

Package ‘nvmix’

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Title Multivariate Normal Variance Mixtures (Including Student's t Distribution for Non-Integer Degrees of Freedom)

Description Functions for working with multivariate normal variance mixture distributions including evaluating their distribution functions, densities and random number generation.

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dnvmix

*Density of Multivariate Normal Variance Mixtures***Description**

Evaluating multivariate normal variance mixture densities (including Student t and normal densities).

Usage

```
dnvmix(x, qmix, loc = rep(0, d), scale = diag(d),
       factor = NULL, method = c("sobol", "ghalton", "PRNG"),
       abstol = 0.001, CI.factor = 3.3, fun.eval = c(2^6, 1e8), B = 12,
       log = FALSE, verbose = TRUE, ...)

dStudent(x, df, loc = rep(0, d), scale = diag(d), factor = NULL,
         log = FALSE, verbose = TRUE, ...)
dNorm(x, loc = rep(0, d), scale = diag(d), factor = NULL,
      log = FALSE, verbose = TRUE, ...)
```

Arguments

x	(n, d)- matrix of evaluation points.
qmix	specification of the mixing variable W ; see McNeil et al. (2015, Chapter 6). Supported are the following types of specification (see also the examples below): character: character string specifying a supported mixing distribution; currently available are "constant" (in which case $W = 1$ and thus the multivariate normal distribution with mean vector <code>loc</code> and covariance matrix <code>scale</code> results) and "inverse.gamma" (in which case W is inverse gamma distributed with shape and rate parameters <code>df/2</code> and thus the multivariate Student t distribution with <code>df</code> degrees of freedom (required to be provided via the ellipsis argument) results). list: list of length at least one, where the first component is a character string specifying the base name of a distribution whose quantile function can be accessed via the prefix "q"; an example is "exp" for which "qexp" exists. If the list is of length larger than one, the remaining elements contain additional parameters of the distribution; for "exp", for example, this can be the parameter rate. function: function interpreted as the quantile function of the mixing variable W .
df	positive degree of freedom; can also be <code>Inf</code> in which case the distribution is interpreted as the multivariate normal distribution with mean vector <code>loc</code> and covariance matrix <code>scale</code>).
loc	location vector of dimension d ; this equals the mean vector of a random vector following the specified normal variance mixture distribution if and only if the latter exists.

scale	scale matrix (a covariance matrix entering the distribution as a parameter) of dimension (d, d) ; this equals the covariance matrix of a random vector following the specified normal variance mixture distribution divided by the expectation of the mixing variable W if and only if the former exists.
factor	(d, d) lower triangular matrix such that <code>factor %*% t(factor)</code> equals <code>scale</code> ; note that for performance reasons, this property is not tested. If not provided, <code>factor</code> is internally determined via <code>t(chol())</code> .
method	character string indicating the method to be used to compute the integral. Available are: "sober": Sobol' sequence "ghalton": generalized Halton sequence "PRNG": plain Monte Carlo based on a pseudo-random number generator
abstol	non-negative numeric providing the absolute precision required. If <code>abstol = 0</code> , the algorithm will typically run until the total number of function evaluations exceeds <code>fun.eval[2]</code> . If $n > 1$ (so <code>x</code> has more than one row), the algorithm runs until the precision requirement is reached for all n density estimates.
CI.factor	multiplier of the Monte Carlo confidence interval bounds. The algorithm runs until <code>CI.factor</code> times the estimated standard error is less than <code>abstol</code> . If <code>CI.factor = 3.3</code> (the default), one can expect the actual absolute error to be less than <code>abstol</code> in 99.9% of the cases.
fun.eval	numeric (2) providing the size of the first point set to be used to estimate the probabilities (typically a power of 2) and the maximal number of function evaluations.
B	number of randomizations for obtaining an error estimate in the randomized quasi-Monte Carlo (RQMC) approach.
log	logical indicating whether the logarithmic density is to be computed.
verbose	logical indicating whether a warning is given if the required precision <code>abstol</code> has not been reached.
...	additional arguments (for example, parameters) passed to the underlying mixing distribution when <code>qmix</code> is a character string or function .

Details

Internally used is `factor`, so `scale` is not required to be provided if `factor` is given.

The default factorization used to obtain `factor` is the Cholesky decomposition via `chol()`. To this end, `scale` needs to have full rank.

The number of rows of `factor` equals the dimension d of the sample. Typically (but not necessarily), `factor` is square.

Internally, an iterative randomized Quasi-Monte Carlo (RQMC) approach is used to estimate the density. It is an iterative algorithm that evaluates the integrand at a randomized Sobol' point-set in each iteration until the pre-specified error tolerance `abstol` is reached for both the density and the log-density. The attribute `"numiter"` gives the worst case number of such iterations needed (over all rows of `x`).

Care should be taken when changing the algorithm-specific parameters, notably `method`, `fun.eval[2]` and `B`. Error estimates will not be reliable for too small `B` and the performance of the algorithm depends heavily on the (quasi-)Monte Carlo point-set used.

If the absolute error tolerance `abstol` cannot be achieved with `fun.eval[2]` function evaluations, an additional warning is thrown.

`dStudent()` and `dNorm()` are wrappers of `dnvmix(, qmix = "inverse.gamma", df = df)` and `dnvmix(, qmix = "constant")`, respectively. In these cases, `dnvmix()` uses the analytical formulas for the density of a multivariate Student t and normal distribution, respectively.

Value

`dnvmix()`, `dStudent()` and `dNorm()` return a `numeric` n -vector with the computed (log-)density values and attributes `"error"` (containing the error estimate of the RQMC estimator) and `"numiter"` (containing the number of iterations). If $n > 1$ (so `x` has more than one row), the error estimate reported is the largest error estimate among the n density estimates.

Author(s)

Erik Hintz, Marius Hofert and Christiane Lemieux.

References

McNeil, A. J., Frey, R. and Embrechts, P. (2015). *Quantitative Risk Management: Concepts, Techniques, Tools*. Princeton University Press.

See Also

`pnvmix()`, `rnvmix()`

Examples

```
### Examples for dnvmix() #####

## Generate a random correlation matrix in three dimensions
d <- 3
set.seed(271)
A <- matrix(runif(d * d), ncol = d)
P <- cov2cor(A %*% t(A))

## Evaluate a t_{3.5} density
df <- 3.5
x <- matrix(1:12/12, ncol = d) # evaluation points
dt1 <- dnvmix(x, qmix = "inverse.gamma", df = df, scale = P)
stopifnot(all.equal(dt1, c(0.013266542, 0.011967156, 0.010760575, 0.009648682),
                    tol = 1e-7, check.attributes = FALSE))

## Here is a version providing the quantile function of the mixing distribution
qW <- function(u, df) 1 / qgamma(u, shape = df/2, rate = df/2)
dt2 <- dnvmix(x, qmix = qW, df = df, scale = P)

## Compare
```

```

stopifnot(all.equal(dt1, dt2, tol = 5e-4, check.attributes = FALSE))

## Evaluate a normal density
dn <- dnmix(x, qmix = "constant", scale = P)
stopifnot(all.equal(dn, c(0.013083858, 0.011141923, 0.009389987, 0.007831596),
                    tol = 1e-7, check.attributes = FALSE))

## Case with missing data
x. <- x
x.[3,2] <- NA
x.[4,3] <- NA
dt <- dnmix(x., qmix = "inverse.gamma", df = df, scale = P)
stopifnot(is.na(dt) == rep(c(FALSE, TRUE), each = 2))

## Univariate case
x.. <- cbind(1:10/10) # (n = 10, 1)-matrix; note: vectors are taken as rows in dnmix()
dt1 <- dnmix(x.., qmix = "inverse.gamma", df = df, factor = 1)
dt2 <- dt(as.vector(x..), df = df)
stopifnot(all.equal(dt1, dt2, check.attributes = FALSE))

### Examples for dStudent() and dNorm() #####

## Evaluate a t_{3.5} density
dt <- dStudent(x, df = df, scale = P)
stopifnot(all.equal(dt, c(0.013266542, 0.011967156, 0.010760575, 0.009648682),
                    tol = 1e-7, check.attributes = FALSE))

## Evaluate a normal density
dn <- dNorm(x, scale = P)
stopifnot(all.equal(dn, c(0.013083858, 0.011141923, 0.009389987, 0.007831596),
                    tol = 1e-7, check.attributes = FALSE))

```

pnvmix

Distribution Function of Multivariate Normal Variance Mixtures

Description

Evaluating multivariate normal variance mixture distribution functions (including Student t and normal distributions).

Usage

```

pnvmix(upper, lower = matrix(-Inf, nrow = n, ncol = d), qmix, mean.sqrt.mix = NULL,
       loc = rep(0, d), scale = diag(d), standardized = FALSE,
       method = c("sobol", "ghalton", "PRNG"), precondition = TRUE,
       abstol = 1e-3, CI.factor = 3.3, fun.eval = c(2^6, 1e8),
       increment = c("doubling", "num.init"), B = 12, verbose = TRUE, ...)

pStudent(upper, lower = rep(-Inf, d), df, loc = rep(0, d), scale = diag(d),

```

```

        standardized = FALSE, method = c("sobol", "ghalton", "PRNG"),
        precondition = TRUE, abstol = 1e-3, CI.factor = 3.3, fun.eval = c(2^6, 1e8),
        B = 12, verbose = TRUE)
pNorm(upper, lower = rep(-Inf, d), loc = rep(0, d), scale = diag(d),
      standardized = FALSE, method = c("sobol", "ghalton", "PRNG"),
      precondition = TRUE, abstol = 1e-3, CI.factor = 3.3, fun.eval = c(2^6, 1e8),
      B = 12, verbose = TRUE)

```

Arguments

upper	(n, d) -matrix of upper integration limits; each row represents a d -vector of upper integration limits.
lower	(n, d) -matrix of lower integration limits (componentwise less than or equal to upper); each row represents a d -vector of lower integration limits.
qmix	specification of the mixing variable W ; see McNeil et al. (2015, Chapter 6). Supported are the following types of specification (see also the examples below): character: <code>character</code> string specifying a supported distribution; currently available are "constant" (in which case $W = 1$ and thus the multivariate normal distribution with mean vector <code>loc</code> and covariance matrix <code>scale</code> results) and "inverse.gamma" (in which case W is inverse gamma distributed with shape and rate parameters <code>df/2</code> and thus the multivariate Student t distribution with <code>df</code> degrees of freedom (required to be provided via the <code>ellipsis</code> argument) results). list: <code>list</code> of length at least one, where the first component is a <code>character</code> string specifying the base name of a distribution whose quantile function can be accessed via the prefix "q"; an example is "exp" for which "qexp" exists. If the list is of length larger than one, the remaining elements contain additional parameters of the distribution; for "exp", for example, this can be the parameter <code>rate</code> . function: <code>function</code> interpreted as the quantile function of the mixing variable W .
mean.sqrt.mix	expectation of the square root \sqrt{W} of the mixing variable W . If NULL, it will be estimated via QMC; this is only needed for determining the reordering of the integration bounds, so a rather crude approximation is fine.
df	positive degree of freedom; can also be <code>Inf</code> in which case the distribution is interpreted as the multivariate normal distribution with mean vector <code>loc</code> and covariance matrix <code>scale</code>).
loc	location vector of dimension d ; this equals the mean vector of a random vector following the specified normal variance mixture distribution if and only if the latter exists.
scale	scale matrix (a covariance matrix entering the distribution as a parameter) of dimension (d, d) ; this equals the covariance matrix of a random vector following the specified normal variance mixture distribution divided by the expectation of the mixing variable W if and only if the former exists.
standardized	logical indicating whether <code>scale</code> is assumed to be a correlation matrix.

method	character string indicating the method to be used to compute the integral. Available are: "sobol": Sobol' sequence "ghalton": generalized Halton sequence "PRNG": plain Monte Carlo based on a pseudo-random number generator
precond	logical indicating whether preconditioning is applied, that is, reordering of the integration variables. If TRUE, integration limits as well as scale are internally re-ordered in a way such that the overall variance of the integrand is usually smaller than with the original ordering; this usually leads smaller run-times.
abstol	non-negative numeric providing the absolute precision required; if <code>abstol = NULL</code> , the algorithm will run until the total number of function evaluations <code>fun.eval[2]</code> is reached.
CI.factor	multiplier of the Monte Carlo confidence interval bounds. The algorithm runs until <code>CI.factor</code> times the estimated standard error is less than <code>abstol</code> . If <code>CI.factor = 3.3</code> (the default), one can expect the actual absolute error to be less than <code>abstol</code> in 99.9% of the cases.
fun.eval	numeric (2) providing the size of the first point set to be used to estimate the probabilities (typically a power of 2) and the maximal number of function evaluations.
increment	character string indicating how the sample size should be increased in each iteration. Available are: "doubling": next iteration has as many sample points as all the previous iterations combined. "num.init": all iterations use an additional <code>fun.eval[1]</code> -many points.
B	number of randomizations for obtaining an error estimate in the randomized quasi-Monte Carlo (RQMC) approach.
verbose	logical indicating whether a warning is given if the required precision <code>abstol</code> has not been reached; can also be an integer in which case 0 is FALSE, 1 is TRUE and 2 stands for producing a more verbose warning (for each set of provided integration bounds).
...	additional arguments (for example, parameters) passed to the underlying mixing distribution when <code>qmix</code> is a character string or function .

Details

One should highlight that evaluating normal variance mixtures is a non-trivial tasks which, at the time of development of **pnvmix**, was not available in R before, not even the special case of a multivariate Student t distribution for non-integer degrees of freedom, which frequently appears in applications in finance, insurance and risk management after estimating such distributions.

Note that the procedures call underlying C code. Currently, dimensions $d \geq 16510$ are not supported for the default method `sobol`.

Internally, an iterative randomized Quasi-Monte Carlo (RQMC) approach is used to estimate the probabilities. It is an iterative algorithm that evaluates the integrand at a point-set with size as specified by `increment`) in each iteration until the pre-specified error tolerance `abstol` is reached. The attribute `"num.iter"` gives the number of such iterations needed.

Care should be taken when changing the algorithm-specific parameters, notably `method`, `precond`, `fun.eval[2]` and `B`. Error estimates will not be reliable for too small `B` and the performance of the algorithm depends heavily on the (quasi-)Monte Carlo point-set used.

If the absolute error tolerance `abstol` cannot be achieved with `fun.eval[2]` function evaluations, an additional warning is thrown.

`pStudent()` and `pNorm()` are wrappers of `pnmix(, qmix = "inverse.gamma", df = df)` and `pnmix(, qmix = "constant")`, respectively. In the univariate case, the functions `pt()` and `pnorm()` are used.

Value

`pnmix()`, `pStudent()` and `pNorm()` return a `numeric` n -vector with the computed probabilities and corresponding attributes `"error"` (error estimates of the RQMC estimator) and `"numiter"` (number of iterations).

Author(s)

Erik Hintz, Marius Hofert and Christiane Lemieux

References

McNeil, A. J., Frey, R. and Embrechts, P. (2015). *Quantitative Risk Management: Concepts, Techniques, Tools*. Princeton University Press.

Genz, A. and Bretz, F. (1999). Numerical computation of multivariate t-probabilities with application to power calculation of multiple contrasts. *Journal of Statistical Computation and Simulation* 63(4), 103–117.

Genz, A. and Bretz, F. (2002). Comparison of methods for the computation of multivariate t probabilities. *Journal of Computational and Graphical Statistics* 11(4), 950–971.

See Also

`dnvmix()`, `rnvmix()`

Examples

```
### Examples for pnmix() #####

## Generate a random correlation matrix in d dimensions
d <- 3
set.seed(157)
A <- matrix(runif(d * d), ncol = d)
P <- cov2cor(A %*% t(A))

## Evaluate a t_{1/2} distribution function
a <- -3 * runif(d) * sqrt(d) # random lower limit
b <- 3 * runif(d) * sqrt(d) # random upper limit
df <- 0.5 # note that this is *non-integer*
set.seed(1)
pt1 <- pnmix(b, lower = a, qmix = "inverse.gamma", df = df, scale = P)
```

```

## Here is a version providing the quantile function of the mixing distribution
qW <- function(u, df) 1 / qgamma(u, shape = df/2, rate = df/2)
mean.sqrt.mix <- sqrt(df) * gamma(df/2) / (sqrt(2) * gamma((df+1) / 2))
set.seed(1)
pt2 <- pnmix(b, lower = a, qmix = qW, mean.sqrt.mix = mean.sqrt.mix, df = df,
            scale = P)

## Compare
stopifnot(all.equal(pt1, pt2, tol = 7e-4, check.attributes = FALSE))

## mean.sqrt.mix will be approximated by QMC internally if not provided,
## so the results will differ slightly.
set.seed(1)
pt3 <- pnmix(b, lower = a, qmix = qW, df = df, scale = P)
stopifnot(all.equal(pt3, pt1, tol = 7e-4, check.attributes = FALSE))

## Case with missing data and a matrix of lower and upper bounds
a. <- matrix(rep(a, each = 4), ncol = d)
b. <- matrix(rep(b, each = 4), ncol = d)
a.[2,1] <- NA
b.[3,2] <- NA
pt <- pnmix(b., lower = a., qmix = "inverse.gamma", df = df, scale = P)
stopifnot(is.na(pt) == c(FALSE, TRUE, TRUE, FALSE))

## Case where upper = (Inf,...,Inf) and lower = (-Inf,...,-Inf)
stopifnot(all.equal(pnmix(upper = rep(Inf, d), qmix = "constant"), 1,
                    check.attributes = FALSE))

### Examples for pStudent() and pNorm() #####

## Evaluate a t_{3.5} distribution function
set.seed(271)
pt <- pStudent(b, lower = a, df = 3.5, scale = P)
stopifnot(all.equal(pt, 0.6180, tol = 5e-5, check.attributes = FALSE))

## Evaluate a normal distribution function
set.seed(271)
pn <- pNorm(b, lower = a, scale = P)
stopifnot(all.equal(pn, 0.7001, tol = 1e-4, check.attributes = FALSE))

## pStudent deals correctly with df = Inf:
set.seed(1)
p.St.dfInf <- pStudent(b, df = Inf, scale = P)
set.seed(1)
p.Norm <- pNorm(b, scale = P)
stopifnot(all.equal(p.St.dfInf, p.Norm, check.attributes = FALSE))

```

Description

Generate vectors of random variates from multivariate normal variance mixtures (including Student t and normal distributions).

Usage

```
rnmix(n, rmix = NULL, qmix = NULL, loc = rep(0, d), scale = diag(2),
      factor = NULL, method = c("PRNG", "sobol", "ghalton"),
      skip = 0, ...)
```

```
rStudent(n, df, loc = rep(0, d), scale = diag(2), factor = NULL,
         method = c("PRNG", "sobol", "ghalton"), skip = 0)
```

```
rNorm(n, loc = rep(0, d), scale = diag(2), factor = NULL,
      method = c("PRNG", "sobol", "ghalton"), skip = 0)
```

Arguments

- n** sample size n (positive integer).
- rmix** specification of the mixing variable W , see McNeil et al. (2015, Chapter 6), via a random number generator. This argument is ignored for `method = "sobol"` and `method = "ghalton"`. Supported are the following types of specification (see also the examples below):
- character:** `character` string specifying a supported distribution; currently available are `"constant"` (in which case $W = 1$ and thus a sample from the multivariate normal distribution with mean vector `loc` and covariance matrix `scale` results) and `"inverse.gamma"` (in which case W is inverse gamma distributed with shape and rate parameters `df/2` and thus the multivariate Student t distribution with `df` degrees of freedom (required to be provided via the `ellipsis` argument) results).
 - list:** `list` of length at least one, where the first component is a `character` string specifying the base name of a distribution which can be sampled via prefix `"r"`; an example is `"exp"` for which `"rexp"` exists for sampling. If the list is of length larger than one, the remaining elements contain additional parameters of the distribution; for `"exp"`, for example, this can be the parameter `rate`.
 - function:** `function` interpreted as a random number generator of the mixing variable W ; additional arguments (such as parameters) can be passed via the `ellipsis` argument.
 - numeric:** `numeric` vector of length `n` providing a random sample of the mixing variable W .
- qmix** specification of the mixing variable W via a quantile function; see McNeil et al. (2015, Chapter 6). This argument is required for `method = "sobol"` and `method = "ghalton"`. Supported are the following types of specification (see also the examples below):
- character:** `character` string specifying a supported distribution; currently available are `"constant"` (in which case $W = 1$ and thus a sample from the multivariate normal distribution with mean vector `loc` and covariance

matrix scale results) and "inverse.gamma" (in which case W is inverse gamma distributed with shape and rate parameters $df/2$ and thus the multivariate Student t distribution with df degrees of freedom (required to be provided via the ellipsis argument) results).

list: `list` of length at least one, where the first component is a `character` string specifying the base name of a distribution which can be sampled via prefix "q"; an example is "exp" for which "qexp" exists for sampling. If the list is of length larger than one, the remaining elements contain additional parameters of the distribution; for "exp", for example, this can be the parameter `rate`.

function: `function` interpreted as the quantile function of the mixing variable W ; internally, sampling is then done with the inversion method by applying the provided function to $U(0,1)$ random variates.

<code>df</code>	positive degree of freedom; can also be <code>Inf</code> in which case the distribution is interpreted as the multivariate normal distribution with mean vector <code>loc</code> and covariance matrix <code>scale</code>).
<code>loc</code>	location vector of dimension d ; this equals the mean vector of a random vector following the specified normal variance mixture distribution if and only if the latter exists.
<code>scale</code>	scale matrix (a covariance matrix entering the distribution as a parameter) of dimension (d, d) (defaults to $d = 2$); this equals the covariance matrix of a random vector following the specified normal variance mixture distribution divided by the expectation of the mixing variable W if and only if the former exists. Note that <code>scale</code> must be positive definite; sampling from singular normal variance mixtures can be achieved by providing <code>factor</code> .
<code>factor</code>	(d, k) -matrix such that <code>factor %*% t(factor)</code> equals <code>scale</code> ; the non-square case $k \neq d$ can be used to sample from singular normal variance mixtures. Note that this notation coincides with McNeil et al. (2015, Chapter 6). If not provided, <code>factor</code> is internally determined via <code>chol()</code> (and multiplied from the right to an (n, k) -matrix of independent standard normals to obtain a sample from a multivariate normal with zero mean vector and covariance matrix <code>scale</code>).
<code>method</code>	<p><code>character</code> string indicating the method to be used to obtain the sample. Available are:</p> <ul style="list-style-type: none"> "PRNG": pseudo-random numbers "sobol": Sobol' sequence "ghalton": generalized Halton sequence <p>If <code>method = "PRNG"</code>, either <code>qmix</code> or <code>rmix</code> can be provided. If both are provided, <code>rmix</code> is used and <code>qmix</code> ignored. For the other two methods, sampling is done via inversion, hence <code>qmix</code> has to be provided and <code>rmix</code> is ignored.</p>
<code>skip</code>	<code>integer</code> specifying the number of points to be skipped when <code>method = "sobol"</code> , see also example below.
<code>...</code>	additional arguments (for example, parameters) passed to the underlying mixing distribution when <code>rmix</code> or <code>qmix</code> is a <code>character</code> string or <code>function</code> .

Details

Internally used is `factor`, so `scale` is not required to be provided if `factor` is given.

The default factorization used to obtain `factor` is the Cholesky decomposition via `chol()`. To this end, `scale` needs to have full rank.

Sampling from a singular normal variance mixture distribution can be achieved by providing `scale`.

The number of rows of `factor` equals the dimension d of the sample. Typically (but not necessarily), `factor` is square.

`rStudent()` and `rNorm()` are wrappers of `rnmix(, qmix = "inverse.gamma", df = df)` and `rnmix(, qmix = "constant", df = df)`, respectively.

Value

`rnmix()` returns an (n, d) -matrix containing n samples of the specified (via `mix`) d -dimensional multivariate normal variance mixture with location vector `loc` and scale matrix `scale` (a covariance matrix).

`rStudent()` returns samples from the d -dimensional multivariate Student t distribution with location vector `loc` and scale matrix `scale`.

`rNorm()` returns samples from the d -dimensional multivariate normal distribution with mean vector `loc` and covariance matrix `scale`.

Author(s)

Erik Hintz, Marius Hofert and Christiane Lemieux

References

McNeil, A. J., Frey, R. and Embrechts, P. (2015). *Quantitative Risk Management: Concepts, Techniques, Tools*. Princeton University Press.

See Also

`dnvmix()`, `pnvmix()`

Examples

```
### Examples for rnmix() #####

## Generate a random correlation matrix in d dimensions
d <- 3
set.seed(157)
A <- matrix(runif(d * d), ncol = d)
P <- cov2cor(A %*% t(A))

## Draw random variates and compare
df <- 3.5
n <- 1000
set.seed(271)
X <- rnmix(n, rmix = "inverse.gamma", df = df, scale = P) # providing scale
```

```

set.seed(271)
X. <- rnmix(n, rmix = "inverse.gamma", df = df, factor = t(chol(P))) # providing the factor
stopifnot(all.equal(X, X.))

## Checking df = Inf
set.seed(271)
X <- rnmix(n, rmix = "constant", scale = P) # normal
set.seed(271)
X. <- rnmix(n, rmix = "inverse.gamma", scale = P, df = Inf) # t_infinity
stopifnot(all.equal(X, X.))

## Univariate case (dimension = number of rows of 'factor' = 1 here)
set.seed(271)
X.1d <- rnmix(n, rmix = "inverse.gamma", df = df, factor = 1/2)
set.seed(271)
X.1d. <- rnmix(n, rmix = "inverse.gamma", df = df, factor = 1)/2 # manual scaling
stopifnot(all.equal(X.1d, X.1d.))

## Checking different ways of providing 'mix'
## 1) By providing a character string (and corresponding ellipsis arguments)
set.seed(271)
X.mix1 <- rnmix(n, rmix = "inverse.gamma", df = df, scale = P)
## 2) By providing a list; the first element has to be an existing distribution
## with random number generator available with prefix "r"
rinverse.gamma <- function(n, df) 1 / rgamma(n, shape = df/2, rate = df/2)
set.seed(271)
X.mix2 <- rnmix(n, rmix = list("inverse.gamma", df = df), scale = P)
## 3) The same without extra arguments (need the extra list() here to
## distinguish from Case 1))
rinverseGamma <- function(n) 1 / rgamma(n, shape = df/2, rate = df/2)
set.seed(271)
X.mix3 <- rnmix(n, rmix = list("inverseGamma"), scale = P)
## 4) By providing a quantile function
## Note:  $P(1/Y \leq x) = P(Y \geq 1/x) = 1 - F_Y(1/x) = y \Leftrightarrow x = 1/F_Y^{-(1-y)}$ 
set.seed(271)
X.mix4 <- rnmix(n, qmix = function(p) 1/qgamma(1-p, shape = df/2, rate = df/2),
               scale = P)
## 5) By providing random variates
set.seed(271) # if seed is set here, results are comparable to the above methods
W <- rinverse.gamma(n, df = df)
X.mix5 <- rnmix(n, rmix = W, scale = P)
## Compare (note that X.mix4 is not 'all equal' with X.mix1 or the other samples)
## since rgamma() != qgamma(runif()) (or qgamma(1-runif()))
stopifnot(all.equal(X.mix2, X.mix1),
          all.equal(X.mix3, X.mix1),
          all.equal(X.mix5, X.mix1))

## For a singular normal variance mixture:
## Need to provide 'factor'
A <- matrix( c(1, 0, 0, 1, 0, 1), ncol = 2, byrow = TRUE)
stopifnot(all.equal(dim(rnmix(n, rmix = "constant", factor = A)), c(n, 3)))
stopifnot(all.equal(dim(rnmix(n, rmix = "constant", factor = t(A))), c(n, 2)))

```

```

## Using 'skip'. Need to reset the seed everytime to get the same shifts in "sobol".
## Note that when using method = "sobol", we have to provide 'qmix' instead of 'rmix'.
set.seed(271)
X.skip0 <- rnmix(n, qmix = "inverse.gamma", df = df, scale = P, method = "sobol")
set.seed(271)
X.skip1 <- rnmix(n, qmix = "inverse.gamma", df = df, scale = P, method = "sobol",
                skip = n)
set.seed(271)
X.wo.skip <- rnmix(2*n, qmix = "inverse.gamma", df = df, scale = P, method = "sobol")
X.skip <- rbind(X.skip0, X.skip1)
all.equal(X.wo.skip, X.skip)

### Examples for rStudent() and rNorm() #####

## Draw  $N(0, P)$  random variates by providing scale or factor and compare
n <- 1000
set.seed(271)
X.n <- rNorm(n, scale = P) # providing scale
set.seed(271)
X.n. <- rNorm(n, factor = t(chol(P))) # providing the factor
stopifnot(all.equal(X.n, X.n.))

## Univariate case (dimension = number of rows of 'factor' = 1 here)
set.seed(271)
X.n.1d <- rNorm(n, factor = 1/2)
set.seed(271)
X.n.1d. <- rNorm(n, factor = 1)/2 # manual scaling
stopifnot(all.equal(X.n.1d, X.n.1d.))

## Draw  $t_{3.5}$  random variates by providing scale or factor and compare
df <- 3.5
n <- 1000
set.seed(271)
X.t <- rStudent(n, df = df, scale = P) # providing scale
set.seed(271)
X.t. <- rStudent(n, df = df, factor = t(chol(P))) # providing the factor
stopifnot(all.equal(X.t, X.t.))

## Univariate case (dimension = number of rows of 'factor' = 1 here)
set.seed(271)
X.t.1d <- rStudent(n, df = df, factor = 1/2)
set.seed(271)
X.t.1d. <- rStudent(n, df = df, factor = 1)/2 # manual scaling
stopifnot(all.equal(X.t.1d, X.t.1d.))

## Check  $df = \text{Inf}$ 
set.seed(271)
X.t <- rStudent(n, df = Inf, scale = P)
set.seed(271)
X.n <- rNorm(n, scale = P)
stopifnot(all.equal(X.t, X.n))

```

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