

# Package ‘RTMBdist’

May 7, 2026

**Type** Package

**Title** Distributions Compatible with Automatic Differentiation by 'RTMB'

**Version** 1.0.4

## Description

Extends the functionality of the 'RTMB' <<https://kaskr.r-universe.dev/RTMB>> package by providing a collection of non-standard probability distributions compatible with automatic differentiation (AD). While 'RTMB' enables flexible and efficient modelling, including random effects, its built-in support is limited to standard distributions. The package adds additional AD-compatible distributions, broadening the range of models that can be implemented and estimated using 'RTMB'. Automatic differentiation and Laplace approximation are described in Kristensen et al. (2016) <[doi:10.18637/jss.v070.i05](https://doi.org/10.18637/jss.v070.i05)>.

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**Encoding** UTF-8

**Imports** stats, Matrix, gamlss.dist, circular, sn, statmod, movMF

**Depends** R (>= 3.5.0), RTMB (>= 1.9.0)

**RoxygenNote** 7.3.3

**Suggests** knitr, rmarkdown, testthat (>= 3.0.0), TMB, gamlss.data, mvtnorm, LaMa (>= 2.1.0)

**Config/testthat/edition** 3

**URL** <https://janolefi.github.io/RTMBdist/>

**VignetteBuilder** knitr

**NeedsCompilation** no

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**Repository** CRAN

**Date/Publication** 2026-04-09 09:00:38 UTC

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abs_smooth	<i>Smooth approximation to the absolute value function</i>
------------	--

---

## Description

Smooth approximation to the absolute value function

## Usage

```
abs_smooth(x, epsilon = 1e-06)
```

## Arguments

x	vector of evaluation points
epsilon	smoothing constant

## Details

We approximate the absolute value here as

$$|x| \approx \sqrt{x^2 + \epsilon}$$

**Value**

Smooth absolute value of  $x$ .

**Examples**

```
abs(0)
abs_smooth(0, 1e-4)
```

---

 bccg

---

*Box–Cox Cole and Green distribution (BCCG)*


---

**Description**

Density, distribution function, quantile function, and random generation for the Box–Cox Cole and Green distribution.

**Usage**

```
dbccg(x, mu = 1, sigma = 0.1, nu = 1, log = FALSE)

pbccg(q, mu = 1, sigma = 0.1, nu = 1, lower.tail = TRUE, log.p = FALSE)

qbccg(p, mu = 1, sigma = 0.1, nu = 1, lower.tail = TRUE, log.p = FALSE)

rbccg(n, mu = 1, sigma = 0.1, nu = 1)
```

**Arguments**

<code>x, q</code>	vector of quantiles
<code>mu</code>	location parameter, must be positive.
<code>sigma</code>	scale parameter, must be positive.
<code>nu</code>	skewness parameter (real).
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
<code>p</code>	vector of probabilities
<code>n</code>	number of random values to return

**Details**

This implementation of `dbccg` and `pbccg` allows for automatic differentiation with RTMB while the other functions are imported from `gamlss.dist` package. See `gamlss.dist::BCCG` for more details.

**Value**

dbccg gives the density, pbccg gives the distribution function, qbccg gives the quantile function, and rbccg generates random deviates.

**References**

Rigby, R. A., Stasinopoulos, D. M., Heller, G. Z., and De Bastiani, F. (2019) Distributions for modeling location, scale, and shape: Using GAMLSS in R, Chapman and Hall/CRC, doi:10.1201/9780429298547. An older version can be found in <https://www.gamlss.com/>.

**Examples**

```
x <- rbccg(5, mu = 10, sigma = 0.2, nu = 0.5)
d <- dbccg(x, mu = 10, sigma = 0.2, nu = 0.5)
p <- pbccg(x, mu = 10, sigma = 0.2, nu = 0.5)
q <- qbccg(p, mu = 10, sigma = 0.2, nu = 0.5)
```

---

bcpe

*Box-Cox Power Exponential distribution (BCPE)*


---

**Description**

Density, distribution function, quantile function, and random generation for the Box-Cox Power Exponential distribution.

**Usage**

```
dbcpe(x, mu = 5, sigma = 0.1, nu = 1, tau = 2, log = FALSE)
pbcpe(q, mu = 5, sigma = 0.1, nu = 1, tau = 2, lower.tail = TRUE, log.p = FALSE)
qbcpe(p, mu = 5, sigma = 0.1, nu = 1, tau = 2, lower.tail = TRUE, log.p = FALSE)
rbcpe(n, mu = 5, sigma = 0.1, nu = 1, tau = 2)
```

**Arguments**

x, q	vector of quantiles
mu	location parameter, must be positive.
sigma	scale parameter, must be positive.
nu	vector of nu parameter values.
tau	vector of tau parameter values, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
p	vector of probabilities
n	number of random values to return

**Details**

This implementation of `dbcpe` and `pbcpce` allows for automatic differentiation with RTMB while the other functions are imported from `gamlss.dist` package. See `gamlss.dist:BCPE` for more details.

**Value**

`dbcpe` gives the density, `pbcpce` gives the distribution function, `qbcpe` gives the quantile function, and `rbcpe` generates random deviates.

**References**

Rigby, R. A., Stasinopoulos, D. M., Heller, G. Z., and De Bastiani, F. (2019) Distributions for modeling location, scale, and shape: Using GAMLSS in R, Chapman and Hall/CRC, doi:10.1201/9780429298547. An older version can be found in <https://www.gamlss.com/>.

**Examples**

```
x <- rbcpe(1, mu = 5, sigma = 0.1, nu = 1, tau = 1)
d <- dbcpe(x, mu = 5, sigma = 0.1, nu = 1, tau = 1)
p <- pbcpce(x, mu = 5, sigma = 0.1, nu = 1, tau = 1)
q <- qbcpe(p, mu = 5, sigma = 0.1, nu = 1, tau = 1)
```

---

bct

*Box–Cox t distribution (BCT)*


---

**Description**

Density, distribution function, quantile function, and random generation for the Box–Cox t distribution.

**Usage**

```
dbct(x, mu = 5, sigma = 0.1, nu = 1, tau = 2, log = FALSE)
pbct(q, mu = 5, sigma = 0.1, nu = 1, tau = 2, lower.tail = TRUE, log.p = FALSE)
qbct(p, mu = 5, sigma = 0.1, nu = 1, tau = 2, lower.tail = TRUE, log.p = FALSE)
rbct(n, mu = 5, sigma = 0.1, nu = 1, tau = 2)
```

**Arguments**

<code>x, q</code>	vector of quantiles
<code>mu</code>	location parameter, must be positive.
<code>sigma</code>	scale parameter, must be positive.
<code>nu</code>	skewness parameter (real).

tau	degrees of freedom, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
p	vector of probabilities
n	number of random values to return

### Details

This implementation of `dbct` and `pbct` allows for automatic differentiation with RTMB while the other functions are imported from `gamlss.dist` package. See `gamlss.dist`: [:BCT](#) for more details.

### Value

`dbct` gives the density, `pbct` gives the distribution function, `qbct` gives the quantile function, and `rbct` generates random deviates.

### References

Rigby, R. A., Stasinopoulos, D. M., Heller, G. Z., and De Bastiani, F. (2019) Distributions for modeling location, scale, and shape: Using GAMLSS in R, Chapman and Hall/CRC, doi:10.1201/9780429298547. An older version can be found in <https://www.gamlss.com/>.

### Examples

```
x <- rbct(1, mu = 10, sigma = 0.2, nu = 0.5, tau = 4)
d <- dbct(x, mu = 10, sigma = 0.2, nu = 0.5, tau = 4)
p <- pbct(x, mu = 10, sigma = 0.2, nu = 0.5, tau = 4)
q <- qbct(p, mu = 10, sigma = 0.2, nu = 0.5, tau = 4)
```

---

beta2	<i>Reparameterised beta distribution</i>
-------	--

---

### Description

Density, distribution function, quantile function, and random generation for the beta distribution reparameterised in terms of mean and concentration.

### Usage

```
dbeta(x, shape1, shape2, log = FALSE, eps = 0)

dbeta2(x, mu, phi, log = FALSE, eps = 0)

pbeta2(q, mu, phi, lower.tail = TRUE, log.p = FALSE)

qbeta2(p, mu, phi, lower.tail = TRUE, log.p = FALSE)

rbeta2(n, mu, phi)
```

**Arguments**

<code>x, q</code>	vector of quantiles
<code>shape1, shape2</code>	non-negative parameters
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>eps</code>	for internal use only, don't change.
<code>mu</code>	mean parameter, must be in the interval from 0 to 1.
<code>phi</code>	concentration parameter, must be positive.
<code>lower.tail</code>	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
<code>p</code>	vector of probabilities
<code>n</code>	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

Currently, `dbeta` masks `RTMB::dbeta` because the latter has a numerically unstable gradient.

**Value**

`dbeta2` gives the density, `pbeta2` gives the distribution function, `qbeta2` gives the quantile function, and `rbeta2` generates random deviates.

**Examples**

```
set.seed(123)
x <- rbeta2(1, 0.5, 1)
d <- dbeta2(x, 0.5, 1)
p <- pbeta2(x, 0.5, 1)
q <- qbeta2(p, 0.5, 1)
```

---

betabinom

*Beta-binomial distribution*


---

**Description**

Density and random generation for the beta-binomial distribution.

**Usage**

```
betabinom(x, size, shape1, shape2, log = FALSE)
```

```
rbetabinom(n, size, shape1, shape2)
```

**Arguments**

x	vector of non-negative counts.
size	vector of total counts (number of trials). Needs to be $\geq x$ .
shape1	positive shape parameter 1 of the Beta prior.
shape2	positive shape parameter 2 of the Beta prior.
log	logical; if TRUE, densities are returned on the log scale.
n	number of random values to return (for rbetabinom).

**Details**

This implementation of dbetabinom allows for automatic differentiation with RTMB.

**Value**

dbetabinom gives the density and rbetabinom generates random samples.

**Examples**

```
set.seed(123)
x <- rbetabinom(1, 10, 2, 5)
d <- dbetabinom(x, 10, 2, 5)
```

---

betaprime

*Beta prime distribution*

---

**Description**

Density, distribution function, quantile function, and random generation for the Beta prime distribution.

**Usage**

```
dbetaprime(x, shape1, shape2, log = FALSE)

pbetaprime(q, shape1, shape2, lower.tail = TRUE, log.p = FALSE)

qbetaprime(p, shape1, shape2, lower.tail = TRUE, log.p = FALSE)

rbetaprime(n, shape1, shape2)
```

**Arguments**

<code>x, q</code>	vector of quantiles
<code>shape1, shape2</code>	non-negative shape parameters of the corresponding Beta distribution
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
<code>p</code>	vector of probabilities
<code>n</code>	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

If  $X \sim \text{Beta}(\alpha, \beta)$ , then  $\frac{X}{1-X} \sim \text{Betaprime}(\alpha, \beta)$

**Value**

`dbetaprime` gives the density, `pbetaprime` gives the distribution function, `qbetaprime` gives the quantile function, and `rbetaprime` generates random deviates.

**Examples**

```
x <- rbetaprime(1, 2, 1)
d <- dbetaprime(x, 2, 1)
p <- pbetaprime(x, 2, 1)
q <- qbetaprime(p, 2, 1)
```

---

cclayton

*Clayton copula constructors*


---

**Description**

Construct a function that computes the log density or CDF of the bivariate Clayton copula, intended to be used with [dcopula](#).

**Usage**

```
cclayton(theta)
```

```
Cclayton(theta)
```

**Arguments**

<code>theta</code>	Positive dependence parameter ( $\theta > 0$ ).
--------------------	---

**Details**

The Clayton copula density is

$$c(u, v; \theta) = (1 + \theta)(uv)^{-(1+\theta)} (u^{-\theta} + v^{-\theta} - 1)^{-(2\theta+1)/\theta}, \quad \theta > 0.$$

**Value**

A function of two arguments (u,v) returning log copula **density** (cclayton) or copula **CDF** (Cclayton).

**See Also**

[cgaussian\(\)](#), [cgumbel\(\)](#), [cfrank\(\)](#)

**Examples**

```
x <- c(0.5, 1); y <- c(0.2, 0.8)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dbeta(y, 2, 1, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pbeta(y, 2, 1)
dcopula(d1, d2, p1, p2, copula = cclayton(2), log = TRUE)

# CDF version (for discrete copulas)
Cclayton(1.5)(0.5, 0.4)
```

---

cfrank

*Frank copula constructor*


---

**Description**

Returns a function computing the log density of the bivariate Frank copula, intended to be used with [dcopula](#).

**Usage**

```
cfrank(theta)
```

```
Cfrank(theta)
```

**Arguments**

theta            Dependence parameter ( $\theta \neq 0$ ).

**Details**

The Frank copula density is

$$c(u, v; \theta) = \frac{\theta(1 - e^{-\theta})e^{-\theta(u+v)}}{[(e^{-\theta u} - 1)(e^{-\theta v} - 1) + (1 - e^{-\theta})]^2}, \quad \theta \neq 0.$$

**Value**

A function of two arguments (u, v) returning either the log copula density (cfrank) or the copula CDF (Cfrank).

**See Also**

[cgaussian\(\)](#), [cclayton\(\)](#), [cgumbel\(\)](#)

**Examples**

```
x <- c(0.5, 1); y <- c(1, 2)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dexp(y, 2, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pexp(y, 2)
dcopula(d1, d2, p1, p2, copula = cfrank(2), log = TRUE)
```

---

cgaussian

*Gaussian copula constructor*

---

**Description**

Returns a function computing the log density of the bivariate Gaussian copula, intended to be used with [dcopula](#).

**Usage**

```
cgaussian(rho = 0)
```

**Arguments**

rho                      Correlation parameter ( $-1 < rho < 1$ ).

**Value**

Function of two arguments (u,v) returning log copula density.

The Gaussian copula density is

$$c(u, v; \rho) = \frac{1}{\sqrt{1 - \rho^2}} \exp \left\{ -\frac{1}{2(1 - \rho^2)} (z_1^2 - 2\rho z_1 z_2 + z_2^2) + \frac{1}{2} (z_1^2 + z_2^2) \right\},$$

where  $z_1 = \Phi^{-1}(u)$ ,  $z_2 = \Phi^{-1}(v)$ , and  $-1 < \rho < 1$ .

**See Also**

[cclayton\(\)](#), [cgumbel\(\)](#), [cfrank\(\)](#)

**Examples**

```
x <- c(0.5, 1); y <- c(1, 2)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dexp(y, 2, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pexp(y, 2)
dcopula(d1, d2, p1, p2, copula = cgaussian(0.5), log = TRUE)
```

---

cgmrf	<i>Multivariate Gaussian copula constructor parameterised by inverse correlation matrix</i>
-------	---

---

**Description**

Returns a function computing the log density of the multivariate Gaussian copula, parameterised by the inverse correlation matrix.

**Usage**

```
cgmrf(Q)
```

**Arguments**

Q                      Inverse of a positive definite correlation matrix with unit diagonal. Can either be sparse or dense matrix.

**Details**

**Caution:** Parameterising the inverse correlation directly is difficult, as inverting it needs to yield a positive definite matrix with **unit diagonal**. Hence we still advise parameterising the correlation matrix R and computing its inverse. This function is useful when you need access to the precision (i.e. inverse correlation) in your likelihood function.

**Value**

Function with matrix argument U returning log copula density.

**See Also**

[cmvgauss\(\)](#)

**Examples**

```
x <- c(0.5, 1); y <- c(1, 2); z <- c(0.2, 0.8)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dexp(y, 2, log = TRUE); d3 <- dbeta(z, 2, 1, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pexp(y, 2); p3 <- pbeta(z, 2, 1)
R <- matrix(c(1,0.5,0.3,0.5,1,0.4,0.3,0.4,1), nrow = 3)

## Based on correlation matrix
dmvcopula(cbind(d1, d2, d3), cbind(p1, p2, p3), copula = cmvgauss(R), log = TRUE)
```

```
## Based on precision matrix
Q <- solve(R)
dmvcopula(cbind(d1, d2, d3), cbind(p1, p2, p3), copula = cgmrf(Q), log = TRUE)

## Parameterisation inside a model
# using RTMB::unstructured to get a valid correlation matrix
library(RTMB)
d <- 5 # dimension
cor_func <- unstructured(d)
npar <- length(cor_func$params())
R <- cor_func$corr(rep(0.1, npar))
```

---

cgumbel

*Gumbel copula constructors*


---

### Description

Construct functions that compute either the log density or the CDF of the bivariate Gumbel copula, intended for use with [dcopula](#).

### Usage

```
cgumbel(theta)
```

```
Cgumbel(theta)
```

### Arguments

theta                    Dependence parameter ( $\theta \geq 1$ ).

### Details

The Gumbel copula density

$$c(u, v; \theta) = \exp \left[ - \left( (-\log u)^\theta + (-\log v)^\theta \right)^{1/\theta} \right] \cdot h(u, v; \theta),$$

where  $h(u, v; \theta)$  contains the derivative terms ensuring the function is a density.

### Value

A function of two arguments (u, v) returning either the log copula density (cgumbel) or the copula CDF (Cgumbel).

### See Also

[cgaussian\(\)](#), [cclayton\(\)](#), [cfrank\(\)](#)

**Examples**

```
x <- c(0.5, 1); y <- c(0.2, 0.4)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dbeta(y, 2, 1, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pbeta(y, 2, 1)
dcopula(d1, d2, p1, p2, copula = cgumbel(1.5), log = TRUE)

# CDF version (for discrete copulas)
Cgumbel(1.5)(0.5, 0.4)
```

---

cmvgauss

*Multivariate Gaussian copula constructor*


---

**Description**

Returns a function computing the log density of the multivariate Gaussian copula, intended to be used with [dmvcopula](#).

**Usage**

```
cmvgauss(R)
```

**Arguments**

R                    Positive definite correlation matrix (unit diagonal)

**Value**

Function with matrix argument U returning log copula density.

**See Also**

[cgmrf\(\)](#)

**Examples**

```
x <- c(0.5, 1); y <- c(1, 2); z <- c(0.2, 0.8)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dexp(y, 2, log = TRUE); d3 <- dbeta(z, 2, 1, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pexp(y, 2); p3 <- pbeta(z, 2, 1)
R <- matrix(c(1,0.5,0.3,0.5,1,0.4,0.3,0.4,1), nrow = 3)

## Based on correlation matrix
dmvcopula(cbind(d1, d2, d3), cbind(p1, p2, p3), copula = cmvgauss(R), log = TRUE)

## Based on precision matrix
Q <- solve(R)
dmvcopula(cbind(d1, d2, d3), cbind(p1, p2, p3), copula = cgmrf(Q), log = TRUE)

## Parameterisation inside a model
# using RTMB::unstructured to get a valid correlation matrix
```

```
library(RTMB)
d <- 5 # dimension
cor_func <- unstructured(d)
npar <- length(cor_func$params())
R <- cor_func$corr(rep(0.1, npar))
```

---

dcopula

*Joint density under a bivariate copula*


---

### Description

Computes the joint density (or log-density) of a bivariate distribution constructed from two arbitrary margins combined with a specified copula.

### Usage

```
dcopula(d1, d2, p1, p2, copula = cgaussian(0), log = FALSE)
```

### Arguments

d1, d2	Marginal density values. If log = TRUE, supply the log-density. If log = FALSE, supply the raw density.
p1, p2	Marginal CDF values. Need not be supplied on log scale.
copula	A function of two arguments (u, v) returning the log copula density $\log c(u, v)$ . You can either construct this yourself or use the copula constructors available (see details)
log	Logical; if TRUE, return the log joint density. In this case, d1 and d2 must be on the log scale.

### Details

The joint density is

$$f(x, y) = c(F_1(x), F_2(y)) f_1(x) f_2(y),$$

where  $F_i$  are the marginal CDFs,  $f_i$  are the marginal densities, and  $c$  is the copula density.

The marginal densities d1, d2 and CDFs p1, p2 must be differentiable for automatic differentiation (AD) to work.

Available copula constructors are:

- [cgaussian](#) (Gaussian copula)
- [cclayton](#) (Clayton copula)
- [cgumbel](#) (Gumbel copula)
- [cfrank](#) (Frank copula)

### Value

Joint density (or log-density) under the bivariate copula.

**See Also**

[ddcopula\(\)](#), [dmvcopula\(\)](#)

**Examples**

```
# Normal + Exponential margins with Gaussian copula
x <- c(0.5, 1); y <- c(1, 2)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dexp(y, 2, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pexp(y, 2)
ddcopula(d1, d2, p1, p2, copula = cgaussian(0.5), log = TRUE)

# Normal + Beta margins with Clayton copula
x <- c(0.5, 1); y <- c(0.2, 0.8)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dbeta(y, 2, 1, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pbeta(y, 2, 1)
ddcopula(d1, d2, p1, p2, copula = cclayton(2), log = TRUE)

# Normal + Beta margins with Gumbel copula
x <- c(0.5, 1); y <- c(0.2, 0.4)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dbeta(y, 2, 1, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pbeta(y, 2, 1)
ddcopula(d1, d2, p1, p2, copula = cgumbel(1.5), log = TRUE)

# Normal + Exponential margins with Frank copula
x <- c(0.5, 1); y <- c(1, 2)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dexp(y, 2, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pexp(y, 2)
ddcopula(d1, d2, p1, p2, copula = cfrank(2), log = TRUE)
```

---

ddcopula

*Joint probability under a discrete bivariate copula*


---

**Description**

Computes the joint probability mass function of two **discrete** margins combined with a copula CDF.

**Usage**

```
ddcopula(d1, d2, p1, p2, Copula, log = FALSE)
```

**Arguments**

d1, d2	Marginal p.m.f. values at the observed points, <b>not</b> on log-scale.
p1, p2	Marginal CDF values at the observed points.
Copula	A function of two arguments returning the copula CDF.
log	Logical; if TRUE, return the log joint density. In this case, d1 and d2 must be on the log scale.

**Details**

The joint probability mass function for two discrete margins is

$$\Pr(Y_1 = y_1, Y_2 = y_2) = C(F_1(y_1), F_2(y_2)) - C(F_1(y_1-1), F_2(y_2)) - C(F_1(y_1), F_2(y_2-1)) + C(F_1(y_1-1), F_2(y_2-1)),$$

where  $F_i$  are the marginal CDFs, and  $C$  is the copula CDF.

Available copula CDF constructors are:

- [Cclayton](#) (Clayton copula)
- [Cgumbel](#) (Gumbel copula)
- [Cfrank](#) (Frank copula)

**Value**

Joint probability (or log-probability) under chosen copula

**See Also**

[dcopula\(\)](#), [dmvcopula\(\)](#)

**Examples**

```
x <- c(3,5); y <- c(2,4)
d1 <- dpois(x, 4); d2 <- dpois(y, 3)
p1 <- ppois(x, 4); p2 <- ppois(y, 3)
ddcopula(d1, d2, p1, p2, Copula = Cclayton(2), log = FALSE)
```

---

dirichlet

*Dirichlet distribution*

---

**Description**

Density and random generation for the Dirichlet distribution.

**Usage**

```
ddirichlet(x, alpha, log = FALSE)
```

```
rdirichlet(n, alpha)
```

**Arguments**

x	vector or matrix of quantiles. If x is a vector, it needs to sum to one. If x is a matrix, each row should sum to one.
alpha	vector or matrix of positive shape parameters
log	logical; if TRUE, densities $p$ are returned as $\log(p)$ .
n	number of random values to return.

**Details**

This implementation of `ddirichlet` allows for automatic differentiation with RTMB.

**Value**

`ddirichlet` gives the density, `rdirichlet` generates random deviates.

**Examples**

```
# single alpha
alpha <- c(1,2,3)
x <- rdirichlet(1, alpha)
d <- ddirichlet(x, alpha)
# vectorised over alpha
alpha <- rbind(alpha, 2*alpha)
x <- rdirichlet(2, alpha)
```

---

dirmult

*Dirichlet-multinomial distribution*


---

**Description**

Density and random generation for the Dirichlet-multinomial distribution.

**Usage**

```
ddirmult(x, size, alpha, log = FALSE)
```

```
rdirmult(n, size, alpha)
```

**Arguments**

<code>x</code>	vector or matrix of non-negative counts, where rows are observations and columns are categories.
<code>size</code>	vector of total counts for each observation. Needs to match the row sums of <code>x</code> .
<code>alpha</code>	vector or matrix of positive shape parameters
<code>log</code>	logical; if TRUE, densities $p$ are returned as $\log(p)$ .
<code>n</code>	number of random values to return.

**Details**

This implementation of `ddirmult` allows for automatic differentiation with RTMB.

**Value**

`ddirmult` gives the density and `rdirmult` generates random samples.

**Examples**

```
# single alpha
alpha <- c(1,2,3)
size <- 10
x <- rdirmult(1, size, alpha)
d <- ddirmult(x, size, alpha)
# vectorised over alpha and size
alpha <- rbind(alpha, 2*alpha)
size <- c(size, 3*size)
x <- rdirmult(2, size, alpha)
```

dmvcopula

*Joint density under a multivariate copula***Description**

Computes the joint density (or log-density) of a distribution constructed from any number of arbitrary margins combined with a specified copula.

**Usage**

```
dmvcopula(D, P, copula = cmvgauss(diag(ncol(D))), log = FALSE)
```

**Arguments**

D	Matrix of marginal density values of with rows corresponding to observations and columns corresponding to dimensions. If log = TRUE, supply the log-densities. If log = FALSE, supply the raw densities
P	Matrix of marginal CDF values of the same dimension as D. Need not be supplied on log scale.
copula	A function of a matrix argument U returning the log copula density $\log c(u_1, \dots, u_d)$ . The columns of U correspond to dimensions. You can either construct this yourself or use the copula constructors available (see details)
log	Logical; if TRUE, return the log joint density. In this case, D must be on the log scale.

**Details**

The joint density is

$$f(x_1, \dots, x_d) = c(F_1(x_1), \dots, F_d(x_d)) f_1(x_1) \dots f_d(x_d),$$

where  $F_i$  are the marginal CDFs,  $f_i$  are the marginal densities, and  $c$  is the copula density.

The marginal densities  $d_1, \dots, d_d$  and CDFs  $p_1, \dots, p_d$  must be differentiable for automatic differentiation (AD) to work.

Available multivariate copula constructors are:

- [cmvgauss](#) (Multivariate Gaussian copula)
- [cgmrf](#) (Multivariate Gaussian copula parameterised by precision (inverse correlation) matrix)

**Value**

Joint density (or log-density) under the chosen copula.

**See Also**

[dcopula\(\)](#), [ddcopula\(\)](#)

**Examples**

```
x <- c(0.5, 1); y <- c(1, 2); z <- c(0.2, 0.8)
d1 <- dnorm(x, 1, log = TRUE); d2 <- dexp(y, 2, log = TRUE); d3 <- dbeta(z, 2, 1, log = TRUE)
p1 <- pnorm(x, 1); p2 <- pexp(y, 2); p3 <- pbeta(z, 2, 1)
R <- matrix(c(1,0.5,0.3,0.5,1,0.4,0.3,0.4,1), nrow = 3)

### Multivariate Gaussian copula
## Based on correlation matrix
dmvcopula(cbind(d1, d2, d3), cbind(p1, p2, p3), copula = cmvgauss(R), log = TRUE)

## Based on precision matrix
Q <- solve(R)
dmvcopula(cbind(d1, d2, d3), cbind(p1, p2, p3), copula = cgmrf(Q), log = TRUE)

## Parameterisation inside a model
# using RTMB::unstructured to get a valid correlation matrix
library(RTMB)
d <- 5 # dimension
cor_func <- unstructured(d)
npar <- length(cor_func$params())
R <- cor_func$corr(rep(0.1, npar))
```

---

erf

*AD-compatible error function and complementary error function*


---

**Description**

AD-compatible error function and complementary error function

**Usage**

```
erf(x)
```

```
erfc(x)
```

**Arguments**

x                      vector of evaluation points

**Value**

erf(x) returns the error function and erfc(x) returns the complementary error function.

**Examples**

```
erf(1)
erfc(1)
```

---

exgauss

*Exponentially modified Gaussian distribution*

---

**Description**

Density, distribution function, quantile function, and random generation for the exponentially modified Gaussian distribution.

**Usage**

```
dexgauss(x, mu = 0, sigma = 1, lambda = 1, log = FALSE)
pexgauss(q, mu = 0, sigma = 1, lambda = 1, lower.tail = TRUE, log.p = FALSE)
qexgauss(p, mu = 0, sigma = 1, lambda = 1, lower.tail = TRUE, log.p = FALSE)
rexgauss(n, mu = 0, sigma = 1, lambda = 1)
```

**Arguments**

x, q	vector of quantiles
mu	mean parameter of the Gaussian part
sigma	standard deviation parameter of the Gaussian part, must be positive.
lambda	rate parameter of the exponential part, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
n	number of random values to return

**Details**

This implementation of dexgauss and pexgauss allows for automatic differentiation with RTMB. qexgauss and rexgauss are reparameterised imports from gamlss.dist::exGAUS.

If  $X \sim N(\mu, \sigma^2)$  and  $Y \sim \text{Exp}(\lambda)$ , then  $Z = X + Y$  follows the exponentially modified Gaussian distribution with parameters  $\mu$ ,  $\sigma$ , and  $\lambda$ .

**Value**

dexgauss gives the density, pexgauss gives the distribution function, qexgauss gives the quantile function, and rexgauss generates random deviates.

## References

Rigby, R. A., Stasinopoulos, D. M., Heller, G. Z., and De Bastiani, F. (2019) Distributions for modeling location, scale, and shape: Using GAMLSS in R, Chapman and Hall/CRC, doi:10.1201/9780429298547. An older version can be found in <https://www.gamlss.com/>.

## Examples

```
x <- rexgauss(1, 1, 2, 2)
d <- dexgauss(x, 1, 2, 2)
p <- pexgauss(x, 1, 2, 2)
q <- qexgauss(p, 1, 2, 2)
```

---

foldnorm

*Folded normal distribution*

---

## Description

Density, distribution function, and random generation for the folded normal distribution.

## Usage

```
dfoldnorm(x, mu = 0, sigma = 1, log = FALSE)

pfoldnorm(q, mu = 0, sigma = 1, lower.tail = TRUE, log.p = FALSE)

rfoldnorm(n, mu = 0, sigma = 1)
```

## Arguments

x, q	vector of quantiles
mu	location parameter
sigma	scale parameter, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return
p	vector of probabilities

## Details

This implementation of `dfoldnorm` allows for automatic differentiation with RTMB.

## Value

`dfoldnorm` gives the density, `pfoldnorm` gives the distribution function, and `rfoldnorm` generates random deviates.

**Examples**

```
x <- rfoldnorm(1, 1, 2)
d <- dfoldnorm(x, 1, 2)
p <- pfoldnorm(x, 1, 2)
```

---

gamma2

*Reparameterised gamma distribution*


---

**Description**

Density, distribution function, quantile function, and random generation for the gamma distribution reparameterised in terms of mean and standard deviation.

**Usage**

```
dgamma2(x, mean = 1, sd = 1, log = FALSE)

pgamma2(q, mean = 1, sd = 1, lower.tail = TRUE, log.p = FALSE)

qgamma2(p, mean = 1, sd = 1, lower.tail = TRUE, log.p = FALSE)

rgamma2(n, mean = 1, sd = 1)
```

**Arguments**

x, q	vector of quantiles
mean	mean parameter, must be positive.
sd	standard deviation parameter, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
n	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

dgamma2 gives the density, pgamma2 gives the distribution function, qgamma2 gives the quantile function, and rgamma2 generates random deviates.

**Examples**

```
x <- rgamma2(1)
d <- dgamma2(x)
p <- pgamma2(x)
q <- qgamma2(p)
```

---

gengamma	<i>Generalised Gamma distribution (GG)</i>
----------	--

---

### Description

Density, distribution function, quantile function, and random generation for the generalised Gamma distribution.

### Usage

```
dgengamma(x, mu = 1, sigma = 0.5, nu = 1, log = FALSE)
```

```
pgengamma(q, mu = 1, sigma = 0.5, nu = 1, lower.tail = TRUE, log.p = FALSE)
```

```
qgengamma(p, mu = 1, sigma = 0.5, nu = 1, lower.tail = TRUE, log.p = FALSE)
```

```
rgengamma(n, mu = 1, sigma = 0.5, nu = 1)
```

### Arguments

x, q	vector of quantiles
mu	location parameter, must be positive.
sigma	scale parameter, must be positive.
nu	skewness parameter (real).
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
p	vector of probabilities
n	number of random values to return

### Details

This implementation of `dgengamma`, `pgengamma`, and `qgengamma` allows for automatic differentiation with RTMB.

### Value

`dgengamma` gives the density, `pgengamma` gives the distribution function, `qgengamma` gives the quantile function, and `rgengamma` generates random deviates.

### References

Rigby, R. A., Stasinopoulos, D. M., Heller, G. Z., and De Bastiani, F. (2019) Distributions for modeling location, scale, and shape: Using GAMLSS in R, Chapman and Hall/CRC, doi:10.1201/9780429298547. An older version can be found in <https://www.gamlss.com/>.

**Examples**

```
x <- rgengamma(5, mu = 4, sigma = 0.5, nu = 0.5)
d <- dgengamma(x, mu = 4, sigma = 0.5, nu = 0.5)
p <- pgengamma(x, mu = 4, sigma = 0.5, nu = 0.5)
q <- qgengamma(p, mu = 4, sigma = 0.5, nu = 0.5)
```

genpois

*Generalised Poisson distribution***Description**

Probability mass function, distribution function, and random generation for the generalised Poisson distribution.

**Usage**

```
dgenpois(x, lambda = 1, phi = 1, log = FALSE)

pgenpois(q, lambda = 1, phi = 1, lower.tail = TRUE, log.p = FALSE)

qgenpois(p, lambda = 1, phi = 1,
         lower.tail = TRUE, log.p = FALSE, max.value = 10000)

rgenpois(n, lambda = 1, phi = 1, max.value = 10000)
```

**Arguments**

x, q	integer vector of counts
lambda	vector of positive means
phi	vector of non-negative dispersion parameters
log, log.p	logical; return log-density if TRUE
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
max.value	a constant, set to the default value of 10000 for how far the algorithm should look for q.
n	number of random values to return.

**Details**

This implementation of `dgenpois` allows for automatic differentiation with RTMB. The other functions are imported from `gamlss.dist::GPO`.

The distribution has mean  $\lambda$  and variance  $\lambda(1 + \phi\lambda)^2$ . For  $\phi = 0$  it reduces to the Poisson distribution, however  $\phi$  must be strictly positive here.

**Value**

dgenpois gives the probability mass function, pgenpois gives the distribution function, qgenpois gives the quantile function, and rgenpois generates random deviates.

**Examples**

```
set.seed(123)
x <- rgenpois(1, 2, 3)
d <- dgenpois(x, 2, 3)
p <- pgenpois(x, 2, 3)
q <- qgenpois(p, 2, 3)
```

gumbel

*Gumbel distribution***Description**

Density, distribution function, quantile function, and random generation for the Gumbel distribution.

**Usage**

```
dgumbel(x, location = 0, scale = 1, log = FALSE)
pgumbel(q, location = 0, scale = 1, lower.tail = TRUE, log.p = FALSE)
qgumbel(p, location = 0, scale = 1, lower.tail = TRUE, log.p = FALSE)
rgumbel(n, location = 0, scale = 1)
```

**Arguments**

x, q	vector of quantiles
location	location parameter
scale	scale parameter, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
n	number of random values to return

**Details**

This implementation of dgumbel allows for automatic differentiation with RTMB.

**Value**

dgumbel gives the density, pgumbel gives the distribution function, qgumbel gives the quantile function, and rgumbel generates random deviates.

**Examples**

```
x <- rgumbel(1, 0.5, 2)
d <- dgumbel(x, 0.5, 2)
p <- pgumbel(x, 0.5, 2)
q <- qgumbel(p, 0.5, 2)
```

---

 invchisq

*Inverse Chi-squared distribution*


---

**Description**

Density, distribution function, quantile function, and random generation for the inverse Chi-squared distribution.

**Usage**

```
dinvchisq(x, df, scale = 1/df, log = FALSE)
pinvchisq(q, df, scale = 1/df, lower.tail = TRUE, log.p = FALSE)
qinvchisq(p, df, scale = 1/df, lower.tail = TRUE, log.p = FALSE)
rinvchisq(n, df, scale = 1/df)
```

**Arguments**

x, q	vector of quantiles, must be positive.
df	degrees of freedom ( $\nu > 0$ )
scale	optional positive scale parameter. Default value of 1/df corresponds to standard inverse gamma
log, log.p	logical; if TRUE, probabilities/densities are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
n	number of random values to return

**Details**

If  $X \sim \text{Chisq}(\nu)$ , then  $1/X \sim \text{invChisq}(\nu)$ .

The inverse Chi-squared distribution with  $\nu$  degrees of freedom has density

$$f(x) = \frac{(\nu/2)^{\nu/2}}{\Gamma(\nu/2)} x^{-(\nu/2+1)} \exp(-\nu/(2x)), \quad x > 0.$$

This implementation of `dinvchisq`, `pinvchisq`, and `qinvchisq` allows for automatic differentiation with RTMB.

**Value**

`dinvchisq` gives the density, `pinvchisq` gives the distribution function, `qinvchisq` gives the quantile function, and `rinvchisq` generates random deviates.

**Examples**

```
x <- rinvchisq(1, df = 5)
d <- dinvchisq(x, df = 5)
p <- pinvchisq(x, df = 5)
q <- qinvchisq(p, df = 5)
```

---

 invgamma

*Inverse Gamma distribution*


---

**Description**

Density, distribution function, and random generation for the inverse Gamma distribution.

**Usage**

```
dinvgamma(x, shape, rate, scale = 1/rate, log = FALSE)
pinvgamma(q, shape, rate, scale = 1/rate, lower.tail = TRUE, log.p = FALSE)
qinvgamma(p, shape, rate, scale = 1/rate, lower.tail = TRUE, log.p = FALSE)
rinvgamma(n, shape, rate, scale = 1/rate)
```

**Arguments**

<code>x, q</code>	vector of quantiles, must be positive.
<code>shape, rate, scale</code>	positive parameters of corresponding gamma distribution
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
<code>p</code>	vector of probabilities
<code>n</code>	number of random values to return

**Details**

This implementation of `dinvgamma`, `pinvgamma`, and `qinvgamma` allows for automatic differentiation with RTMB.

If  $X \sim \Gamma(\alpha, \beta)$ , then  $1/X \sim \text{InvGamma}(\alpha, \beta)$ .

**Value**

`dinvgamma` gives the density, `pinvgamma` gives the distribution function, `qinvgamma` gives the quantile function, and `rinvgamma` generates random deviates.

**Examples**

```
x <- rinvgamma(1, 1, 0.5)
d <- dinvgamma(x, 1, 0.5)
p <- pinvgamma(x, 1, 0.5)
q <- qinvgamma(p, 1, 0.5)
```

---

 invgauss

*Inverse Gaussian distribution*


---

**Description**

Density, distribution function, and random generation for the inverse Gaussian distribution.

**Usage**

```
dinvgauss(x, mean = 1, shape = 1, log = FALSE)

pinvgauss(q, mean = 1, shape = 1, lower.tail = TRUE, log.p = FALSE)

qinvgauss(p, mean = 1, shape = 1, lower.tail = TRUE, log.p = FALSE, ...)

rinvgauss(n, mean = 1, shape = 1)
```

**Arguments**

<code>x, q</code>	vector of quantiles, must be positive.
<code>mean</code>	location parameter
<code>shape</code>	shape parameter, must be positive.
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
<code>p</code>	vector of probabilities
<code>...</code>	additional parameter passed to <code>statmod::qinvgauss</code> for numerical evaluation of the quantile function.
<code>n</code>	number of random values to return

**Details**

This implementation of `dinvgauss` allows for automatic differentiation with RTMB. `qinvgauss` and `rinvgauss` are imported from the `statmod` package.

**Value**

`dinvgauss` gives the density, `pinvgauss` gives the distribution function, `qinvgauss` gives the quantile function, and `rinvgauss` generates random deviates.

**Examples**

```
x <- rinvgauss(1, 1, 0.5)
d <- dinvgauss(x, 1, 0.5)
p <- pinvgauss(x, 1, 0.5)
q <- qinvgauss(p, 1, 0.5)
```

---

kumar

*Kumaraswamy distribution*


---

**Description**

Density, distribution function, quantile function, and random generation for the Kumaraswamy distribution.

**Usage**

```
dkumar(x, a, b, log = FALSE)

pkumar(q, a, b, lower.tail = TRUE, log.p = FALSE)

qkumar(p, a, b, lower.tail = TRUE, log.p = FALSE)

rkumar(n, a, b)
```

**Arguments**

<code>x, q</code>	vector of quantiles in $(0, 1)$
<code>a, b</code>	positive shape parameters
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
<code>p</code>	vector of probabilities
<code>n</code>	number of random values to return.

**Value**

`dkumar` gives the density, `pkumar` gives the distribution function, `qkumar` gives the quantile function, and `rkumar` generates random deviates.

**Examples**

```
x <- rkumar(1, a = 1, b = 2)
d <- dkumar(x, a = 1, b = 2)
p <- pkumar(x, a = 1, b = 2)
q <- qkumar(p, a = 1, b = 2)
```

---

laplace

*Laplace distribution*


---

**Description**

Density, distribution function, quantile function, and random generation for the Laplace distribution.

**Usage**

```
dlaplace(x, mu = 0, b = 1, eps = NULL, log = FALSE)

plaplace(q, mu = 0, b = 1, lower.tail = TRUE, log.p = FALSE)

qlaplace(p, mu = 0, b = 1, lower.tail = TRUE, log.p = FALSE)

rlaplace(n, mu = 0, b = 1)
```

**Arguments**

x, q	vector of quantiles
mu	location parameter
b	scale parameter, must be positive.
eps	optional smoothing parameter for <code>dlaplace</code> to smooth the absolute value function. See <a href="#">abs_smooth</a> for details. It is recommended to set this to a small constant like 1e-6 for numerical optimisation.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
n	number of random values to return

**Details**

This implementation of `dlaplace` allows for automatic differentiation with RTMB.

**Value**

`dlaplace` gives the density, `plaplace` gives the distribution function, `qlaplace` gives the quantile function, and `rlaplace` generates random deviates.

**Examples**

```
x <- rlaplace(1, 1, 1)
d <- dlaplace(x, 1, 1)
p <- plaplace(x, 1, 1)
q <- qlaplace(p, 1, 1)
```

---

mcreport

*Sample parameters from approximate Gaussian posterior distribution*


---

**Description**

Efficient Monte Carlo sampling of parameters (and REPORTed quantities) from the approximate posterior of an RTMB model. See [sdreport](#) for details on posterior variance-covariance in random effects models.

**Usage**

```
mcreport(
  obj,
  nSamples = 1000,
  include_random_pars = TRUE,
  report = FALSE,
  Q = NULL,
  ...
)
```

**Arguments**

obj	Optimised RTMB object generated by <a href="#">MakeADFun</a>
nSamples	Number of samples to draw
include_random_pars	Logical; Should random parameters be included in the output?
report	Logical; Should REPORTed quantities be sampled as well? Defaults to FALSE because this may be slow depending on your model.
Q	Optional precalculated sparse precision matrix returned by <code>sdreport(..., getJointPrecision = TRUE)\$jointPrecision</code> . If not provided, computed internally using <code>sdreport</code> . Only used for models with random effects.
...	For internal use only

**Details**

**Caution:** This has nothing to do with Bayesian posterior sampling. It is simple Monte Carlo sampling from a Gaussian distribution with mean at the MLE and covariance given by the inverse of the Hessian (for fixed effects models) or the joint precision matrix (for random effects models). This is a common approach to get an approximate idea of parameter uncertainty around the MLE, but it relies on large-sample asymptotics.

Sampling random effects from their posterior as compared to calculating marginal standard deviations like `sdreport` is particularly useful for multidimensional random effects (e.g. for locations  $x$  and  $y$ ) where pointwise confidence intervals (e.g. along a path) based on standard deviations are not possible.

## Value

A list structured like the original parameter list used in the `MakeADFun` call (potentially including additional REPORTed quantities). Each entry is a list with `nSamples` entries.

## Examples

```
set.seed(123)

### Example 1: Without random effects ###

# simulate data
x <- rgumbel(100, location = 5, scale = 2)

# negative log-likelihood function
nll <- function(par) {
  getAll(par)
  x <- OBS(x) # mark x as the response
  scale <- exp(log_scale)
  ADREPORT(scale); REPORT(scale)
  -sum(dgumbel(x, loc, scale, log = TRUE))
}

# RTMB AD object
par <- list(loc = 5, log_scale = log(2))
obj <- MakeADFun(nll, par, silent = TRUE)

# model fitting using AD gradient
opt <- nlminb(obj$par, obj$fn, obj$gr)

# sample from asymptotic distribution of the MLE
samples <- mcreport(obj, report = TRUE)
# parameters
mean(unlist(samples$loc)); sd(unlist(samples$loc))
# reported transformed parameters
mean(unlist(samples$scale)); sd(unlist(samples$scale))

# compare to delta method sd from sdreport
sdr <- sdreport(obj)
summary(sdr)

### Example 2: With random effects ###

# simulate random effects
id <- rep(1:10, each = 100) # 10 groups with 100 observations each
```

```

true_gamma <- rnorm(10, 0, 2) # random coefficients
true_probs <- plogis(1 + true_gamma) # linear predictor

# simulate Bernoulli data conditional on random coefficients
x <- rbinom(1000, 1, true_probs[id])

# negative log-likelihood function
nll <- function(par) {
  getAll(par)
  x <- OBS(x) # mark x as the response

  # random effect likelihood
  sd_gamma <- exp(log_sd_gamma)
  ADREPORT(sd_gamma); REPORT(sd_gamma)
  nll <- -sum(dnorm(gamma, 0, sd_gamma, log = TRUE))

  # data likelihood
  probs <- plogis(beta0 + gamma) # linear predictor
  ADREPORT(probs); REPORT(probs)
  nll <- nll - sum(dbinom(x, 1, probs[id], log = TRUE))
  return(nll)
}

# RTMB Laplace approximation
par <- list(beta0 = 1, log_sd_gamma = log(2), gamma = rep(0,10))
obj <- MakeADFun(nll, par, random = "gamma", silent = TRUE)

# model fitting using AD gradient
opt <- nlmminb(obj$par, obj$fn, obj$gr)

# draw samples
samples <- mcreport(obj, report = TRUE)

# run sdreport
sdr <- sdreport(obj)

# standard deviation using delta method
as.list(sdr, "Std", report = TRUE)$probs

# standard deviation from samples
apply(simplify2array(samples$probs), 1, sd)

```

## Description

Density and random generation for the multivariate t distribution

**Usage**

```
dmvt(x, mu, Sigma, df, log = FALSE)
```

```
rmvt(n, mu, Sigma, df)
```

**Arguments**

x	vector or matrix of quantiles
mu	vector or matrix of location parameters (mean if $df > 1$ )
Sigma	positive definite scale matrix (proportional to the covariance matrix if $df > 2$ )
df	degrees of freedom; must be positive
log	logical; if TRUE, densities $p$ are returned as $\log(p)$ .
n	number of random values to return.

**Details**

This implementation of `dmvt` allows for automatic differentiation with RTMB.

Note: for  $df \leq 1$  the mean is undefined, and for  $df \leq 2$  the covariance is infinite. For  $df > 2$ , the covariance is  $df/(df-2) * \text{Sigma}$ .

**Value**

`dmvt` gives the density, `rmvt` generates random deviates.

**Examples**

```
# single mu
mu <- c(1,2,3)
Sigma <- diag(c(1,1,1))
df <- 5
x <- rmvt(2, mu, Sigma, df)
d <- dmvt(x, mu, Sigma, df)
# vectorised over mu
mu <- rbind(c(1,2,3), c(0, 0.5, 1))
x <- rmvt(2, mu, Sigma, df)
d <- dmvt(x, mu, Sigma, df)
```

---

nbinom2

*Reparameterised negative binomial distribution*


---

**Description**

Probability mass function, distribution function, quantile function, and random generation for the negative binomial distribution reparameterised in terms of mean and size.

**Usage**

```
dnbinom2(x, mu, size, log = FALSE)
pnbinom2(q, mu, size, lower.tail = TRUE, log.p = FALSE)
qnbinom2(p, mu, size, lower.tail = TRUE, log.p = FALSE)
rnbinom2(n, mu, size)
pnbinom(q, size, prob, lower.tail = TRUE, log.p = FALSE)
```

**Arguments**

x, q	vector of quantiles
mu	mean parameter, must be positive.
size	size parameter, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
n	number of random values to return.
prob	probability of success in each trial. $0 < \text{prob} \leq 1$ .

**Details**

This implementation allows for automatic differentiation with RTMB.

pnbinom is an AD-compatible implementation of the standard parameterisation of the CDF, missing from RTMB.

**Value**

dnbinom2 gives the density, pnbinom2 gives the distribution function, qnbinom2 gives the quantile function, and rnbinom2 generates random deviates.

**Examples**

```
set.seed(123)
x <- rnbinom2(1, 1, 2)
d <- dnbinom2(x, 1, 2)
p <- pnbinom2(x, 1, 2)
q <- qnbinom2(p, 1, 2)
```

---

`oibeta`*One-inflated beta distribution*

---

**Description**

Density, distribution function, and random generation for the one-inflated beta distribution.

**Usage**

```
doibeta(x, shape1, shape2, oneprob = 0, log = FALSE)
```

```
poibeta(q, shape1, shape2, oneprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
roibeta(n, shape1, shape2, oneprob = 0)
```

**Arguments**

<code>x, q</code>	vector of quantiles
<code>shape1, shape2</code>	non-negative shape parameters of the beta distribution
<code>oneprob</code>	zero-inflation probability between 0 and 1.
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
<code>n</code>	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

`doibeta` gives the density, `poibeta` gives the distribution function, and `roibeta` generates random deviates.

**Examples**

```
set.seed(123)
x <- roibeta(1, 2, 2, 0.5)
d <- doibeta(x, 2, 2, 0.5)
p <- poibeta(x, 2, 2, 0.5)
```

---

`oibeta2`*Reparameterised one-inflated beta distribution*

---

### Description

Density, distribution function, and random generation for the one-inflated beta distribution reparameterised in terms of mean and concentration.

### Usage

```
doibeta2(x, mu, phi, oneprob = 0, log = FALSE)
```

```
poibeta2(q, mu, phi, oneprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
roibeta2(n, mu, phi, oneprob = 0)
```

### Arguments

<code>x, q</code>	vector of quantiles
<code>mu</code>	mean parameter, must be in the interval from 0 to 1.
<code>phi</code>	concentration parameter, must be positive.
<code>oneprob</code>	zero-inflation probability between 0 and 1.
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
<code>n</code>	number of random values to return.

### Details

This implementation allows for automatic differentiation with RTMB.

### Value

`doibeta2` gives the density, `poibeta2` gives the distribution function, and `roibeta2` generates random deviates.

### Examples

```
set.seed(123)
x <- roibeta2(1, 0.6, 2, 0.5)
d <- doibeta2(x, 0.6, 2, 0.5)
p <- poibeta2(x, 0.6, 2, 0.5)
```

---

pareto *Pareto distribution*

---

### Description

Density, distribution function, quantile function, and random generation for the pareto distribution.

### Usage

```
dpareto(x, mu = 1, log = FALSE)
```

```
ppareto(q, mu = 1, lower.tail = TRUE, log.p = FALSE)
```

```
qpareto(p, mu = 1, lower.tail = TRUE, log.p = FALSE)
```

```
rpareto(n, mu = 1)
```

### Arguments

<code>x, q</code>	vector of quantiles
<code>mu</code>	location parameter, must be positive.
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
<code>p</code>	vector of probabilities
<code>n</code>	number of random values to return

### Details

This implementation of `dpareto` and `ppareto` allows for automatic differentiation with RTMB while the other functions are imported from `gamlss.dist` package. See `gamlss.dist::PARETO` for more details.

### Value

`dpareto` gives the density, `ppareto` gives the distribution function, `qpareto` gives the quantile function, and `rpareto` generates random deviates.

### References

Rigby, R. A., Stasinopoulos, D. M., Heller, G. Z., and De Bastiani, F. (2019) Distributions for modeling location, scale, and shape: Using GAMLSS in R, Chapman and Hall/CRC, doi:10.1201/9780429298547. An older version can be found in <https://www.gamlss.com/>.

**Examples**

```

set.seed(123)
x <- rpareto(1, mu = 5)
d <- dpareto(x, mu = 5)
p <- ppareto(x, mu = 5)
q <- qpareto(p, mu = 5)

```

---

pgweibull

*Power generalized Weibull distribution*


---

**Description**

Survival, hazard, cumulative distribution, density, quantile and sampling function for the power generalized Weibull (PgW) distribution with parameters scale, shape and powershape.

**Usage**

```

spgweibull(q, scale = 1, shape = 1, powershape = 1, log = FALSE)
hpgweibull(x, scale = 1, shape = 1, powershape = 1, log = FALSE)
ppgweibull(q, scale = 1, shape = 1, powershape = 1,
           lower.tail = TRUE, log.p = FALSE)
dpgweibull(x, scale = 1, shape = 1, powershape = 1, log = FALSE)
qpgweibull(p, scale = 1, shape = 1, powershape = 1)
rpgweibull(n, scale = 1, shape = 1, powershape = 1)

```

**Arguments**

scale	positive scale parameter
shape	positive shape parameter
powershape	positive power shape parameter
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
x, q	vector of quantiles
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
p	vector of probabilities
n	number of observations

**Details**

The survival function of the PgW distribution is:

$$S(x) = \exp \left\{ 1 - \left[ 1 + \left( \frac{x}{\theta} \right)^\nu \right]^{\frac{1}{\gamma}} \right\}.$$

The hazard function is

$$\frac{\nu}{\gamma\theta^\nu} \cdot x^{\nu-1} \cdot \left[ 1 + \left( \frac{x}{\theta} \right)^\nu \right]^{\frac{1}{\gamma-1}}$$

The cumulative distribution function is then  $F(x) = 1 - S(x)$  and the density function is  $S(x) \cdot h(x)$ .

If both shape parameters equal 1, the PgW distribution reduces to the exponential distribution (see [dexp](#)) with rate = 1/scale. If the power shape parameter equals 1, the PgW distribution simplifies to the Weibull distribution (see [dweibull](#)) with the same parametrization.

**Value**

`dpgweibull` gives the density, `ppgweibull` gives the distribution function, `qpgweibull` gives the quantile function, and `rpgweibull` generates random deviates. `spgweibull` gives the survival function and `hpgweibull` gives the hazard function.

**Examples**

```
x <- rpgweibull(1, 2, 2, 3)
d <- dpgweibull(x, 2, 2, 3)
p <- ppgweibull(x, 2, 2, 3)
q <- qpgweibull(p, 2, 2, 3)
s <- spgweibull(x, 2, 2, 3)
h <- hpgweibull(x, 2, 2, 3)
```

---

powerexp

*Power Exponential distribution (PE and PE2)*

---

**Description**

Density, distribution function, quantile function, and random generation for the Power Exponential distribution (two versions).

**Usage**

```
dpowerexp(x, mu = 0, sigma = 1, nu = 2, log = FALSE)

ppowerexp(q, mu = 0, sigma = 1, nu = 2, lower.tail = TRUE, log.p = FALSE)

qpowerexp(p, mu = 0, sigma = 1, nu = 2, lower.tail = TRUE, log.p = FALSE)

rpowerexp(n, mu = 0, sigma = 1, nu = 2)
```

```

dpowerexp2(x, mu = 0, sigma = 1, nu = 2, log = FALSE)
ppowerexp2(q, mu = 0, sigma = 1, nu = 2, lower.tail = TRUE, log.p = FALSE)
qpowerexp2(p, mu = 0, sigma = 1, nu = 2, lower.tail = TRUE, log.p = FALSE)
rpowerexp2(n, mu = 0, sigma = 1, nu = 2)

```

### Arguments

x, q	vector of quantiles
mu	location parameter
sigma	scale parameter, must be positive
nu	shape parameter (real)
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$
p	vector of probabilities
n	number of random values to return

### Details

This implementation of the densities and distribution functions allow for automatic differentiation with RTMB while the other functions are imported from `gamlss.dist` package.

For `powerexp`, `mu` is the mean and `sigma` is the standard deviation while this does not hold for `powerexp2`.

See `gamlss.dist::PE` for more details.

### Value

`dpowerexp` gives the density, `ppowerexp` gives the distribution function, `qpowerexp` gives the quantile function, and `rpowerexp` generates random deviates.

### References

Rigby, R. A., Stasinopoulos, D. M., Heller, G. Z., and De Bastiani, F. (2019) Distributions for modeling location, scale, and shape: Using GAMLSS in R, Chapman and Hall/CRC, doi:10.1201/9780429298547. An older version can be found in <https://www.gamlss.com/>.

### Examples

```

# PE
x <- rpowerexp(1, mu = 0, sigma = 1, nu = 2)
d <- dpowerexp(x, mu = 0, sigma = 1, nu = 2)
p <- ppowerexp(x, mu = 0, sigma = 1, nu = 2)
q <- qpowerexp(p, mu = 0, sigma = 1, nu = 2)

# PE2

```

```
x <- rpowerexp2(1, mu = 0, sigma = 1, nu = 2)
d <- dpowerexp2(x, mu = 0, sigma = 1, nu = 2)
p <- ppowerexp2(x, mu = 0, sigma = 1, nu = 2)
q <- qpowerexp2(p, mu = 0, sigma = 1, nu = 2)
```

---

 rgmrf

*Sample from a multivariate Gaussian with a sparse precision matrix*


---

### Description

Draw samples from a multivariate Gaussian distribution specified by a sparse precision matrix. This is numerically efficient for high-dimensional but sparse systems.

### Usage

```
rgmrf(n, mean = 0, Q)
```

### Arguments

n	Number of samples to draw.
mean	Mean vector (or scalar, which will be recycled to match the dimension of Q).
Q	Sparse precision matrix ( $\Sigma^{-1}$ ).

### Value

A matrix of samples with rows corresponding to samples and columns to dimensions.

### Examples

```
rgmrf(3, mean = c(1, 2, 3), Q = Matrix::Diagonal(3))
```

---

 skellam

*Skellam distribution*


---

### Description

Probability mass function, distribution function, and random generation for the Skellam distribution.

### Usage

```
dskellam(x, mu1, mu2, log = FALSE)
```

```
rskellam(n, mu1, mu2)
```

**Arguments**

x	integer vector of counts
mu1, mu2	Poisson means
log	logical; return log-density if TRUE
n	number of random values to return.

**Details**

The Skellam distribution is the distribution of the difference of two Poisson random variables. Specifically, if  $X_1 \sim \text{Pois}(\mu_1)$  and  $X_2 \sim \text{Pois}(\mu_2)$ , then  $X_1 - X_2 \sim \text{Skellam}(\mu_1, \mu_2)$ .

This implementation of `dskellam` allows for automatic differentiation with RTMB.

**Value**

`dskellam` gives the probability mass function and `rskellam` generates random deviates.

**Examples**

```
x <- rskellam(1, 2, 3)
d <- dskellam(x, 2, 3)
```

---

skewnorm

*Skew normal distribution*


---

**Description**

Density, distribution function, quantile function, and random generation for the skew normal distribution.

**Usage**

```
dskewnorm(x, xi = 0, omega = 1, alpha = 0, log = FALSE)
pskewnorm(q, xi = 0, omega = 1, alpha = 0, ...)
qskewnorm(p, xi = 0, omega = 1, alpha = 0, ...)
rskewnorm(n, xi = 0, omega = 1, alpha = 0)
```

**Arguments**

x, q	vector of quantiles
xi	location parameter
omega	scale parameter, must be positive.
alpha	skewness parameter, +/- Inf is allowed.

log	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
...	additional parameters to be passed to the sn package functions for pskewnorm and qskewnorm.
p	vector of probabilities
n	number of random values to return

### Details

This implementation of dskewnorm allows for automatic differentiation with RTMB while the other functions are imported from the sn package. See `sn::dsn` for more details.

### Value

dskewnorm gives the density, pskewnorm gives the distribution function, qskewnorm gives the quantile function, and rskewnorm generates random deviates.

### See Also

[skewnorm2](#), [skewt](#), [skewt2](#)

### Examples

```
# alpha is skew parameter
x <- rskewnorm(1, alpha = 1)
d <- dskewnorm(x, alpha = 1)
p <- pskewnorm(x, alpha = 1)
q <- qskewnorm(p, alpha = 1)
```

---

skewnorm2

*Reparameterised skew normal distribution*

---

### Description

Density, distribution function, quantile function and random generation for the skew normal distribution reparameterised in terms of mean, standard deviation and skew magnitude

### Usage

```
dskewnorm2(x, mean = 0, sd = 1, alpha = 0, log = FALSE)
```

```
pskewnorm2(q, mean = 0, sd = 1, alpha = 0, ...)
```

```
qskewnorm2(p, mean = 0, sd = 1, alpha = 0, ...)
```

```
rskewnorm2(n, mean = 0, sd = 1, alpha = 0)
```

**Arguments**

x, q	vector of quantiles
mean	mean parameter
sd	standard deviation, must be positive.
alpha	skewness parameter, +/- Inf is allowed.
log	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
...	additional parameters to be passed to the sn package functions for pskewnorm and qskewnorm.
p	vector of probabilities
n	number of random values to return

**Details**

This implementation of dskewnorm2 allows for automatic differentiation with RTMB while the other functions are imported from the sn package.

**Value**

dskewnorm2 gives the density, pskewnorm2 gives the distribution function, qskewnorm2 gives the quantile function, and rskewnorm2 generates random deviates.

**See Also**

[skewnorm](#), [skewt](#), [skewt2](#)

**Examples**

```
# alpha is skew parameter
x <- rskewnorm2(1, alpha = 1)
d <- dskewnorm2(x, alpha = 1)
p <- pskewnorm2(x, alpha = 1)
q <- qskewnorm2(p, alpha = 1)
```

---

skewt

*Skewed students t distribution*


---

**Description**

Density, distribution function, quantile function, and random generation for the skew t distribution (type 2).

**Usage**

```
dskewt(x, mu = 0, sigma = 1, skew = 0, df = 100, log = FALSE)
```

```
pskewt(q, mu = 0, sigma = 1, skew = 0, df = 100,
       method = 0, lower.tail = TRUE, log.p = FALSE)
```

```
qskewt(p, mu = 0, sigma = 1, skew = 0, df = 100,
       tol = 1e-8, method = 0)
```

```
rskewt(n, mu = 0, sigma = 1, skew = 0, df = 100)
```

**Arguments**

x, q	vector of quantiles
mu	location parameter
sigma	scale parameter, must be positive.
skew	skewness parameter, can be positive or negative.
df	degrees of freedom, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
method	an integer value between 0 and 5 which selects the computing method; see ‘Details’ in the <a href="#">pst</a> documentation below for the meaning of these values. If method=0 (default value), an automatic choice is made among the four actual computing methods, depending on the other arguments.
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
tol	a scalar value which regulates the accuracy of the result of qsn, measured on the probability scale.
n	number of random values to return.

**Details**

This corresponds to the skew t type 2 distribution in GAMLSS ([ST2](#)), see pp. 411-412 of Rigby et al. (2019) and the version implemented in the `sn` package. This implementation of `dskewt` allows for automatic differentiation with RTMB while the other functions are imported from the `sn` package. See `sn::dst` for more details.

**Caution:** In a numerical optimisation, the skew parameter should NEVER be initialised with exactly zero. This will cause the initial and all subsequent derivatives to be exactly zero and hence the parameter will remain at its initial value.

**Value**

`dskewt` gives the density, `pskewt` gives the distribution function, `qskewt` gives the quantile function, and `rskewt` generates random deviates.

**See Also**

[skewt2](#), [skewnorm](#), [skewnorm2](#)

**Examples**

```
x <- rskewt(1, 1, 2, 5, 2)
d <- dskewt(x, 1, 2, 5, 2)
p <- pskewt(x, 1, 2, 5, 2)
q <- qskewt(p, 1, 2, 5, 2)
```

---

 skewt2

*Moment-parameterised skew t distribution*


---

**Description**

Density, distribution function, quantile function, and random generation for the skew t distribution reparameterised so that mean and sd correspond to the *true* mean and standard deviation.

**Usage**

```
dskewt2(x, mean = 0, sd = 1, skew = 0, df = 100, log = FALSE)
```

```
pskewt2(q, mean = 0, sd = 1, skew = 0, df = 100,
        method = 0, lower.tail = TRUE, log.p = FALSE)
```

```
qskewt2(p, mean = 0, sd = 1, skew = 0, df = 100, tol = 1e-08, method = 0)
```

```
rskewt2(n, mean = 0, sd = 1, skew = 0, df = 100)
```

**Arguments**

x, q	vector of quantiles
mean	mean parameter
sd	standard deviation parameter, must be positive.
skew	skewness parameter, can be positive or negative.
df	degrees of freedom, must be positive.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
method	an integer value between 0 and 5 which selects the computing method; see ‘Details’ in the <a href="#">pst</a> documentation below for the meaning of these values. If method=0 (default value), an automatic choice is made among the four actual computing methods, depending on the other arguments.
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
tol	a scalar value which regulates the accuracy of the result of qsn, measured on the probability scale.
n	number of random values to return.

### Details

This corresponds to the skew t type 2 distribution in GAMLSS (ST2), see pp. 411-412 of Rigby et al. (2019) and the version implemented in the `sn` package. However, it is reparameterised in terms of a standard deviation parameter `sd` rather than just a scale parameter `sigma`. Details of this reparameterisation are given below. This implementation of `dskewt` allows for automatic differentiation with RTMB while the other functions are imported from the `sn` package. See `sn:dst` for more details.

**Caution:** In a numerical optimisation, the skew parameter should NEVER be initialised with exactly zero. This will cause the initial and all subsequent derivatives to be exactly zero and hence the parameter will remain at its initial value.

For given skew =  $\alpha$  and df =  $\nu$ , define

$$\delta = \alpha / \sqrt{1 + \alpha^2}, \quad b_\nu = \sqrt{\nu/\pi} \Gamma((\nu - 1)/2) / \Gamma(\nu/2),$$

then

$$E(X) = \mu + \sigma \delta b_\nu, \quad Var(X) = \sigma^2 \left( \frac{\nu}{\nu - 2} - \delta^2 b_\nu^2 \right).$$

### Value

`dskewt2` gives the density, `pskewt2` gives the distribution function, `qskewt2` gives the quantile function, and `rskewt2` generates random deviates.

### See Also

[skewt](#), [skewnorm](#), [skewnorm2](#)

### Examples

```
x <- rskewt2(1, 1, 2, 5, 5)
d <- dskewt2(x, 1, 2, 5, 5)
p <- pskewt2(x, 1, 2, 5, 5)
q <- qskewt2(p, 1, 2, 5, 5)
```

### Description

Density, distribution function, quantile function, and random generation for the t distribution with location and scale parameters.

**Usage**

```
dt2(x, mu, sigma, df, log = FALSE)
```

```
pt2(q, mu, sigma, df)
```

```
rt2(n, mu, sigma, df)
```

```
qt2(p, mu, sigma, df)
```

```
pt(q, df)
```

**Arguments**

x, q	vector of quantiles
mu	location parameter
sigma	scale parameter, must be positive.
df	degrees of freedom, must be positive.
log	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
n	number of random values to return.
p	vector of probabilities

**Details**

This implementation of dt2 allows for automatic differentiation with RTMB.

**Value**

dt2 gives the density, pt2 gives the distribution function, qt2 gives the quantile function, and rt2 generates random deviates.

**Examples**

```
x <- rt2(1, 1, 2, 5)
d <- dt2(x, 1, 2, 5)
p <- pt2(x, 1, 2, 5)
q <- qt2(p, 1, 2, 5)
```

---

truncnorm

*Truncated normal distribution*

---

**Description**

Density, distribution function, quantile function, and random generation for the truncated normal distribution.

**Usage**

```
dtruncnorm(x, mean = 0, sd = 1, min = -Inf, max = Inf, log = FALSE)

ptruncnorm(q, mean = 0, sd = 1, min = -Inf, max = Inf,
           lower.tail = TRUE, log.p = FALSE)

qtruncnorm(p, mean = 0, sd = 1, min = -Inf, max = Inf,
           lower.tail = TRUE, log.p = FALSE)

rtruncnorm(n, mean = 0, sd = 1, min = -Inf, max = Inf)
```

**Arguments**

x, q	vector of quantiles
mean	mean parameter, must be positive.
sd	standard deviation parameter, must be positive.
min, max	truncation bounds.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
p	vector of probabilities
n	number of random values to return.

**Details**

This implementation of `dtruncnorm` allows for automatic differentiation with RTMB.

**Value**

`dtruncnorm` gives the density, `ptruncnorm` gives the distribution function, `qtruncnorm` gives the quantile function, and `rtruncnorm` generates random deviates.

**Examples**

```
x <- rtruncnorm(1, mean = 2, sd = 2, min = -1, max = 5)
d <- dtruncnorm(x, mean = 2, sd = 2, min = -1, max = 5)
p <- ptruncnorm(x, mean = 2, sd = 2, min = -1, max = 5)
q <- qtruncnorm(p, mean = 2, sd = 2, min = -1, max = 5)
```

---

trunct	<i>Truncated t distribution</i>
--------	---------------------------------

---

### Description

Density, distribution function, quantile function, and random generation for the truncated t distribution.

### Usage

```
dtrunct(x, df, min = -Inf, max = Inf, log = FALSE)
ptrunct(q, df, min = -Inf, max = Inf, lower.tail = TRUE, log.p = FALSE)
qtrunct(p, df, min = -Inf, max = Inf, lower.tail = TRUE, log.p = FALSE)
rtrunct(n, df, min = -Inf, max = Inf)
```

### Arguments

x, q	vector of quantiles
df	degrees of freedom parameter, must be positive.
min, max	truncation bounds.
log, log.p	logical; if TRUE, probabilities/densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
p	vector of probabilities
n	number of random values to return.

### Details

This implementation of `dtrunct` allows for automatic differentiation with RTMB.

### Value

`dtrunct` gives the density, `ptrunct` gives the distribution function, `qtrunct` gives the quantile function, and `rtrunct` generates random deviates.

### Examples

```
x <- rtrunct(1, df = 5, min = -1, max = 5)
d <- dtrunct(x, df = 5, min = -1, max = 5)
p <- ptrunct(x, df = 5, min = -1, max = 5)
q <- qtrunct(p, df = 5, min = -1, max = 5)
```

---

trunct2	<i>Truncated t distribution with location and scale</i>
---------	---

---

### Description

Density, distribution function, quantile function, and random generation for the truncated t distribution with location  $\mu$  and scale  $\sigma$ .

### Usage

```
dtrunct2(x, df, mu = 0, sigma = 1, min = -Inf, max = Inf, log = FALSE)
```

```
ptrunct2(q, df, mu = 0, sigma = 1, min = -Inf, max = Inf,
         lower.tail = TRUE, log.p = FALSE)
```

```
qtrunct2(p, df, mu = 0, sigma = 1, min = -Inf, max = Inf,
         lower.tail = TRUE, log.p = FALSE)
```

```
rtrunct2(n, df, mu = 0, sigma = 1, min = -Inf, max = Inf)
```

### Arguments

x, q	vector of quantiles
df	degrees of freedom parameter, must be positive.
mu	location parameter.
sigma	scale parameter, must be positive.
min, max	truncation bounds.
log, log.p	logical; if TRUE, probabilities/densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
p	vector of probabilities
n	number of random values to return.

### Details

This implementation of `dtrunct2` allows for automatic differentiation with RTMB.

### Value

`dtrunct2` gives the density, `ptrunct2` gives the distribution function, `qtrunct2` gives the quantile function, and `rtrunct2` generates random deviates.

### Examples

```
x <- rtrunct2(1, df = 5, mu = 2, sigma = 3, min = -1, max = 5)
d <- dtrunct2(x, df = 5, mu = 2, sigma = 3, min = -1, max = 5)
p <- ptrunct2(x, df = 5, mu = 2, sigma = 3, min = -1, max = 5)
q <- qtrunct2(p, df = 5, mu = 2, sigma = 3, min = -1, max = 5)
```

---

vm	<i>von Mises distribution</i>
----	-------------------------------

---

**Description**

Density, distribution function, and random generation for the von Mises distribution.

**Usage**

```
dvm(x, mu = 0, kappa = 1, log = FALSE)
pvm(q, mu = 0, kappa = 1, from = NULL, tol = 1e-20)
rvm(n, mu = 0, kappa = 1, wrap = TRUE)
```

**Arguments**

x, q	vector of angles measured in radians at which to evaluate the density function.
mu	mean direction of the distribution measured in radians.
kappa	non-negative numeric value for the concentration parameter of the distribution.
log	logical; if TRUE, densities are returned on the log scale.
from	value from which the integration for CDF starts. If NULL, is set to $\mu - \pi$ .
tol	the precision in evaluating the distribution function
n	number of random values to return.
wrap	logical; if TRUE, generated angles are wrapped to the interval from $-\pi$ to $\pi$ .

**Details**

This implementation of `dvm` allows for automatic differentiation with RTMB. `rvm` and `pvm` are simply wrappers of the corresponding functions from `circular`.

**Value**

`dvm` gives the density, `pvm` gives the distribution function, and `rvm` generates random deviates.

**Examples**

```
set.seed(1)
x <- rvm(10, 0, 1)
d <- dvm(x, 0, 1)
p <- pvm(x, 0, 1)
```

---

vmf *von Mises-Fisher distribution*

---

### Description

Density, distribution function, and random generation for the von Mises-Fisher distribution.

### Usage

```
dvmf(x, mu, kappa, log = FALSE)
```

```
rvmf(n, mu, kappa)
```

### Arguments

x	unit vector or matrix (with each row being a unit vector) of evaluation points
mu	unit mean vector
kappa	non-negative numeric value for the concentration parameter of the distribution.
log	logical; if TRUE, densities are returned on the log scale.
n	number of random values to return.

### Details

This implementation of `dvmf` allows for automatic differentiation with RTMB. `rvmf` is a reparameterised import from `movMF::rmovMF`.

### Value

`dvmf` gives the density and `rvm` generates random deviates.

### Examples

```
set.seed(123)
# single parameter set
mu <- rep(1, 3) / sqrt(3)
kappa <- 4
x <- rvmf(1, mu, kappa)
d <- dvmf(x, mu, kappa)

# vectorised over parameters
mu <- matrix(mu, nrow = 1)
mu <- mu[rep(1,10), ]
kappa <- rep(kappa, 10)
x <- rvmf(10, mu, kappa)
d <- dvmf(x, mu, kappa)
```

---

 vmf2

*Reparameterised von Mises-Fisher distribution*


---

**Description**

Density, distribution function, and random generation for the von Mises-Fisher distribution.

**Usage**

```
dvmf2(x, theta, log = FALSE)
```

```
rvmf2(n, theta)
```

**Arguments**

x	unit vector or matrix (with each row being a unit vector) of evaluation points
theta	direction and concentration vector. The direction of theta determines the mean direction on the sphere. The norm of theta is the concentration parameter of the distribution.
log	logical; if TRUE, densities are returned on the log scale.
n	number of random values to return.

**Details**

In this parameterisation,  $\theta = \kappa\mu$ , where  $\mu$  is a unit vector and  $\kappa$  is the concentration parameter.

dvmf2 allows for automatic differentiation with RTMB. rvmf2 is imported from movMF: : rmovMF.

**Value**

dvmf gives the density and rvm generates random deviates.

**Examples**

```
set.seed(123)
# single parameter set
theta <- c(1,2,3)
x <- rvmf2(1, theta)
d <- dvmf2(x, theta)

# vectorised over parameters
theta <- matrix(theta, nrow = 1)
theta <- theta[rep(1,10), ]
x <- rvmf2(10, theta)
d <- dvmf2(x, theta)
```

---

 wishart

*Wishart distribution*


---

**Description**

Density and random generation for the wishart distribution

**Usage**

```
dwishart(x, nu, Sigma, log = FALSE)
```

```
rwishart(n, nu, Sigma)
```

**Arguments**

x	positive definite $p \times p$ matrix or array of such matrices of dimension $p \times p \times n$ (for $n$ density evaluations)
nu	degrees of freedom, needs to be greater than $p - 1$
Sigma	scale matrix, needs to be positive definite and match the dimension of $x$ .
log	logical; if TRUE, densities $p$ are returned as $\log(p)$ .
n	number of random deviates to return

**Value**

dwishart gives the density, rwishart generates random deviates (matrix for  $n = 1$ , array with  $n$  slices for  $n > 1$ )

**Examples**

```
# single input: matrix-valued
x <- rwishart(1, nu = 5, Sigma = diag(3))
d <- dwishart(x, nu = 5, Sigma = diag(3))

# multiple inputs: array of matrices
x <- rwishart(4, nu = 5, Sigma = diag(3))
d <- dwishart(x, nu = 5, Sigma = diag(3))

# multiple inputs for x, nu and Sigma
nu <- c(7,5,8,9)
Sigma <- array(dim = c(3,3,4))
for(i in 1:4) Sigma[, , i] <- (i + 3) * diag(3)
x <- rwishart(4, nu, Sigma)
d <- dwishart(x, nu, Sigma)
```

---

wrpcauchy	<i>wrapped Cauchy distribution</i>
-----------	------------------------------------

---

**Description**

Density and random generation for the wrapped Cauchy distribution.

**Usage**

```
dwrpcauchy(x, mu = 0, rho, log = FALSE)
```

```
rwrpcauchy(n, mu = 0, rho, wrap = TRUE)
```

**Arguments**

x	vector of angles measured in radians at which to evaluate the density function.
mu	mean direction of the distribution measured in radians.
rho	concentration parameter of the distribution, must be in the interval from 0 to 1.
log	logical; if TRUE, densities are returned on the log scale.
n	number of random values to return.
wrap	logical; if TRUE, generated angles are wrapped to the interval from $-\pi$ to $\pi$ .

**Details**

This implementation of `dwrpcauchy` allows for automatic differentiation with RTMB. `rwrpcauchy` is simply a wrapper for `rwrappedcauchy` imported from `circular`.

**Value**

`dwrpcauchy` gives the density and `rwrpcauchy` generates random deviates.

**Examples**

```
set.seed(1)
x <- rwrpcauchy(10, 0, 0.5)
d <- dwrpcauchy(x, 0, 0.5)
```

---

zero_inflate	<i>Zero-inflated density constructor</i>
--------------	--

---

**Description**

Constructs a zero-inflated density function from a given probability density function

**Usage**

```
zero_inflate(dist, discrete = NULL)
```

**Arguments**

dist	either a probability density function or a probability mass function
discrete	logical; if TRUE, the density for $x = 0$ will be <code>zeroprob + (1-zeroprob) * dist(0, ...)</code> . Otherwise it will just be <code>zeroprob</code> . In standard cases, this will be determined automatically. For non-standard cases, set this to TRUE or FALSE depending on the type of <code>dist</code> . See details.

**Details**

The definition of zero-inflation is different for discrete and continuous distributions. For discrete distributions with p.m.f.  $f$  and zero-inflation probability  $p$ , we have

$$\Pr(X = 0) = p + (1 - p) \cdot f(0),$$

and

$$\Pr(X = x) = (1 - p) \cdot f(x), \quad x > 0.$$

For continuous distributions with p.d.f.  $f$ , we have

$$f_{\text{zinfl}}(x) = p \cdot \delta_0(x) + (1 - p) \cdot f(x),$$

where  $\delta_0$  is the Dirac delta function at zero.

**Value**

zero-inflated density function with first argument  $x$ , second argument `zeroprob`, and additional arguments ... that will be passed to `dist`.

**Examples**

```
# Zero-inflated normal distribution
dzinorm <- zero_inflate(dnorm)
dzinorm(c(NA, 0, 2), 0.5, mean = 1, sd = 1)

# Zero-inflated Poisson distribution
zipois <- zero_inflate(dpois)
zipois(c(NA, 0, 1), 0.5, 1)
```

```
# Non-standard case: Zero-inflated reparametrised beta distribution
dzibeta2 <- zero_inflate(dbeta2, discrete = FALSE)
```

---

zibeta	<i>Zero-inflated beta distribution</i>
--------	--

---

### Description

Density, distribution function, and random generation for the zero-inflated beta distribution.

### Usage

```
dzibeta(x, shape1, shape2, zeroprob = 0, log = FALSE)

pzibeta(q, shape1, shape2, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)

rzibeta(n, shape1, shape2, zeroprob = 0)
```

### Arguments

x, q	vector of quantiles
shape1, shape2	non-negative shape parameters of the beta distribution
zeroprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

### Details

This implementation allows for automatic differentiation with RTMB.

### Value

dzibeta gives the density, pzibeta gives the distribution function, and rzibeta generates random deviates.

### Examples

```
set.seed(123)
x <- rzibeta(1, 2, 2, 0.5)
d <- dzibeta(x, 2, 2, 0.5)
p <- pzibeta(x, 2, 2, 0.5)
```

zibeta2

*Reparameterised zero-inflated beta distribution***Description**

Density, distribution function, and random generation for the zero-inflated beta distribution reparameterised in terms of mean and concentration.

**Usage**

```
dzibeta2(x, mu, phi, zeroprob = 0, log = FALSE)
```

```
pzibeta2(q, mu, phi, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
rzibeta2(n, mu, phi, zeroprob = 0)
```

**Arguments**

x, q	vector of quantiles
mu	mean parameter, must be in the interval from 0 to 1.
phi	concentration parameter, must be positive.
zeroprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise $P[X > x]$ .
n	number of random values to return.
p	vector of probabilities

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

dzibeta2 gives the density, pzibeta2 gives the distribution function, and rzibeta2 generates random deviates.

**Examples**

```
set.seed(123)
x <- rzibeta2(1, 0.5, 1, 0.5)
d <- dzibeta2(x, 0.5, 1, 0.5)
p <- pzibeta2(x, 0.5, 1, 0.5)
```

---

zibinom	<i>Zero-inflated binomial distribution</i>
---------	--

---

### Description

Probability mass function, distribution function, and random generation for the zero-inflated binomial distribution.

### Usage

```
dzibinom(x, size, prob, zeroprob = 0, log = FALSE)
pzibinom(q, size, prob, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)
rzibinom(n, size, prob, zeroprob = 0)
```

### Arguments

x, q	vector of quantiles
size	number of trials (zero or more).
prob	probability of success on each trial.
zeroprob	zero-inflation probability between 0 and 1
log, log.p	logical; return log-density if TRUE
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

### Details

This implementation allows for automatic differentiation with RTMB.

### Value

dzibinom gives the probability mass function, pzibinom gives the distribution function, and rzibinom generates random deviates.

### Examples

```
set.seed(123)
x <- rzibinom(1, size = 10, prob = 0.5, zeroprob = 0.5)
d <- dzibinom(x, size = 10, prob = 0.5, zeroprob = 0.5)
p <- pzibinom(x, size = 10, prob = 0.5, zeroprob = 0.5)
```

---

zigamma	<i>Zero-inflated gamma distribution</i>
---------	---

---

### Description

Density, distribution function, and random generation for the zero-inflated gamma distribution.

### Usage

```
dzigamma(x, shape, scale, zeroprob = 0, log = FALSE)
```

```
pzigamma(q, shape, scale, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
rzigamma(n, shape, scale, zeroprob = 0)
```

### Arguments

x, q	vector of quantiles
shape	positive shape parameter
scale	positive scale parameter
zeroprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return

### Details

This implementation allows for automatic differentiation with RTMB.

### Value

`dzigamma` gives the density, `pzigamma` gives the distribution function, and `rzigamma` generates random deviates.

### Examples

```
x <- rzigamma(1, 1, 1, 0.5)
d <- dzigamma(x, 1, 1, 0.5)
p <- pzigamma(x, 1, 1, 0.5)
```

---

`zigamma2`*Zero-inflated and reparameterised gamma distribution*

---

### Description

Density, distribution function, and random generation for the zero-inflated gamma distribution reparameterised in terms of mean and standard deviation.

### Usage

```
dzigamma2(x, mean = 1, sd = 1, zeroprob = 0, log = FALSE)
```

```
pzigamma2(q, mean = 1, sd = 1, zeroprob = 0)
```

```
rzigamma2(n, mean = 1, sd = 1, zeroprob = 0)
```

### Arguments

<code>x, q</code>	vector of quantiles
<code>mean</code>	mean parameter, must be positive.
<code>sd</code>	standard deviation parameter, must be positive.
<code>zeroprob</code>	zero-inflation probability between 0 and 1.
<code>log</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>n</code>	number of random values to return

### Details

This implementation allows for automatic differentiation with RTMB.

### Value

`dzigamma2` gives the density, `pzigamma2` gives the distribution function, and `rzigamma` generates random deviates.

### Examples

```
x <- rzigamma2(1, 2, 1, 0.5)
d <- dzigamma2(x, 2, 1, 0.5)
p <- pzigamma2(x, 2, 1, 0.5)
```

---

`ziinvgauss`*Zero-inflated inverse Gaussian distribution*

---

### Description

Density, distribution function, and random generation for the zero-inflated inverse Gaussian distribution.

### Usage

```
dziinvgauss(x, mean = 1, shape = 1, zeroprob = 0, log = FALSE)
```

```
pziinvgauss(q, mean = 1, shape = 1, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
rziinvgauss(n, mean = 1, shape = 1, zeroprob = 0)
```

### Arguments

<code>x, q</code>	vector of quantiles
<code>mean</code>	location parameter
<code>shape</code>	shape parameter, must be positive.
<code>zeroprob</code>	zero-probability, must be in $[0, 1]$ .
<code>log, log.p</code>	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
<code>lower.tail</code>	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
<code>n</code>	number of random values to return

### Details

This implementation of `zidinvgauss` allows for automatic differentiation with RTMB.

### Value

`dziinvgauss` gives the density, `pziinvgauss` gives the distribution function, and `rziinvgauss` generates random deviates.

### Examples

```
x <- rziinvgauss(1, 1, 2, 0.5)
d <- dziinvgauss(x, 1, 2, 0.5)
p <- pziinvgauss(x, 1, 2, 0.5)
```

---

zilnorm	<i>Zero-inflated log normal distribution</i>
---------	--

---

### Description

Density, distribution function, and random generation for the zero-inflated log normal distribution.

### Usage

```
dzilnorm(x, meanlog = 0, sdlog = 1, zeroprob = 0, log = FALSE)

pzilnorm(q, meanlog = 0, sdlog = 1, zeroprob = 0,
         lower.tail = TRUE, log.p = FALSE)

rzilnorm(n, meanlog = 0, sdlog = 1, zeroprob = 0)

plnorm(q, meanlog = 0, sdlog = 1, lower.tail = TRUE, log.p = FALSE)
```

### Arguments

x, q	vector of quantiles
meanlog, sdlog	mean and standard deviation of the distribution on the log scale with default values of 0 and 1 respectively.
zeroprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return

### Details

This implementation allows for automatic differentiation with RTMB.

### Value

dzilnorm gives the density, pzilnorm gives the distribution function, and rzilnorm generates random deviates.

### Examples

```
x <- rzilnorm(1, 1, 1, 0.5)
d <- dzilnorm(x, 1, 1, 0.5)
p <- pzilnorm(x, 1, 1, 0.5)
```

zinbinom

*Zero-inflated negative binomial distribution***Description**

Probability mass function, distribution function, quantile function, and random generation for the zero-inflated negative binomial distribution.

**Usage**

```
dzinbinom(x, size, prob, zeroprof = 0, log = FALSE)
```

```
pzinbinom(q, size, prob, zeroprof = 0, lower.tail = TRUE, log.p = FALSE)
```

```
rzinbinom(n, size, prob, zeroprof = 0)
```

**Arguments**

x, q	vector of (non-negative integer) quantiles
size	size parameter, must be positive.
prob	mean parameter, must be positive.
zeroprof	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.
p	vector of probabilities

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

dzinbinom gives the density, pzinbinom gives the distribution function, and rzinbinom generates random deviates.

**Examples**

```
set.seed(123)
x <- rzinbinom(1, size = 2, prob = 0.5, zeroprof = 0.5)
d <- dzinbinom(x, size = 2, prob = 0.5, zeroprof = 0.5)
p <- pzinbinom(x, size = 2, prob = 0.5, zeroprof = 0.5)
```

---

zinbinom2

*Zero-inflated and reparameterised negative binomial distribution*


---

**Description**

Probability mass function, distribution function, quantile function and random generation for the zero-inflated negative binomial distribution reparameterised in terms of mean and size.

**Usage**

```
dzinbinom2(x, mu, size, zeroprob = 0, log = FALSE)
```

```
pzinbinom2(q, mu, size, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
rzinbinom2(n, mu, size, zeroprob = 0)
```

**Arguments**

x, q	vector of (non-negative integer) quantiles
mu	mean parameter, must be positive.
size	size parameter, must be positive.
zeroprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.
p	vector of probabilities

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

dzinbinom2 gives the density, pzinbinom2 gives the distribution function, and rzinbinom2 generates random deviates.

**Examples**

```
set.seed(123)
x <- rzinbinom2(1, 2, 1, zeroprob = 0.5)
d <- dzinbinom2(x, 2, 1, zeroprob = 0.5)
p <- pzinbinom2(x, 2, 1, zeroprob = 0.5)
```

---

`zipois`*Zero-inflated Poisson distribution*

---

**Description**

Probability mass function, distribution function, and random generation for the zero-inflated Poisson distribution.

**Usage**

```
dzipois(x, lambda, zeroprob = 0, log = FALSE)
```

```
pzipois(q, lambda, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
rzipois(n, lambda, zeroprob = 0)
```

**Arguments**

<code>x, q</code>	integer vector of counts
<code>lambda</code>	vector of (non-negative) means
<code>zeroprob</code>	zero-inflation probability between 0 and 1
<code>log, log.p</code>	logical; return log-density if TRUE
<code>lower.tail</code>	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
<code>n</code>	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

`dzipois` gives the probability mass function, `pzipois` gives the distribution function, and `rzipois` generates random deviates.

**Examples**

```
set.seed(123)
x <- rzipois(1, 0.5, 1)
d <- dzipois(x, 0.5, 1)
p <- pzipois(x, 0.5, 1)
```

---

ziweibull	<i>Zero-inflated Weibull distribution</i>
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### Description

Density, distribution function, and random generation for the zero-inflated Weibull distribution.

### Usage

```
dziweibull(x, shape, scale, zeroprob = 0, log = FALSE)
```

```
pziweibull(q, shape, scale, zeroprob = 0, lower.tail = TRUE, log.p = FALSE)
```

```
rziweibull(n, shape, scale, zeroprob = 0)
```

### Arguments

x, q	vector of quantiles
shape	positive shape parameter
scale	positive scale parameter
zeroprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return

### Details

This implementation allows for automatic differentiation with RTMB.

### Value

dziweibull gives the density, pziweibull gives the distribution function, and rziweibull generates random deviates.

### Examples

```
x <- rziweibull(1, 1, 1, 0.5)
d <- dziweibull(x, 1, 1, 0.5)
p <- pziweibull(x, 1, 1, 0.5)
```

zoibeta

*Zero- and one-inflated beta distribution***Description**

Density, distribution function, and random generation for the zero-one-inflated beta distribution.

**Usage**

```
dzoibeta(x, shape1, shape2, zeroprob = 0, oneprob = 0, log = FALSE)
pzoibeta(q, shape1, shape2, zeroprob = 0, oneprob = 0,
         lower.tail = TRUE, log.p = FALSE)
rzoibeta(n, shape1, shape2, zeroprob = 0, oneprob = 0)
```

**Arguments**

x, q	vector of quantiles
shape1, shape2	non-negative shape parameters of the beta distribution
zeroprob	zero-inflation probability between 0 and 1.
oneprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

dzoibeta gives the density, pzoibeta gives the distribution function, and rzoibeta generates random deviates.

**Examples**

```
set.seed(123)
x <- rzoibeta(1, 2, 2, 0.2, 0.3)
d <- dzoibeta(x, 2, 2, 0.2, 0.3)
p <- pzoibeta(x, 2, 2, 0.2, 0.3)
```

zoibeta2

*Reparameterised zero- and one-inflated beta distribution***Description**

Density, distribution function, and random generation for the zero-one-inflated beta distribution reparameterised in terms of mean and concentration.

**Usage**

```
dzoibeta2(x, mu, phi, zeroprob = 0, oneprob = 0, log = FALSE)
```

```
pzoibeta2(q, mu, phi, zeroprob = 0, oneprob = 0,
          lower.tail = TRUE, log.p = FALSE)
```

```
rzoibeta2(n, mu, phi, zeroprob = 0, oneprob = 0)
```

**Arguments**

x, q	vector of quantiles
mu	mean parameter, must be in the interval from 0 to 1.
phi	concentration parameter, must be positive.
zeroprob	zero-inflation probability between 0 and 1.
oneprob	zero-inflation probability between 0 and 1.
log, log.p	logical; if TRUE, probabilities/ densities $p$ are returned as $\log(p)$ .
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

**Value**

dzoibeta2 gives the density, pzoibeta2 gives the distribution function, and rzoibeta2 generates random deviates.

**Examples**

```
set.seed(123)
x <- rzoibeta2(1, 0.6, 2, 0.2, 0.3)
d <- dzoibeta2(x, 0.6, 2, 0.2, 0.3)
p <- pzoibeta2(x, 0.6, 2, 0.2, 0.3)
```

ztbinom

*Zero-truncated Binomial distribution***Description**

Probability mass function, distribution function, and random generation for the zero-truncated Binomial distribution.

**Usage**

```
dzttbinom(x, size, prob, log = FALSE)
```

```
pztbinom(q, size, prob, lower.tail = TRUE, log.p = FALSE)
```

```
rztbinom(n, size, prob)
```

**Arguments**

x, q	integer vector of counts
size	number of trials
prob	success probability in each trial
log, log.p	logical; return log-density if TRUE
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

By definition, this distribution only has support on the positive integers (1, ..., size). Any zero-truncated distribution is defined as

$$P(X = x | X > 0) = P(X = x) / (1 - P(X = 0)),$$

where  $P(X = x)$  is the probability mass function of the corresponding untruncated distribution.

**Value**

dzttbinom gives the probability mass function, pztbinom gives the distribution function, and rztbinom generates random deviates.

**Examples**

```
set.seed(123)
x <- rztbinom(1, size = 10, prob = 0.3)
d <- dzttbinom(x, size = 10, prob = 0.3)
p <- pztbinom(x, size = 10, prob = 0.3)
```

---

ztnbinom *Zero-truncated Negative Binomial distribution*

---

### Description

Probability mass function, distribution function, and random generation for the zero-truncated Negative Binomial distribution.

### Usage

```
dztnbinom(x, size, prob, log = FALSE)

pztnbinom(q, size, prob, lower.tail = TRUE, log.p = FALSE)

rztnbinom(n, size, prob)
```

### Arguments

x, q	integer vector of counts
size	target for number of successful trials, or dispersion parameter (the shape parameter of the gamma mixing distribution). Must be strictly positive, need not be integer.
prob	probability of success in each trial. $0 < \text{prob} \leq 1$ .
log, log.p	logical; return log-density if TRUE
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

### Details

This implementation allows for automatic differentiation with RTMB.

By definition, this distribution only has support on the positive integers (1, 2, ...). Any zero-truncated distribution is defined as

$$P(X = x | X > 0) = P(X = x) / (1 - P(X = 0)),$$

where  $P(X = x)$  is the probability mass function of the corresponding untruncated distribution.

### Value

dztnbinom gives the probability mass function, pztnbinom gives the distribution function, and rztnbinom generates random deviates.

### Examples

```
set.seed(123)
x <- rztnbinom(1, size = 2, prob = 0.5)
d <- dztnbinom(x, size = 2, prob = 0.5)
p <- pztnbinom(x, size = 2, prob = 0.5)
```

---

ztnbinom2

*Reparameterised zero-truncated negative binomial distribution*


---

**Description**

Probability mass function, distribution function, quantile function, and random generation for the zero-truncated negative binomial distribution reparameterised in terms of mean and size.

**Usage**

```
dztnbinom2(x, mu, size, log = FALSE)
pztnbinom2(q, mu, size, lower.tail = TRUE, log.p = FALSE)
rztnbinom2(n, mu, size)
```

**Arguments**

x, q	integer vector of counts
mu	mean parameter, must be positive
size	size/dispersion parameter, must be positive
log, log.p	logical; return log-density if TRUE
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

By definition, this distribution only has support on the positive integers (1, 2, ...). Any zero-truncated distribution is defined as

$$P(X = x | X > 0) = P(X = x) / (1 - P(X = 0)),$$

where  $P(X = x)$  is the probability mass function of the corresponding untruncated distribution.

**Value**

dztnbinom2 gives the probability mass function, pztnbinom2 gives the distribution function, and rztnbinom2 generates random deviates.

**Examples**

```
set.seed(123)
x <- rztnbinom2(1, mu = 2, size = 1)
d <- dztnbinom2(x, mu = 2, size = 1)
p <- pztnbinom2(x, mu = 2, size = 1)
```

ztpois

*Zero-truncated Poisson distribution***Description**

Probability mass function, distribution function, and random generation for the zero-truncated Poisson distribution.

**Usage**

```
dztpois(x, lambda, log = FALSE)
```

```
pztpois(q, lambda, lower.tail = TRUE, log.p = FALSE)
```

```
rztpois(n, lambda)
```

**Arguments**

x, q	integer vector of counts
lambda	vector of (non-negative) means
log, log.p	logical; return log-density if TRUE
lower.tail	logical; if TRUE, probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .
n	number of random values to return.

**Details**

This implementation allows for automatic differentiation with RTMB.

By definition, this distribution only has support on the positive integers (1, 2, ...). Any zero-truncated distribution is defined as

$$P(X = x | X > 0) = P(X = x) / (1 - P(X = 0)),$$

where  $P(X = x)$  is the probability mass function of the corresponding untruncated distribution.

**Value**

dztpois gives the probability mass function, pztpois gives the distribution function, and rztpois generates random deviates.

**Examples**

```
set.seed(123)
x <- rztpois(1, 0.5)
d <- dztpois(x, 0.5)
p <- pztpois(x, 0.5)
```

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