--------------|---|
| logLik.mixEM | <i>Log-Likelihood of 'mixEM' Object</i> |
|--------------|---|

---

**Description**

To obtain the log-Likelihood of 'mixEM' object, based on [mixtools](#) 2020-02-05.

**Usage**

```
## S3 method for class 'mixEM'
logLik(object, ...)
```

**Arguments**

|        |  |
|--------|--|
| object | 'mixEM' object, currently only the returned value of <a href="#">normalmixEM</a> and <a href="#">gam-mamixEM</a> are supported |
| ...    | additional parameters, currently not in use  |

**Value**

Function [logLik.mixEM](#) returns a [logLik](#) object.

---

|     |  |
|-----|--|
| MaP | <i>Maximum a Posteriori clustering</i> |
|-----|--|

---

**Description**

..

**Usage**

```
MaP(x, dist, ...)
```

**Arguments**

|      |                               |
|------|-------------------------------|
| x    | numeric vector                |
| dist | an <a href="#">fmx</a> object |
| ...  | ..                            |

**Value**

Function [MaP](#) returns an [integer vector](#).



**Examples**

```
x = rnorm(1e2L, sd = 2)
m = fmx('norm', mean = c(-1.5, 1.5), w = c(1, 2))
library(ggplot2)
ggplot() + geom_function(fun = dfmx, args = list(dist = m)) +
  geom_point(mapping = aes(x = x, y = .05, color = factor(MaP(x, dist = m)))) +
  labs(color = 'Maximum a Posteriori\nClustering')
```

---

|            |   |
|------------|---|
| mixEM_pars | <i>Names of Distribution Parameters of 'mixEM' Object</i> |
|------------|---|

---

**Description**

Names of distribution parameters of 'mixEM' object, based on [mixtools](#) 2020-02-05.

**Usage**

```
mixEM_pars(object)
```

**Arguments**

object            'mixEM' object, currently only the returned value of [normalmixEM](#) and [gammamixEM](#) are supported

**Value**

Function [mixEM\\_pars](#) returns a [character vector](#)

**See Also**

[normalmixEM](#) [gammamixEM](#)

---

|        |   |
|--------|---|
| mlogis | <i>Multinomial Probabilities &amp; Logits</i> |
|--------|---|

---

**Description**

Performs transformation between [vectors](#) of multinomial probabilities and multinomial logits.

This transformation is a generalization of [plogis](#) which converts scalar logit into probability and [qlogis](#) which converts probability into scalar logit.

**Usage**

```
qmlogis_first(p)
```

```
qmlogis_last(p)
```

```
pmlogis_first(q)
```

```
pmlogis_last(q)
```

**Arguments**

`p` **numeric vector**, multinomial probabilities, adding up to 1

`q` **numeric vector**, multinomial logits

**Details**

Functions `pmlogis_first` and `pmlogis_last` take a length  $k - 1$  **numeric vector** of multinomial logits  $q$  and convert them into length  $k$  multinomial probabilities  $p$ , regarding the first or last category as reference, respectively.

Functions `qmlogis_first` and `qmlogis_last` take a length  $k$  **numeric vector** of multinomial probabilities  $p$  and convert them into length  $k - 1$  multinomial logits  $q$ , regarding the first or last category as reference, respectively.

**Value**

Functions `pmlogis_first` and `pmlogis_last` return a **vector** of multinomial probabilities  $p$ .

Functions `qmlogis_first` and `qmlogis_last` return a **vector** of multinomial logits  $q$ .

**See Also**

[plogis](#) [qlogis](#)

**Examples**

```
(a = qmlogis_last(c(2,5,3)))
(b = qmlogis_first(c(2,5,3)))
pmlogis_last(a)
pmlogis_first(b)

q0 = .8300964
(p1 = pmlogis_last(q0))
(q1 = qmlogis_last(p1))

# various exceptions
pmlogis_first(qmlogis_first(c(1, 0)))
pmlogis_first(qmlogis_first(c(0, 1)))
pmlogis_first(qmlogis_first(c(0, 0, 1)))
pmlogis_first(qmlogis_first(c(0, 1, 0, 0)))
pmlogis_first(qmlogis_first(c(1, 0, 0, 0)))
```

```

pmlogis_last(qmlogis_last(c(1, 0)))
pmlogis_last(qmlogis_last(c(0, 1)))
pmlogis_last(qmlogis_last(c(0, 0, 1)))
pmlogis_last(qmlogis_last(c(0, 1, 0, 0)))
pmlogis_last(qmlogis_last(c(1, 0, 0, 0)))

```

---

moment2fmx

*Creates [fmx](#) Object with Given Component-Wise Moments*


---

## Description

Creates [fmx](#) Object with Given Component-Wise Moments

## Usage

```
moment2fmx(distname, w, ...)
```

## Arguments

|          |   |
|----------|---|
| distname | <a href="#">character</a> scalar  |
| w        | <a href="#">numeric</a> vector  |
| ...      | <a href="#">numeric</a> scalars, some or all of mean, sd, skewness and kurtosis (length will be recycled), see <a href="#">moment2param</a> |

## Value

Function [moment2fmx](#) returns a [fmx](#) object.

## Examples

```

m = c(-1.5, 1.5)
s = c(.9, 1.1)
sk = c(.2, -.3)
kt = c(.5, .75)
w = c(2, 3)
(d1 = moment2fmx(distname='GH', w=w, mean=m, sd=s, skewness=sk, kurtosis=kt))
moment_fmx(d1)
(d2 = moment2fmx(distname='st', w=w, mean=m, sd=s, skewness=sk, kurtosis=kt))
moment_fmx(d2)
library(ggplot2)
ggplot() +
  geom_function(aes(color = 'GH'), fun = dfmx, args = list(dist=d1), n = 1001) +
  geom_function(aes(color = 'st'), fun = dfmx, args = list(dist=d1), n = 1001) +
  xlim(-5, 6)
# two curves looks really close, but actually not identical
x = rfmx(n = 1e3L, dist = d1)
range(dfmx(x, dist = d1) - dfmx(x, dist = d2))

```

---

|            |  |
|------------|--|
| moment_fmx | <i>Moment of Each Component in an <a href="#">fmx</a> Object</i> |
|------------|--|

---

**Description**

To find moments of each component in an [fmx](#) object.

**Usage**

```
moment_fmx(object)
```

**Arguments**

object            an [fmx](#) object

**Details**

Function [moment\\_fmx](#) calculates the [moments](#) and distribution characteristics of each mixture component of an S4 [fmx](#) object.

**Value**

Function [moment\\_fmx](#) returns a [moment](#) object.

**Examples**

```
(d2 = fmx('GH', A = c(1,6), B = 2, g = c(0,.3), h = c(.2,0), w = c(1,2)))
moment_fmx(d2)
```

---

|              |   |
|--------------|---|
| nobs.fitdist | <i>Number of Observations in <a href="#">fitdist</a> Object</i> |
|--------------|---|

---

**Description**

..

**Usage**

```
## S3 method for class 'fitdist'
nobs(object, ...)
```

**Arguments**

object            [fitdist](#) object  
 ...               additional parameters, currently not in use

**Value**

Function `nobs.fitdist` returns an `integer` scalar

---

|                       |  |
|-----------------------|--|
| <code>npar.fmx</code> | <i>Number of Parameters of <code>fmx</code> Object</i> |
|-----------------------|--|

---

**Description**

..

**Usage**

```
npar.fmx(dist)
```

**Arguments**

`dist`            `fmx` object

**Details**

Also the degree-of-freedom in `logLik`, as well as `stats::AIC.logLik` and `stats::BIC.logLik`

**Value**

Function `npar.fmx` returns an `integer` scalar.

---

|                        |   |
|------------------------|---|
| <code>print.fmx</code> | <i>S3 <code>print</code> of <code>fmx</code> Object</i> |
|------------------------|---|

---

**Description**

..

**Usage**

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

**Arguments**

`x`                    an `fmx` object  
`...`                additional parameters, not currently in use

**Value**

Function `print.fmx` returns the input `fmx` object invisibly.

---

|               |   |
|---------------|---|
| qfmx_interval | Obtain interval for <i>vuniroot2</i> for Function <i>qfmx</i> |
|---------------|---|

---

**Description**

Obtain interval for *vuniroot2* for Function *qfmx*

**Usage**

```
qfmx_interval(
  dist,
  p = c(1e-06, 1 - 1e-06),
  distname = dist@distname,
  K = dim(pars)[1L],
  pars = dist@pars,
  w = dist@w,
  ...
)
```

**Arguments**

|                      |   |
|----------------------|---|
| dist                 | <i>fmx</i> object                         |
| p                    | length-2 numeric vector                   |
| distname, K, pars, w | (optional) ignored if dist is provided    |
| ...                  | additional parameters, currently not used |

**Value**

Function *qfmx\_interval* returns a length-2 numeric vector.

---

|                  |                        |
|------------------|------------------------|
| show, fmx-method | Show <i>fmx</i> Object |
|------------------|------------------------|

---

**Description**

Print the parameters of an *fmx* object and plot its density curves.

**Usage**

```
## S4 method for signature 'fmx'
show(object)
```

**Arguments**

|        |                      |
|--------|----------------------|
| object | an <i>fmx</i> object |
|--------|----------------------|

**Value**

The [show](#) method for [fmx](#) object does not have a returned value.

---

 sort.mixEM

*Sort 'mixEM' Object by First Parameters*


---

**Description**

To sort a 'mixEM' object by its first parameters, i.e.,  $\mu$ 's for normal mixture,  $\alpha$ 's for  $\gamma$ -mixture, etc.

**Usage**

```
## S3 method for class 'mixEM'
sort(x, decreasing = FALSE, ...)
```

**Arguments**

|            |  |
|------------|--|
| x          | 'mixEM' object   |
| decreasing | <a href="#">logical</a> scalar. Should the sort by <i>mu</i> 's be increasing (FALSE, default) or decreasing (TRUE)? |
| ...        | additional parameters, currently not in use  |

**Details**

[normalmixEM](#) does *not* order the location parameter

**Value**

Function [sort.mixEM](#) returns a 'mixEM' object.

**See Also**

[sort](#)

---

|              |                                       |                |             |   |
|--------------|---------------------------------------|----------------|-------------|---|
| sort_mixsmsn | <i>Sort</i>                           | <i>Objects</i> | <i>from</i> | <i>Rhref<a href="https://CRAN.R-project.org/package=mixsmsn">https://CRAN.R-project.org/package=mixsmsn</a></i> |
|              | <i>mixsmsn by Location Parameters</i> |                |             |   |

---

## Description

To sort an object returned from package **mixsmsn** by its location parameters

## Usage

```
## S3 method for class 'Skew.normal'
sort(x, decreasing = FALSE, ...)

## S3 method for class 'Normal'
sort(x, decreasing = FALSE, ...)

## S3 method for class 'Skew.t'
sort(x, decreasing = FALSE, ...)

## S3 method for class 't'
sort(x, decreasing = FALSE, ...)
```

## Arguments

|            |  |
|------------|--|
| x          | 'Normal', 'Skew.normal', 'Skew.t' object   |
| decreasing | <b>logical</b> scalar. Should the sort the location parameter be increasing (FALSE, default) or decreasing (TRUE)? |
| ...        | additional parameters, currently not in use  |

## Details

[smsn.mix](#) does *not* order the location parameter

## Value

Function [sort.Normal](#) returns a 'Normal' object.

Function [sort.Skew.normal](#) returns a 'Skew.normal' object.

Function [sort.Skew.t](#) returns a 'Skew.t' object.

## See Also

[sort](#)



---

|                 |   |
|-----------------|---|
| user_constraint | <i>Formalize User-Specified Constraint of <code>fmx</code> Object</i> |
|-----------------|---|

---

## Description

Formalize user-specified constraint of `fmx` object

## Usage

```
user_constraint(x, distname, K)
```

## Arguments

|          |   |
|----------|---|
| x        | <b>character vector</b> , constraint(s) to be imposed. For example, for a two-component Tukey <i>g</i> -&- <i>h</i> mixture, <code>c('g1', 'h2')</code> indicates $g_1 = h_2 = 0$ given $A_1 < A_2$ , i.e., the <i>g</i> -parameter for the first component (with smaller location value) and the <i>h</i> -parameter for the second component (with larger mean value) are to be constrained as 0. |
| distname | <b>character</b> scalar, name of distribution   |
| K        | <b>integer</b> scalar, number of components   |

## Value

Function `user_constraint` returns the indices of internal parameters (only applicable to Tukey's *g*-&-*h* mixture distribution, yet) to be constrained, based on the type of distribution `distname`, number of components `K` and a user-specified string (e.g., `c('g2', 'h1')`).

## Examples

```
(d0 = fmx('GH', A = c(1,4), g = c(.2,.1), h = c(.05,.1), w = c(1,1)))
(c0 = fmx_constraint(d0))
user_constraint(distname = 'GH', K = 2L, x = character()) # equivalent

(d1 = fmx('GH', A = c(1,4), g = c(.2,0), h = c(0,.1), w = c(1,1)))
(c1 = fmx_constraint(d1))
user_constraint(distname = 'GH', K = 2L, x = c('g2', 'h1')) # equivalent

(d2 = fmx('GH', A = c(1,4), g = c(.2,0), h = c(.15,.1), w = c(1,1)))
(c2 = fmx_constraint(d2))
user_constraint(distname = 'GH', K = 2L, x = 'g2') # equivalent
```

vcov.fmx

*Variance-Covariance of fmx Object***Description**

..

**Usage**

```
## S3 method for class 'fmx'
vcov(object, internal = FALSE, ...)
```

**Arguments**

|          |  |
|----------|--|
| object   | fmx object   |
| internal | logical scalar, either for the user-friendly parameters (FALSE, default) (e.g., mean, sd for normal mixture, and A, B, g, h for Tukey <i>g</i> -and- <i>h</i> mixture), or for the internal/unconstrained parameters (TRUE). |
| ...      | place holder for S3 naming convention  |

**Details**

Function `vcov.fmx` returns the approximate asymptotic variance-covariance [matrix](#) of the user-friendly parameters via delta-method (`parm = 'user'`), or the asymptotic variance-covariance matrix of the internal/unconstrained parameters (`parm = 'internal'`). When the distribution has constraints on one or more parameters, function `vcov.fmx` does not return the variance/covariance involving the constrained parameters.

**Value**

Function `vcov.fmx` returns a [matrix](#).

[,fmx,ANY,ANY,ANY-method

*Subset of Components in fmx Object***Description**

Taking subset of components in [fmx](#) object

**Usage**

```
## S4 method for signature 'fmx,ANY,ANY,ANY'
x[i]
```

### Arguments

- x [fmx](#) object
- i [integer](#) or [logical vector](#), the row indices of *components* to be chosen, see [

### Details

Using definitions as S3 method dispatch ``[, fmx`` won't work for [fmx](#) objects.

### Value

An [fmx](#) object consisting of a subset of components. information about the observations (e.g. slots `@data` and `@data.name`), will be lost.

### Examples

```
(d = fmx('norm', mean = c(1, 4, 7), w = c(1, 1, 1)))  
d[1:2]
```

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