Package ‘mnormt’

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Author Fortran code by Alan Genz, R code by Adelchi Azzalini
Maintainer Adelchi Azzalini <adelchi.azzalini@unipd.it>
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Description Functions are provided for computing the density and the
distribution function of multivariate normal and ”t” random variables,
and for generating random vectors sampled from these distributions.
Probabilities are computed via non-Monte Carlo methods; different routines
are used in the case d=1, d=2, d>2, if d denotes the number of dimensions.
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The 'mnormt' package: summary information

Description

This package provides functions for computing the density and the distribution function of multivariate normal and multivariate Student’s t variates and for generating random vectors sampled from these distributions.

Details

Probabilities are computed via a non-Monte Carlo method. Different routines are used in the three cases $d=1$, $d=2$, $d>2$, if $d$ denotes the number of dimensions.

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Author(s)

Adelchi Azzalini (R code and package creation) and Alan Genz (Fortran code, see references below; this includes routines of other authors)

References


Genz, A.: Fortran code available at http://www.math.wsu.edu/math/faculty/genz/software/fort77/mvn.f

Multivariate normal distribution

Description

The probability density function, the distribution function and random number generation for the multivariate normal (Gaussian) distribution.
Usage

dmnorm(x, mean = rep(0, d), varcov, log = FALSE)

pmnorm(x, mean = rep(0, d), varcov, ...)

rmnorm(n = 1, mean = rep(0, d), varcov, sqrt=NULL)

sadmvn(lower, upper, mean, varcov, maxpts = 2000*d, abseps = 1e-06, releps = 0)

Arguments

x

either a vector of length d or a matrix with d columns, where d=ncol(varcov),
representing the coordinates of the point(s) where the density must be evaluated;
for pmnorm, d cannot exceed 20.

mean

either a vector of length d, representing the mean value, or (except for rmnorm)
a matrix whose rows represent different mean vectors; in the matrix case, only
allowed for dmnorm and pmnorm, its dimensions must match those of x.

varcov

a symmetric positive-definite matrix representing the variance-covariance ma-
trix of the distribution; a vector of length 1 is also allowed (in this case, d=1 is set).

sqrt

if not NULL (default value is NULL), a square root of the intended varcov matrix;
see ‘Details’ for a full description.

log

a logical value (default value is FALSE); if TRUE, the logarithm of the density is
computed.

... parameters passed to sadmvn, among maxpts, abseps, releps.

n

the number of random vectors to be generated.

lower

a numeric vector of lower integration limits of the density function; must be of
maximal length 20; +Inf and -Inf entries are allowed.

upper

a numeric vector of upper integration limits of the density function; must be of
maximal length 20; +Inf and -Inf entries are allowed.

maxpts

the maximum number of function evaluations (default value: 2000*d).

abseps

absolute error tolerance (default value: 1e-6).

releps

relative error tolerance (default value: 0).

Details

The function pmnorm works by making a suitable call to sadmvn if d^R
or to bivnt.prob if d=2, or to pnorm if d=1. Function sadmvn is an interface to a Fortran-77 routine with the same name written by
Alan Genz, available from his web page, which works using an adaptive integration method.
This Fortran-77 routine makes uses of some auxiliary functions whose authors are documented in
the code.

If sqrt=NULL (default value), the working of rmnorm involves computation of a square root of
varcov via the Cholesky decomposition. If a non-NULL value of sqrt is supplied, it is assumed that
it represents a matrix, R say, such that R'R represents the required variance-covariance matrix of the
distribution; in this case, the argument varcov is ignored. This mechanism is intended primarily for
use in a sequence of calls to rmnorm, all sampling from a distribution with fixed variance matrix; a
suitable matrix sqrt can then be computed only once beforehand, avoiding that the same operation
is repeated multiple times along the sequence of calls; see the examples below. Another use of
sqrt is to supply a different form of square root of the variance-covariance matrix, in place of the Cholesky factor.

For efficiency reasons, rmnorm does not perform checks on the supplied arguments.

If, after setting the same seed value to set.seed, two calls to rmnorm are made with the same arguments except that one generates n1 vectors and the other n2 vectors, with n1<n2, then the n1 vectors of the first call coincide with the initial n2 vectors of the second call.

Value

dmnorm returns a vector of density values (possibly log-transformed); pmnorm returns a vector of probabilities, possibly with attributes on the accuracy in case x is a vector; sadmvn return a single probability with attributes giving details on the achieved accuracy; rmnorm returns a matrix of n rows of random vectors or a vector in case n=1.

Note

The attributes error and status of the probability returned by pmnorm and sadmvn indicate whether the function had a normal termination, achieving the required accuracy. If this is not the case, re-run the function with a higher value of maxpts.

Author(s)

Fortran code of SADMVN and most auxiliary functions by Alan Genz, some additional auxiliary functions by people referred to within his program. Interface to R and additional R code by Adelchi Azzalini.

References


Genz, A.: Fortran code available at [http://www.math.wsu.edu/math/faculty/genz/software/fort77/mvn.f](http://www.math.wsu.edu/math/faculty/genz/software/fort77/mvn.f)

See Also

dnorm, dmt, biv.nt.prob

Examples

```r
x <- seq(-2, 4, length=21)
y <- cos(2*x) + 10
z <- x + sin(3*y)
mu <- c(1,12,2)
Sigma <- matrix(c(1,2,0,2,5,0.5,0,0.5,3), 3, 3)
f <- dmnorm(cbind(x,y,z), mu, Sigma)
f0 <- dmnorm(mu, mu, Sigma)
p1 <- pmnorm(c(2,11,3), mu, Sigma)
```
dmt <- pmnorm(c(2,1,1,3), mu, Sigma, maxpts=10000, abseps=1e-10)
p <- pmnorm(cbind(x,y,z), mu, Sigma)
#
set.seed(123)
x1 <- rmnorm(5, mu, Sigma)
set.seed(123)
x2 <- rmnorm(5, mu, sqrt=chol(Sigma)) # x1=x2
eig <- eigen(Sigma, symmetric = TRUE)
R <- t(eig$vectors)%% diag(sqrt(eig$values))
for(i in 1:50) x <- rmnorm(5, mu, sqrt=R)
#
p <- sadmvn(lower=c(2,1,1,3), upper=rep(Inf,3), mu, Sigma) # upper tail
#
p0 <- pmnorm(c(2,1), mu[1:2], Sigma[1:2,1:2])
p1 <- bivNntNprob(0, lower=rep(-Inf,2), upper=c(2, 11), mu[1:2], Sigma[1:2,1:2])
p2 <- sadmvn(lower=rep(-Inf,2), upper=c(2, 11), mu[1:2], Sigma[1:2,1:2])
c(p0, p1, p2, p0-p1, p0-p2)
#
p1 <- pnorm(0, 1, 3)
p2 <- pmnorm(0, 1, 3^2)

---

dmt

Multivariate t distribution

Description

The probability density function, the distribution function and random number generation for the multivariate Student’s t distribution

Usage

dmt(x, mean = rep(0, d), S, df=Inf, log = FALSE)
pmt(x, mean = rep(0, d), S, df=Inf, ...)
rmt(n = 1, mean = rep(0, d), S, df=Inf, sqrt=NULL)
sadmvt(df, lower, upper, mean, S, maxpts = 2000*d, abseps = 1e-06, releps = 0)
biv.nt.prob(df, lower, upper, mean, S)

Arguments

x

either a vector of length d or a matrix with d columns, where d=ncol(S), giving the coordinates of the point(s) where the density must be evaluated; for pmt, d cannot exceed 20.

mean

either a vector of length d, representing the location parameter (equal to the mean vector when df>1) or a matrix whose rows represent different mean vectors (except for rmt); in the matrix case, its dimensions must match those of x.

S

a symmetric positive-definite matrix representing the scale matrix of the distribution, such that S*df/(df-2) is the variance-covariance matrix when df>2; a vector of length 1 is also allowed (in this case, d=1 is set).
df  degrees of freedom; it must be a positive integer for \( \text{pmt} \), \( \text{sadmvt} \) and \( \text{biv.n.t.prob} \), otherwise a positive number. If \( \text{df} = \text{Inf} \) (default value), the corresponding \*nnorm function is called, unless \( \text{d} = 2 \); in this case \( \text{biv.n.t.prob} \) is used. If \( \text{biv.n.t.prob} \) is called with \( \text{df} = \text{Inf} \), it returns the probability of a rectangle assigned by a bivariate normal distribution.

log  a logical value (default value is \text{FALSE}); if \text{TRUE}, the logarithm of the density is computed.

sqrt  if not \text{NULL} (default value is \text{NULL}), a square root of the intended scale matrix \( S \); see ‘Details’ for a full description.

...  parameters passed to \( \text{sadmvt} \), among \text{maxpts}, \text{absrel}, \text{releps}.

n  the number of random vectors to be generated

lower  a numeric vector of lower integration limits of the density function; must be of maximal length \( 2d \); \(+\text{Inf} \) and \(-\text{Inf} \) entries are allowed.

upper  a numeric vector of upper integration limits of the density function; must be of maximal length \( 2d \); \(+\text{Inf} \) and \(-\text{Inf} \) entries are allowed.

maxpts  the maximum number of function evaluations (default value: \( 2000d \))

abseps  absolute error tolerance (default value: \( 1e-6 \)).

releps  relative error tolerance (default value: \( 0 \)).

Details

The functions \( \text{sadmvt} \) and \( \text{biv.n.t.prob} \) are interfaces to Fortran-77 routines by Alan Genz, and available from his web page; they make use of some auxiliary functions whose authors are documented in the Fortran code. The routine \( \text{sadmvt} \) uses an adaptive integration method. The routine \( \text{biv.n.t.prob} \) is specific for the bivariate case; if \( \text{df} < 1 \) or \( \text{df} = \text{Inf} \), it computes the bivariate normal distribution function using a non-iterative method described in a reference given below. If \( \text{pmt} \) is called with \( \text{d} = 2 \), this is converted into a suitable call to \( \text{sadmvt} \); if \( \text{d} = 1 \), a call to \( \text{biv.n.t.prob} \) is used; if \( \text{d} = 1 \), then \( \text{pt} \) is used.

If \( \text{sqrt} = \text{NULL} \) (default value), the working of \( \text{rmt} \) involves computation of a square root of \( S \) via the Cholesky decomposition. If a non-\text{NULL} value of \( \text{sqrt} \) is supplied, it is assumed that it represents a square root of the scale matrix, otherwise represented by \( S \), whose value is ignored in this case. This mechanism is intended primarily for use in a sequence of calls to \( \text{rmt} \), all sampling from a distribution with fixed scale matrix; a suitable matrix \( \text{sqrt} \) can then be computed only once beforehand, avoiding that the same operation is repeated multiple times along the sequence of calls. For examples of use of this argument, see those in the documentation of \( \text{rmnorm} \). Another use of \( \text{sqrt} \) is to supply a different form of square root of the scale matrix, in place of the Cholesky factor.

For efficiency reasons, \( \text{rmt} \) does not perform checks on the supplied arguments.

Value

\( \text{dmt} \) returns a vector of density values (possibly log-transformed); \( \text{pmt} \) and \( \text{sadmvt} \) return a single probability with attributes giving details on the achieved accuracy, provided \( x \) of \( \text{pmmnorm} \) is a vector; \( \text{rmt} \) returns a matrix of \( n \) rows of random vectors.
Note
The attributes error and status of the probability returned by sadmv and by pmt (the latter only if x is a vector and d>=2) indicate whether the function had a normal termination, achieving the required accuracy. If this is not the case, re-run the function with a higher value of maxpts.

Author(s)
Fortran code of SADMVT and most auxiliary functions by Alan Genz, some additional auxiliary functions by people referred to within his program; interface to R and additional R code by Adelchi Azzalini.

References

See Also
dt, rmnorm for use of argument sqrt

Examples
x <- seq(-2,4,length=21)
y <- 2*x+10
z <- x+cos(y)
mu <- c(1,12,2)
Sigma <- matrix(c(1,2,0,2,5,0,5,0,0.5,3), 3, 3)
df <- 4
f <- dmt(cbind(x,y,z), mu, Sigma, df)
p1 <- pmt(c(2,11,3), mu, Sigma, df)
p2 <- pmt(c(2,11,3), mu, Sigma, df, maxpts=10000, abseps=1e-8)
x <- rmt(10, mu, Sigma, df)
p <- sadmv(df, lower=c(2,11,3), upper=rep(Inf,3), mu, Sigma) # upper tail
#
p0 <- pmt(c(2,11), mu[1:2], Sigma[1:2,1:2], df=5)
p1 <- biv.nt.prob(5, lower=rep(-Inf,2), upper=c(2, 11), mu[1:2], Sigma[1:2,1:2])
p2 <- sadmv(5, lower=rep(-Inf,2), upper=c(2, 11), mu[1:2], Sigma[1:2,1:2])
c(p0, p1, p2, p0-p1, p0-p2)
**Usage**

```r
pd.solve(x, silent = FALSE, log.det=FALSE)
```

**Arguments**

- `x`: a symmetric positive-definite matrix.
- `silent`: a logical value which indicates the action to take in case of an error. If `silent==TRUE` and an error occurs, the function silently returns a NULL value; if `silent==FALSE` (default), an error generates a stop with an error message.
- `log.det`: a logical value to indicate whether the log-determinant of `x` is required (default is FALSE).

**Details**

The function checks that `x` is a symmetric positive-definite matrix. If an error is detected, an action is taken which depends on the value of the argument `silent`.

**Value**

the inverse matrix of `x`; if `log.det==TRUE`, this inverse has an attribute which contains the logarithm of the determinant of `x`.

**Author(s)**

Adelchi Azzalini

**Examples**

```r
x <- toeplitz(rev(1:4))
x.inv <- pd.solve(x)
print(x.inv **% x)
x.inv <- pd.solve(x, log.det=TRUE)
logDet <- attr(x.inv, "log.det")
print(abs(logDet - determinant(x, logarithm=TRUE)%%modulus))
```
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