Package: posterior (via r-universe)

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Title Tools for Working with Posterior Distributions

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Description Provides useful tools for both users and developers of packages for fitting Bayesian models or working with output from Bayesian models. The primary goals of the package are to: (a) Efficiently convert between many different useful formats of draws (samples) from posterior or prior distributions. (b) Provide consistent methods for operations commonly performed on draws, for example, subsetting, binding, or mutating draws. (c) Provide various summaries of draws in convenient formats. (d) Provide lightweight implementations of state of the art posterior inference diagnostics. References: Vehtari et al. (2021) <doi:10.1214/20-BA1221>.

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Contents

`posterior-package` *Tools for working with posterior (and prior) distributions*

Description

The **posterior** package is intended to provide useful tools for both users and developers of packages for fitting Bayesian models or working with output from Bayesian models. The primary goals of the package are to:

- Efficiently convert between many different useful formats of draws (samples) from posterior or prior distributions.
- Provide consistent methods for operations commonly performed on draws, for example, subsetting, binding, or mutating draws.
- Provide various summaries of draws in convenient formats.
- Provide lightweight implementations of state of the art posterior inference diagnostics.

Package options

The following options are used to format and print `draws` objects, as in `print.draws_array()`, `print.draws_df()`, `print.draws_list()`, `print.draws_matrix()`, and `print.draws_rvars()`:

- `posterior.max_draws`: Maximum number of draws to print.
- `posterior.max_iterations`: Maximum number of iterations to print.
- `posterior.max_chains`: Maximum number of chains to print.
- `posterior.max_variables`: Maximum number of variables to print.

The following options are used for formatting the output of `summarize_draws`:

- `posterior.num_args`: Arguments passed to num() for pretty printing of summaries.

The following options are used to format and print `rvar` objects, as in `print.rvar()` and `print.draws_rvars()`:

- `posterior.rvar_summary`: What style of summary to display: `"mean_sd"` displays `mean ± sd`, `"median_mad"` displays `median ± mad`.
- `posterior.digits`: How many significant digits are displayed. This defaults to a smaller value (2) than `getOption("digits")` because `rvar`s print two numbers (point summary and uncertainty) next to each other.

The following option is used to construct new `rvar` objects, as in `rfun()` and `rdo()`:

- `posterior.rvar_ndraws`: The number of draws used to construct new random variables when this number cannot be determined from existing arguments (e.g., other `rvar`s passed to a function).

The following options are used to control warning messages:

- `posterior.warn_on_merge_chains`: (logical) Some operations will trigger an automatic merging of chains, for example, because chains do not match between two objects involved in a binary operation. Whether this causes a warning can be controlled by this option.

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See Also

Useful links:

- https://mc-stan.org/posterior/
- https://discourse.mc-stan.org/
- Report bugs at https://github.com/stan-dev/posterior/issues

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as_rvar *Coerce to a random variable*

Description

Convert `x` to an `rvar` object.

Usage

```
as_rvar(x, dim = NULL, dimnames = NULL, nchains = NULL)

as_rvar_numeric(x, dim = NULL, dimnames = NULL, nchains = NULL)

as_rvar_integer(x, dim = NULL, dimnames = NULL, nchains = NULL)

as_rvar_logical(x, dim = NULL, dimnames = NULL, nchains = NULL)
```

Arguments

x	(multiple options) An object that can be converted to an rvar, such as a vector, array, or an rvar itself.
dim	(integer vector) One or more integers giving the maximal indices in each dimension to override the dimensions of the rvar to be created (see dim()). If NULL (the default), dim is determined by the input. **NOTE:** This argument controls the dimensions of the rvar, not the underlying array, so you cannot change the number of draws using this argument.
dimnames	(list) Character vectors giving the names in each dimension to override the names of the dimensions of the rvar to be created (see dimnames()). If NULL (the default), this is determined by the input. **NOTE:** This argument controls the names of the dimensions of the rvar, not the underlying array.
nchains	(positive integer) The number of chains. The default is 1.

Details

For objects that are already rvars, returns them (with modified dimensions if dim is not NULL).

For numeric or logical vectors or arrays, returns an rvar with a single draw and the same dimensions as x. This is in contrast to the rvar() constructor, which treats the first dimension of x as the draws dimension. As a result, as_rvar() is useful for creating constants.

While as_rvar() attempts to pick the most suitable subtype of rvar based on the type of x (possibly returning an rvar_factor or rvar_ordered), as_rvar_numeric(), as_rvar_integer(), and as_rvar_logical() always coerce the draws of the output rvar to be numeric, integer, or logical (respectively), and always return a base rvar, never a subtype.

Value

An object of class "rvar" (or one of its subtypes) representing a random variable.

See Also

rvar() to construct rvars directly. See rdo(), rfun(), and rvar_rng() for higher-level interfaces for creating rvars.

Examples

```
# You can use as_rvar() to create "constant" rvars (having only one draw):
x <- as_rvar(1)
x

# Such constants can be of arbitrary shape:
as_rvar(1:4)
as_rvar(matrix(1:10, nrow = 5))
as_rvar(array(1:12, dim = c(2, 3, 2)))

# as_rvar_numeric() coerces subtypes of rvar to the base rvar type
y <- as_rvar_factor(c("a", "b", "c"))
y
```

```
as_rvar_numeric(y)
```

`as_rvar_factor` *Coerce to a factor random variable*

Description

Convert `x` to an `rvar_factor` or `rvar_ordered` object.

Usage

```
as_rvar_factor(x, dim = NULL, dimnames = NULL, nchains = NULL, ...)

as_rvar_ordered(x, dim = NULL, dimnames = NULL, nchains = NULL, ...)
```

Arguments

`x`	(multiple options) An object that can be converted to an `rvar`, such as a vector, array, or an `rvar` itself.
`dim`	(integer vector) One or more integers giving the maximal indices in each dimension to override the dimensions of the `rvar` to be created (see `dim()`). If `NULL` (the default), `dim` is determined by the input. **NOTE:** This argument controls the dimensions of the `rvar`, not the underlying array, so you cannot change the number of draws using this argument.
`dimnames`	(list) Character vectors giving the names in each dimension to override the names of the dimensions of the `rvar` to be created (see `dimnames()`). If `NULL` (the default), this is determined by the input. **NOTE:** This argument controls the names of the dimensions of the `rvar`, not the underlying array.
`nchains`	(positive integer) The number of chains. The default is `1`.
`...`	Arguments passed on to `base::factor`
	`levels` an optional vector of the unique values (as character strings) that `x` might have taken. The default is the unique set of values taken by `as.character(x)`, sorted into increasing order *of* `x`. Note that this set can be specified as smaller than `sort(unique(x))`.
	`labels` *either* an optional character vector of labels for the levels (in the same order as `levels` after removing those in `exclude`), *or* a character string of length 1. Duplicated values in `labels` can be used to map different values of `x` to the same factor level.
	`exclude` a vector of values to be excluded when forming the set of levels. This may be factor with the same level set as `x` or should be a `character`.
	`ordered` logical flag to determine if the levels should be regarded as ordered (in the order given).
	`nmax` an upper bound on the number of levels; see 'Details'.

Details

For objects that are already rvars, returns them (with modified dimensions if dim is not NULL), possibly adding levels using the unique values of the draws of the rvar (if the object is not already factor-like).

For numeric, logical, factor, or character vectors or arrays, returns an rvar_factor or rvar_ordered with a single draw and the same dimensions as x. This is in contrast to the rvar_factor() and rvar_ordered() constructors, which treats the first dimension of x as the draws dimension. As a result, as_rvar_factor() and as_rvar_ordered() are useful for creating constants.

Value

An object of class "rvar_factor" or "rvar_ordered" representing a random variable.

See Also

rvar(), rvar_factor(), and rvar_ordered() to construct rvars directly. See rdo(), rfun(), and rvar_rng() for higher-level interfaces for creating rvars.

Examples

```
# You can use as_rvar_factor() to create "constant" rvars (having only one draw):
x <- as_rvar_factor("a")
x

# Such constants can be of arbitrary shape:
as_rvar_factor(letters[1:4])
as_rvar_ordered(matrix(letters[1:10], nrow = 5))
as_rvar_factor(array(letters[1:12], dim = c(2, 3, 2)))
```

---

bind_draws *Bind* draws *objects together*

Description

Bind multiple draws objects together to form a single draws object.

Usage

```
bind_draws(x, ...)

## S3 method for class 'draws_matrix'
bind_draws(x, ..., along = "variable")

## S3 method for class 'draws_array'
bind_draws(x, ..., along = "variable")
```

```
## S3 method for class 'draws_df'
bind_draws(x, ..., along = "variable")

## S3 method for class 'draws_list'
bind_draws(x, ..., along = "variable")

## S3 method for class 'draws_rvars'
bind_draws(x, ..., along = "variable")
```

Arguments

x	(draws) A draws object. The draws format of x will define the format of the returned draws object.
...	(draws) Additional draws objects to bind to x.
along	(string) The dimension along which draws objects should be bound together. Possible values are "variable" (the default), "chain", "iteration", and "draw". Not all options are supported for all input formats.

Value

A draws object of the same class as x.

Examples

```
x1 <- draws_matrix(alpha = rnorm(5), beta = rnorm(5))
x2 <- draws_matrix(alpha = rnorm(5), beta = rnorm(5))
ndraws(x1)
ndraws(x2)
x3 <- bind_draws(x1, x2, along = "draw")
ndraws(x3)

x4 <- draws_matrix(theta = rexp(5))
x5 <- bind_draws(x1, x4, along = "variable")
variables(x5)
```

---

chol.rvar *Cholesky decomposition of random matrix*

Description

Cholesky decomposition of an rvar containing a matrix.

Usage

```
## S3 method for class 'rvar'
chol(x, ...)
```

Arguments

x	(rvar) A 2-dimensional rvar.
...	Additional parameters passed on to chol.tensor()

Value

An rvar containing the upper triangular factor of the Cholesky decomposition, i.e., the matrix R such that $R'R = x$.

---

diag,rvar-method *Matrix diagonals (including for random variables)*

Description

Extract the diagonal of a matrix or construct a matrix, including random matrices (2-dimensional rvars). Makes base::diag() generic.

Usage

```
## S4 method for signature 'rvar'
diag(x = 1, nrow, ncol, names = TRUE)
```

Arguments

x	(numeric,rvar) a matrix, vector, 1D array, missing, or a 1- or 2-dimensional rvar.
nrow, ncol	optional dimensions for the result when x is not a matrix.
names	(when x is a matrix) logical indicating if the resulting vector, the diagonal of x, should inherit names from dimnames(x) if available.

Details

Makes base::diag() into a generic function. See that function's documentation for usage with numerics and for usage of diag<-, which is also supported by rvar.

Value

For rvars, has two modes:

1. x is a matrix-like rvar: it returns the diagonal as a vector-like rvar
2. x is a vector-like rvar: it returns a matrix-like rvar with x as the diagonal and zero for off-diagonal entries.

See Also

base::diag()

Examples

```
# Sigma is a 3x3 covariance matrix
Sigma <- as_draws_rvars(example_draws("multi_normal"))$Sigma
Sigma

diag(Sigma)

diag(Sigma) <- 1:3
Sigma

diag(as_rvar(1:3))
```

`diagnostics` *List of available convergence diagnostics*

Description

A list of available diagnostics and links to their individual help pages.

Details

Function	Description
`ess_basic()`	Basic version of effective sample size
`ess_bulk()`	Bulk effective sample size
`ess_tail()`	Tail effective sample size
`ess_mean()`	Effective sample sizes for the mean
`ess_median()`	Effective sample sizes for the median
`ess_quantile()`	Effective sample sizes for quantiles
`ess_sd()`	Effective sample sizes for the standard deviation
`mcse_mean()`	Monte Carlo standard error for the mean
`mcse_median()`	Monte Carlo standard error for the median
`mcse_quantile()`	Monte Carlo standard error for quantiles
`mcse_sd()`	Monte Carlo standard error for the standard deviation
`pareto_khat()`	Pareto khat diagnostic for tail(s)
`pareto_diags()`	Additional diagnostics related to Pareto khat
`rhat_basic()`	Basic version of Rhat
`rhat()`	Improved, rank-based version of Rhat
`rhat_nested()`	Rhat for use with many short chains
`rstar()`	R* diagnostic

Value

See individual functions for a description of return types.

`dissent` *Dissention*

Description

Dissention, for measuring dispersion in draws from ordinal distributions.

Usage

```
dissent(x)

## Default S3 method:
dissent(x)

## S3 method for class 'rvar'
dissent(x)
```

Arguments

`x`	(multiple options) A vector to be interpreted as draws from an ordinal distribution, such as: • A factor • A numeric (should be integer or integer-like) • An rvar, rvar_factor, or rvar_ordered

Details

Calculates Tastle and Wierman's (2007) *dissention* measure:

$$-\sum_{i=1}^{n} p_i \log_2 \left(1 - \frac{|x_i - \mathrm{E}(x)|}{\max(x) - \min(x)}\right)$$

This ranges from 0 (all probability in one category) through 0.5 (uniform) to 1 (bimodal: all probability split equally between the first and last category).

Value

If `x` is a factor or numeric, returns a length-1 numeric vector with a value between 0 and 1 (inclusive) giving the dissention of `x`.

If `x` is an rvar, returns an array of the same shape as `x`, where each cell is the dissention of the draws in the corresponding cell of `x`.

References

William J. Tastle, Mark J. Wierman (2007). Consensus and dissention: A measure of ordinal dispersion. *International Journal of Approximate Reasoning*. 45(3), 531–545. doi:10.1016/j.ijar.2006.06.024.

Examples

```
set.seed(1234)

levels <- c("lowest", "low", "neutral", "high", "highest")

# a bimodal distribution: high dissention
x <- ordered(
  sample(levels, 4000, replace = TRUE, prob = c(0.45, 0.04, 0.02, 0.04, 0.45)),
  levels = levels
)
dissent(x)

# a unimodal distribution: low dissention
y <- ordered(
  sample(levels, 4000, replace = TRUE, prob = c(0.95, 0.02, 0.015, 0.01, 0.005)),
  levels = levels
)
dissent(y)

# both together, as an rvar
xy <- c(rvar(x), rvar(y))
xy
dissent(xy)
```

---

draws *Transform to* draws *objects*

Description

Try to transform an R object to a format supported by the **posterior** package.

Usage

```
as_draws(x, ...)

is_draws(x)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).

Details

The class "draws" is the parent class of all supported formats, which also have their own subclasses of the form "draws_{format}" (e.g. "draws_array").

Value

If possible, a draws object in the closest supported format to x. The formats are linked to in the **See Also** section below.

See Also

Other formats: draws_array(), draws_df(), draws_list(), draws_matrix(), draws_rvars()

Examples

```
# create some random draws
x <- matrix(rnorm(30), nrow = 10)
colnames(x) <- c("a", "b", "c")
str(x)

# transform to a draws object
y <- as_draws(x)
str(y)

# remove the draws classes from the object
class(y) <- class(y)[-(1:2)]
str(y)
```

---

draws-index *Index* draws *objects*

Description

Index iterations, chains, and draws of draws objects.

Usage

```
iteration_ids(x)

chain_ids(x)

draw_ids(x)

niterations(x)

nchains(x)

ndraws(x)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.

Details

The methods `iteration_ids()`, `chain_ids()`, and `draw_ids()` return vectors of all iterations, chains, and draws, respectively. In contrast, the methods `niterations()`, `nchains()`, and `ndraws()` return the number of variables, iterations, chains, and draws, respectively.

Value

For `iteration_ids()`, `chain_ids()`, and `draw_ids()`, an integer vector.

For `niterations()`, `nchains()`, and `ndraws()`, a scalar integer.

See Also

`variables`, `rename_variables`

Examples

```
x <- example_draws()

iteration_ids(x)
niterations(x)

chain_ids(x)
nchains(x)

draw_ids(x)
ndraws(x)
```

---

draws_array *The* draws_array *format*

Description

The `as_draws_array()` methods convert objects to the `draws_array` format. The `draws_array()` function creates an object of the `draws_array` format based on a set of numeric vectors. See **Details**.

Usage

```
as_draws_array(x, ...)

## Default S3 method:
as_draws_array(x, ...)

## S3 method for class 'draws_array'
as_draws_array(x, ...)

## S3 method for class 'draws_matrix'
```

```
as_draws_array(x, ...)

## S3 method for class 'draws_df'
as_draws_array(x, ...)

## S3 method for class 'draws_list'
as_draws_array(x, ...)

## S3 method for class 'draws_rvars'
as_draws_array(x, ...)

## S3 method for class 'mcmc'
as_draws_array(x, ...)

## S3 method for class 'mcmc.list'
as_draws_array(x, ...)

draws_array(..., .nchains = 1)

is_draws_array(x)
```

Arguments

x	An object to convert to a draws_array object.
...	For as_draws_array(): Arguments passed to individual methods (if applicable). For draws_array(): Named arguments containing numeric vectors each defining a separate variable.
.nchains	(positive integer) The number of chains. The default is 1.

Details

Objects of class "draws_array" are 3-D arrays with dimensions "iteration", "chain", and "variable". See **Examples**.

Value

A draws_array object, which has classes c("draws_array", "draws", "array").

See Also

Other formats: draws, draws_df(), draws_list(), draws_matrix(), draws_rvars()

Examples

```
x1 <- as_draws_array(example_draws())
class(x1)
print(x1)
str(x1)
```

```
x2 <- draws_array(a = rnorm(10), b = rnorm(10), c = 1)
class(x2)
print(x2)
str(x2)
```

---

draws_df *The* draws_df *format*

Description

The as_draws_df() methods convert objects to the draws_df format. The draws_df() function creates an object of the draws_df format based on a set of numeric vectors. See **Details**.

Usage

```
as_draws_df(x, ...)

## Default S3 method:
as_draws_df(x, ...)

## S3 method for class 'data.frame'
as_draws_df(x, ...)

## S3 method for class 'draws_df'
as_draws_df(x, ...)

## S3 method for class 'draws_matrix'
as_draws_df(x, ...)

## S3 method for class 'draws_array'
as_draws_df(x, ...)

## S3 method for class 'draws_list'
as_draws_df(x, ...)

## S3 method for class 'draws_rvars'
as_draws_df(x, ...)

## S3 method for class 'mcmc'
as_draws_df(x, ...)

## S3 method for class 'mcmc.list'
as_draws_df(x, ...)

draws_df(..., .nchains = 1)

is_draws_df(x)
```

Arguments

x	An object to convert to a draws_df object.
...	For as_draws_df(): Arguments passed to individual methods (if applicable). For draws_df(): Named arguments containing numeric vectors each defining a separate variable.
.nchains	(positive integer) The number of chains. The default is 1.

Details

Objects of class "draws_df" are tibble data frames. They have one column per variable as well as additional metadata columns ".iteration", ".chain", and ".draw". The difference between the ".iteration" and ".draw" columns is that the former is relative to the MCMC chain while the latter ignores the chain information and has all unique values. See **Examples**.

If a data.frame-like object is supplied to as_draws_df that contains columns named ".iteration" or ".chain", they will be treated as iteration and chain indices, respectively. See **Examples**.

Value

A draws_df object, which has classes c("draws_df", "draws", class(tibble::tibble())).

See Also

Other formats: draws, draws_array(), draws_list(), draws_matrix(), draws_rvars()

Examples

```
x1 <- as_draws_df(example_draws())
class(x1)
print(x1)
str(x1)

x2 <- draws_df(a = rnorm(10), b = rnorm(10), c = 1)
class(x2)
print(x2)
str(x2)

# the difference between iteration and draw is clearer when contrasting
# the head and tail of the data frame
print(head(x1), reserved = TRUE, max_variables = 2)
print(tail(x1), reserved = TRUE, max_variables = 2)

# manually supply chain information
xnew <- data.frame(mu = rnorm(10), .chain = rep(1:2, each = 5))
xnew <- as_draws_df(xnew)
print(xnew)
```

draws_list *The* draws_list *format*

Description

The as_draws_list() methods convert objects to the draws_list format. The draws_list() function creates an object of the draws_list format based on a set of numeric vectors. See **Details**.

Usage

```
as_draws_list(x, ...)

## Default S3 method:
as_draws_list(x, ...)

## S3 method for class 'draws_list'
as_draws_list(x, ...)

## S3 method for class 'draws_matrix'
as_draws_list(x, ...)

## S3 method for class 'draws_array'
as_draws_list(x, ...)

## S3 method for class 'draws_df'
as_draws_list(x, ...)

## S3 method for class 'draws_rvars'
as_draws_list(x, ...)

## S3 method for class 'mcmc'
as_draws_list(x, ...)

## S3 method for class 'mcmc.list'
as_draws_list(x, ...)

draws_list(..., .nchains = 1)

is_draws_list(x)
```

Arguments

x	An object to convert to a draws_list object.
...	For as_draws_list(): Arguments passed to individual methods (if applicable). For draws_list(): Named arguments containing numeric vectors each defining a separate variable.
.nchains	(positive integer) The number of chains. The default is 1.

Details

Objects of class "draws_list" are lists with one element per MCMC chain. Each of these elements is itself a named list of numeric vectors with one vector per variable. The length of each vector is equal to the number of saved iterations per chain. See **Examples**.

Value

A draws_list object, which has classes c("draws_list", "draws", "list").

See Also

Other formats: draws, draws_array(), draws_df(), draws_matrix(), draws_rvars()

Examples

```
x1 <- as_draws_list(example_draws())
class(x1)
print(x1)
str(x1)

x2 <- draws_list(a = rnorm(10), b = rnorm(10), c = 1)
class(x2)
print(x2)
str(x2)
```

---

draws_matrix *The* draws_matrix *format*

Description

The as_draws_matrix() methods convert objects to the draws_matrix format. The draws_matrix() function creates an object of the draws_matrix format based on a set of numeric vectors. See **Details**.

Usage

```
as_draws_matrix(x, ...)

## Default S3 method:
as_draws_matrix(x, ...)

## S3 method for class 'draws_matrix'
as_draws_matrix(x, ...)

## S3 method for class 'draws_array'
as_draws_matrix(x, ...)

## S3 method for class 'draws_df'
```

```
as_draws_matrix(x, ...)

## S3 method for class 'draws_list'
as_draws_matrix(x, ...)

## S3 method for class 'draws_rvars'
as_draws_matrix(x, ...)

## S3 method for class 'mcmc'
as_draws_matrix(x, ...)

## S3 method for class 'mcmc.list'
as_draws_matrix(x, ...)

draws_matrix(..., .nchains = 1)

is_draws_matrix(x)
```

Arguments

x	An object to convert to a draws_matrix object.
...	For as_draws_matrix(): Arguments passed to individual methods (if applicable). For draws_matrix(): Named arguments containing numeric vectors each defining a separate variable.
.nchains	(positive integer) The number of chains. The default is 1.

Details

Objects of class "draws_matrix" are matrices (2-D arrays) with dimensions "draw" and "variable". See **Examples**.

Value

A draws_matrix object, which has classes c("draws_matrix", "draws", "matrix").

See Also

Other formats: draws, draws_array(), draws_df(), draws_list(), draws_rvars()

Examples

```
x1 <- as_draws_matrix(example_draws())
class(x1)
print(x1)
str(x1)

x2 <- draws_matrix(a = rnorm(10), b = rnorm(10), c = 1)
class(x2)
print(x2)
str(x2)
```

draws_of *Get/set array of draws underlying a random variable*

Description

Gets/sets the array-representation that backs an rvar. Should be used rarely.

Usage

```
draws_of(x, with_chains = FALSE)

draws_of(x, with_chains = FALSE) <- value
```

Arguments

x	(rvar) An rvar object.
with_chains	(logical) Should the array of draws include a dimension for chains? If FALSE (the default), chains are not included and the array has dimension c(ndraws(x), dim(x)). If TRUE, chains are included and the array has dimension c(niterations(x), nchains(x), dim(x)).
value	(array) An array of values to use as the backing array of x.

Details

While rvars implement fast versions of basic math operations (including matrix multiplication), sometimes you may need to bypass the rvar abstraction to do what you need to do more efficiently. draws_of() allows you to get / set the underlying array of draws in order to do that.

rvars represent draws internally using arrays of arbitrary dimension, which is returned by draws_of(x) and can be set using draws_of(x) <- value. The **first** dimension of these arrays is the index of the draws. If with_chains = TRUE, then the dimensions of the returned array are modified so that the first dimension is the index of the iterations and the second dimension is the index of the chains.

Value

If with_chains = FALSE, an array with dimensions c(ndraws(x), dim(x)).

If with_chains = TRUE, an array with dimensions c(niterations(x), nchains(x), dim(x)).

Examples

```
x <- rvar(1:10, nchains = 2)
x

# draws_of() without arguments will return the array of draws without
# chain information (first dimension is draw)
draws_of(x)

# draws_of() with with_chains = TRUE will reshape the returned array to
```

```
# include chain information in the second dimension
draws_of(x, with_chains = TRUE)

# you can also set draws using draws_of(). When with_chains = FALSE the
# existing chain information will be retained ...
draws_of(x) <- 2:11
x

# when with_chains = TRUE the chain information will be set by the
# second dimension of the assigned array
draws_of(x, with_chains = TRUE) <- array(2:11, dim = c(2,5))
x
```

---

draws_rvars *The* draws_rvars *format*

Description

The as_draws_rvars() methods convert objects to the draws_rvars format. The draws_rvars() function creates an object of the draws_rvars format based on a set of numeric vectors. See **Details**.

Usage

```
as_draws_rvars(x, ...)

## Default S3 method:
as_draws_rvars(x, ...)

## S3 method for class 'draws_rvars'
as_draws_rvars(x, ...)

## S3 method for class 'list'
as_draws_rvars(x, ...)

## S3 method for class 'draws_matrix'
as_draws_rvars(x, ...)

## S3 method for class 'draws_array'
as_draws_rvars(x, ...)

## S3 method for class 'draws_df'
as_draws_rvars(x, ...)

## S3 method for class 'draws_list'
as_draws_rvars(x, ...)
```

```
## S3 method for class 'mcmc'
as_draws_rvars(x, ...)

## S3 method for class 'mcmc.list'
as_draws_rvars(x, ...)

draws_rvars(..., .nchains = 1)

is_draws_rvars(x)
```

Arguments

x	An object to convert to a draws_rvars object.
...	For as_draws_rvars(): Arguments passed to individual methods (if applicable). For draws_rvars(): Named arguments containing numeric vectors each defining a separate variable.
.nchains	(positive integer) The number of chains. The default is 1.

Details

Objects of class "draws_rvars" are lists of rvar objects. See **Examples**.

Value

A draws_rvars object, which has classes c("draws_rvars", "draws", "list").

See Also

Other formats: draws, draws_array(), draws_df(), draws_list(), draws_matrix()

Examples

```
x1 <- as_draws_rvars(example_draws())
class(x1)
print(x1)
str(x1)

x2 <- draws_rvars(a = rnorm(10), b = rnorm(10), c = 1)
class(x2)
print(x2)
str(x2)
```

draws_summary *Summaries of* draws *objects*

Description

The summarise_draws() (and summarize_draws()) methods provide a quick way to get a table of summary statistics and diagnostics. These methods will convert an object to a draws object if it isn't already. For convenience, a summary() method for draws and rvar objects are also provided as an alias for summarise_draws() if the input object is a draws or rvar object.

Usage

```
summarise_draws(.x, ...)

summarize_draws(.x, ...)

## S3 method for class 'draws'
summarise_draws(
  .x,
  ...,
  .args = list(),
  .num_args = getOption("posterior.num_args", list()),
  .cores = 1
)

## S3 method for class 'draws'
summary(object, ...)

## S3 method for class 'rvar'
summarise_draws(.x, ...)

## S3 method for class 'rvar'
summary(object, ...)

default_summary_measures()

default_convergence_measures()

default_mcse_measures()
```

Arguments

.x, object	(draws) A draws object or one coercible to a draws object.
...	Name-value pairs of summary or diagnostic functions. The provided names will be used as the names of the columns in the result *unless* the function returns a named vector, in which case the latter names are used. The functions can be specified in any format supported by as_function(). See **Examples**.

.args	(named list) Optional arguments passed to the summary functions.
.num_args	(named list) Optional arguments passed to num() for pretty printing of summaries. Can be controlled globally via the posterior.num_args option.
.cores	(positive integer) The number of cores to use for computing summaries for different variables in parallel. Coerced to integer if possible, otherwise errors. The default is .cores = 1, in which case no parallelization is implemented. By default, a socket cluster is used on Windows and forks otherwise.

Details

The default summary functions used are the ones specified by default_summary_measures() and default_convergence_measures():

default_summary_measures()

- mean()
- median()
- sd()
- mad()
- quantile2()

default_convergence_measures()

- rhat()
- ess_bulk()
- ess_tail()

The var() function should not be used to compute variances due to its inconsistent behavior with matrices. Instead, please use distributional::variance().

Value

The summarise_draws() methods return a tibble data frame. The first column ("variable") contains the variable names and the remaining columns contain summary statistics and diagnostics.

The functions default_summary_measures(), default_convergence_measures(), and default_mcse_measures() return character vectors of names of the default measures.

See Also

diagnostics for a list of available diagnostics and links to their individual help pages.

Examples

```
x <- example_draws("eight_schools")
class(x)
str(x)

summarise_draws(x)
summarise_draws(x, "mean", "median")
```

```
summarise_draws(x, mean, mcse = mcse_mean)
summarise_draws(x, ~quantile(.x, probs = c(0.4, 0.6)))

# using default_*_meaures()
summarise_draws(x, default_summary_measures())
summarise_draws(x, default_convergence_measures())
summarise_draws(x, default_mcse_measures())

# compute variance of variables
summarise_draws(x, var = distributional::variance)

# illustrate use of '.args'
ws <- rexp(ndraws(x))
summarise_draws(x, weighted.mean, .args = list(w = ws))

# adjust how numerical summaries are printed
summarise_draws(x, .num_args = list(sigfig = 2, notation = "dec"))
```

`drop,rvar-method` *Drop redundant dimensions*

Description

Delete the dimensions of an `rvar` which are of size one. See `base::drop()`

Usage

```
## S4 method for signature 'rvar'
drop(x)
```

Arguments

`x`	(rvar) an `rvar`.

Value

An `rvar` with the same length as `x`, but where any entry equal to 1 in `dim(x)` has been removed. The exception is if `dim(x) == 1`, in which case `dim(drop(x)) == 1` as well (this is because `rvar`s, unlike `numeric`s, never have `NULL` dimensions).

Examples

```
# Sigma is a 3x3 covariance matrix
Sigma <- as_draws_rvars(example_draws("multi_normal"))$Sigma
Sigma

Sigma[1, ]

drop(Sigma[1, ])
```

```
# equivalently ...
Sigma[1, drop = TRUE]
```

`entropy` *Normalized entropy*

Description

Normalized entropy, for measuring dispersion in draws from categorical distributions.

Usage

```
entropy(x)

## Default S3 method:
entropy(x)

## S3 method for class 'rvar'
entropy(x)
```

Arguments

`x`	(multiple options) A vector to be interpreted as draws from a categorical distribution, such as: • A factor • A numeric (should be integer or integer-like) • An rvar, rvar_factor, or rvar_ordered

Details

Calculates the normalized Shannon entropy of the draws in `x`. This value is the entropy of `x` divided by the maximum entropy of a distribution with `n` categories, where `n` is `length(unique(x))` for numeric vectors and `length(levels(x))` for factors:

$$-\frac{\sum_{i=1}^{n} p_i \log(p_i)}{\log(n)}$$

This scales the output to be between 0 (all probability in one category) and 1 (uniform). This form of normalized entropy is referred to as H_{REL} in Wilcox (1967).

Value

If `x` is a factor or numeric, returns a length-1 numeric vector with a value between 0 and 1 (inclusive) giving the normalized Shannon entropy of `x`.

If `x` is an rvar, returns an array of the same shape as `x`, where each cell is the normalized Shannon entropy of the draws in the corresponding cell of `x`.

References

Allen R. Wilcox (1967). *Indices of Qualitative Variation* (No. ORNL-TM-1919). Oak Ridge National Lab., Tenn.

Examples

```
set.seed(1234)

levels <- c("a", "b", "c", "d", "e")

# a uniform distribution: high normalized entropy
x <- factor(
  sample(levels, 4000, replace = TRUE, prob = c(0.2, 0.2, 0.2, 0.2, 0.2)),
  levels = levels
)
entropy(x)

# a unimodal distribution: low normalized entropy
y <- factor(
  sample(levels, 4000, replace = TRUE, prob = c(0.95, 0.02, 0.015, 0.01, 0.005)),
  levels = levels
)
entropy(y)

# both together, as an rvar
xy <- c(rvar(x), rvar(y))
xy
entropy(xy)
```

---

ess_basic *Basic version of the effective sample size*

Description

Compute the basic effective sample size (ESS) estimate for a single variable as described in Gelman et al. (2013) with some changes according to Vehtari et al. (2021). For practical applications, we strongly recommend the improved ESS convergence diagnostics implemented in `ess_bulk()` and `ess_tail()`. See Vehtari (2021) for an in-depth comparison of different effective sample size estimators.

Usage

```
ess_basic(x, ...)

## Default S3 method:
ess_basic(x, split = TRUE, ...)

## S3 method for class 'rvar'
ess_basic(x, split = TRUE, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See extract_variable_matrix(). • An rvar.
...	Arguments passed to individual methods (if applicable).
split	(logical) Should the estimate be computed on split chains? The default is TRUE.

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, NA, NaN, Inf, or -Inf, the returned output will be (numeric) NA. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) NA as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an rvar, returns an array of the same dimensions as the rvar, where each element is equal to the value that would be returned by passing the draws array for that element of the rvar to this function.

References

Andrew Gelman, John B. Carlin, Hal S. Stern, David B. Dunson, Aki Vehtari and Donald B. Rubin (2013). *Bayesian Data Analysis, Third Edition*. Chapman and Hall/CRC.

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

Aki Vehtari (2021). Comparison of MCMC effective sample size estimators. Retrieved from https://avehtari.github.io/rhat_ess/ess_comparison.html

See Also

Other diagnostics: ess_bulk(), ess_quantile(), ess_sd(), ess_tail(), mcse_mean(), mcse_quantile(), mcse_sd(), pareto_diags(), pareto_khat(), rhat(), rhat_basic(), rhat_nested(), rstar()

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
ess_basic(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
ess_basic(d$Sigma)
```

`ess_bulk` *Bulk effective sample size (bulk-ESS)*

Description

Compute a bulk effective sample size estimate (bulk-ESS) for a single variable. Bulk-ESS is useful as a diagnostic for the sampling efficiency in the bulk of the posterior. It is defined as the effective sample size for rank normalized values using split chains. For the tail effective sample size see `ess_tail()`. See Vehtari (2021) for an in-depth comparison of different effective sample size estimators.

Usage

```
ess_bulk(x, ...)

## Default S3 method:
ess_bulk(x, ...)

## S3 method for class 'rvar'
ess_bulk(x, ...)
```

Arguments

`x`	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
`...`	Arguments passed to individual methods (if applicable).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

Aki Vehtari (2021). Comparison of MCMC effective sample size estimators. Retrieved from https://avehtari.github.io/rhat_ess/ess_comparison.html

See Also

Other diagnostics: ess_basic(), ess_quantile(), ess_sd(), ess_tail(), mcse_mean(), mcse_quantile(), mcse_sd(), pareto_diags(), pareto_khat(), rhat(), rhat_basic(), rhat_nested(), rstar()

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
ess_bulk(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
ess_bulk(d$Sigma)
```

---

ess_mean *Effective sample size for the mean*

Description

Compute an effective sample size estimate for a mean (expectation) estimate of a single variable.

Usage

```
ess_mean(x, ...)

## Default S3 method:
ess_mean(x, ...)

## S3 method for class 'rvar'
ess_mean(x, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See extract_variable_matrix(). • An rvar.
...	Arguments passed to individual methods (if applicable).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Andrew Gelman, John B. Carlin, Hal S. Stern, David B. Dunson, Aki Vehtari and Donald B. Rubin (2013). *Bayesian Data Analysis, Third Edition*. Chapman and Hall/CRC.

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
ess_mean(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
ess_mean(d$Sigma)
```

---

ess_quantile *Effective sample sizes for quantiles*

Description

Compute effective sample size estimates for quantile estimates of a single variable.

Usage

```
ess_quantile(x, probs = c(0.05, 0.95), ...)

## Default S3 method:
ess_quantile(x, probs = c(0.05, 0.95), names = TRUE, ...)

## S3 method for class 'rvar'
ess_quantile(x, probs = c(0.05, 0.95), names = TRUE, ...)

ess_median(x, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
probs	(numeric vector) Probabilities in `[0, 1]`.
...	Arguments passed to individual methods (if applicable).
names	(logical) Should the result have a `names` attribute? The default is `TRUE`, but use `FALSE` for improved speed if there are many values in `probs`.

Value

If the input is an array, returns a numeric vector with one element per quantile. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be a vector of (numeric) `NA` values. Also, if all draws of a variable are the same (constant), the returned output will be a vector of (numeric) `NA` values as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar` and `length(probs) == 1`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function. If `length(probs) > 1`, the first dimension of the result indexes the input probabilities; i.e. the result has dimension `c(length(probs), dim(x))`.

References

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

See Also

Other diagnostics: `ess_basic()`, `ess_bulk()`, `ess_sd()`, `ess_tail()`, `mcse_mean()`, `mcse_quantile()`, `mcse_sd()`, `pareto_diags()`, `pareto_khat()`, `rhat()`, `rhat_basic()`, `rhat_nested()`, `rstar()`

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
ess_quantile(mu, probs = c(0.1, 0.9))

d <- as_draws_rvars(example_draws("multi_normal"))
ess_quantile(d$mu, probs = c(0.1, 0.9))
```

`ess_sd` *Effective sample size for the standard deviation*

Description

Compute an effective sample size estimate for the standard deviation (SD) estimate of a single variable. This is defined as the effective sample size estimate for the absolute deviation from mean.

Usage

```
ess_sd(x, ...)

## Default S3 method:
ess_sd(x, ...)

## S3 method for class 'rvar'
ess_sd(x, ...)
```

Arguments

`x`	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
`...`	Arguments passed to individual methods (if applicable).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

See Also

Other diagnostics: ess_basic(), ess_bulk(), ess_quantile(), ess_tail(), mcse_mean(), mcse_quantile(), mcse_sd(), pareto_diags(), pareto_khat(), rhat(), rhat_basic(), rhat_nested(), rstar()

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
ess_sd(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
ess_sd(d$Sigma)
```

---

ess_tail *Tail effective sample size (tail-ESS)*

Description

Compute a tail effective sample size estimate (tail-ESS) for a single variable. Tail-ESS is useful as a diagnostic for the sampling efficiency in the tails of the posterior. It is defined as the minimum of the effective sample sizes for 5% and 95% quantiles. For the bulk effective sample size see ess_bulk(). See Vehtari (2021) for an in-depth comparison of different effective sample size estimators.

Usage

```
ess_tail(x, ...)

## Default S3 method:
ess_tail(x, ...)

## S3 method for class 'rvar'
ess_tail(x, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See extract_variable_matrix(). • An rvar.
...	Arguments passed to individual methods (if applicable).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

Aki Vehtari (2021). Comparison of MCMC effective sample size estimators. Retrieved from https://avehtari.github.io/rhat_ess/ess_comparison.html

See Also

Other diagnostics: `ess_basic()`, `ess_bulk()`, `ess_quantile()`, `ess_sd()`, `mcse_mean()`, `mcse_quantile()`, `mcse_sd()`, `pareto_diags()`, `pareto_khat()`, `rhat()`, `rhat_basic()`, `rhat_nested()`, `rstar()`

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
ess_tail(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
ess_tail(d$Sigma)
```

---

example_draws *Example* draws *objects*

Description

Objects for use in examples, vignettes, and tests.

Usage

```
example_draws(example = "eight_schools")
```

Arguments

example	(string) The name of the example draws object. See **Details** for available options.

Details

The following example draws objects are available.

eight_schools: A draws_array object with 100 iterations from each of 4 Markov chains obtained by fitting the eight schools model described in Gelman et al. (2013) with Stan. The variables are:

- mu: Overall mean of the eight schools
- tau: Standard deviation between schools
- theta: Individual means of each of the eight schools

multi_normal: A draws_array object with 100 iterations from each of the 4 Markov chains obtained by fitting a 3-dimensional multivariate normal model to 100 simulated observations. The variables are:

- mu: Mean parameter vector of length 3
- Sigma: Covariance matrix of dimension 3 x 3

Value

A draws object.

Note

These objects are only intended to be used in demonstrations and tests. They contain fewer iterations and chains than recommended for performing actual inference.

References

Andrew Gelman, John B. Carlin, Hal S. Stern, David B. Dunson, Aki Vehtari and Donald B. Rubin (2013). Bayesian Data Analysis, Third Edition. Chapman and Hall/CRC.

Examples

```
draws_eight_schools <- example_draws("eight_schools")
summarise_draws(draws_eight_schools)

draws_multi_normal <- example_draws("multi_normal")
summarise_draws(draws_multi_normal)
```

`extract_list_of_variable_arrays` *Extract arrays of multiple variables*

Description

Extract arrays of draws for multiple variables, returning them as a named list of arrays. Each array has the same structure as returned by `extract_variable_array`.

Usage

```
extract_list_of_variable_arrays(x, variables = NULL, ...)

## Default S3 method:
extract_list_of_variable_arrays(x, variables = NULL, ...)

## S3 method for class 'draws'
extract_list_of_variable_arrays(x, variables = NULL, ...)
```

Arguments

`x`	(draws) A `draws` object or another R object for which the method is defined.
`variables`	A character vector of variable names to extract, or NULL to extract all variables. To extract all dimensions from variables with indices (e.g. `"x[1]"`), provide the base variable names (e.g. `"x"`).
`...`	Arguments passed to individual methods (if applicable).

Value

A named list of arrays, where each array has dimension `niterations(x)` x `nchains(x)` x any remaining dimensions determined by the indices of the variable `x` (if `with_chains = TRUE`) or dimension `niterations(x) * nchains(x)` x any remaining dimensions (if `with_chains = FALSE`).

See Also

Other variable extraction methods: `extract_variable()`, `extract_variable_array()`, `extract_variable_matrix()`

Examples

```
x <- example_draws(example = "multi_normal")

# Extract multiple variables at once
vars <- extract_list_of_variable_arrays(x, c("mu", "Sigma"))
str(vars)

# Extract all variables (uses base variable names)
all_vars <- extract_list_of_variable_arrays(x)
str(all_vars)
```

```
# Extract specific indexed variables
vars2 <- extract_list_of_variable_arrays(x, c("mu[1]", "mu[2]"))
str(vars2)
```

`extract_variable` *Extract draws of a single variable*

Description

Extract a vector of draws of a single variable.

Usage

```
extract_variable(x, variable, ...)

## Default S3 method:
extract_variable(x, variable, ...)

## S3 method for class 'draws'
extract_variable(x, variable, ...)

## S3 method for class 'draws_df'
extract_variable(x, variable, ...)

## S3 method for class 'draws_list'
extract_variable(x, variable, ...)

## S3 method for class 'draws_rvars'
extract_variable(x, variable, ...)
```

Arguments

`x`	(draws) A `draws` object or another R object for which the method is defined.
`variable`	(string) The name of the variable to extract. The name must correspond to a scalar variable or must include indices for array variables (e.g. `"x[1]"`, `"y[1,2]"`) that result in a scalar variable. To extract all dimensions from variables with indices, use `extract_variable_array()`.
`...`	Arguments passed to individual methods (if applicable).

Value

A vector of length equal to the number of draws.

See Also

Other variable extraction methods: `extract_list_of_variable_arrays()`, `extract_variable_array()`, `extract_variable_matrix()`

Examples

```
x <- example_draws()
mu <- extract_variable(x, variable = "mu")
str(mu)
```

`extract_variable_array`

Extract array of a single (possibly indexed) variable

Description

Extract an array of draws of a single variable, including any dimensions of variables with indices.

Usage

```
extract_variable_array(x, variable, with_chains = TRUE, ...)

## Default S3 method:
extract_variable_array(x, variable, with_chains = TRUE, ...)

## S3 method for class 'draws'
extract_variable_array(x, variable, with_chains = TRUE, ...)
```

Arguments

`x`	(draws) A `draws` object or another R object for which the method is defined.
`variable`	(string) The name of the variable to extract. To extract all dimensions from variables with indices (e.g. `"x[1]"`), provide the base variable name (e.g. `"x"`).
`with_chains`	(logical) Should the array of draws include a dimension for chains? If `TRUE` (the default), chains are included and the array has dimension `c(niterations(x), nchains(x), ...)`. If `FALSE`, chains are not included and the array has dimension `c(niterations(x) * nchains(x), ...)`.
`...`	Arguments passed to individual methods (if applicable).

Value

An `array` with dimension `niterations(x)` x `nchains(x)` x any remaining dimensions determined by the indices of the variable `x` (if `with_chains` = `TRUE`) or dimension `niterations(x)` * `nchains(x)` x any remaining dimensions (if `with_chains` = `FALSE`).

See Also

Other variable extraction methods: `extract_list_of_variable_arrays()`, `extract_variable()`, `extract_variable_matrix()`

Examples

```
x <- example_draws(example = "multi_normal")

mu <- extract_variable_array(x, variable = "mu")
str(mu)

# With chains collapsed
mu_no_chains <- extract_variable_array(x, variable = "mu", with_chains = FALSE)
str(mu_no_chains)

mu1 <- extract_variable_array(x, variable = "mu[1]")
str(mu1)

Sigma <- extract_variable_array(x, variable = "Sigma")
str(Sigma)
```

---

extract_variable_matrix

Extract matrix of a single variable

Description

Extract an iterations x chains matrix of draws of a single variable. This is primarily used for convergence diagnostic functions such as `rhat()`.

Usage

```
extract_variable_matrix(x, variable, ...)

## Default S3 method:
extract_variable_matrix(x, variable, ...)

## S3 method for class 'draws'
extract_variable_matrix(x, variable, ...)

## S3 method for class 'draws_df'
extract_variable_matrix(x, variable, ...)

## S3 method for class 'draws_list'
extract_variable_matrix(x, variable, ...)

## S3 method for class 'draws_rvars'
extract_variable_matrix(x, variable, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
variable	(string) The name of the variable to extract. The name must correspond to a scalar variable or must include indices for array variables (e.g. "x[1]", "y[1,2]") that result in a scalar variable. To extract all dimensions from variables with indices, use extract_variable_array().
...	Arguments passed to individual methods (if applicable).

Value

A matrix with dimension iterations x chains.

See Also

Other variable extraction methods: extract_list_of_variable_arrays(), extract_variable(), extract_variable_array()

Examples

```
x <- example_draws()
mu <- extract_variable_matrix(x, variable = "mu")
dim(mu)
rhat(mu)
```

---

for_each_draw *Loop over draws*

Description

Executes an expression once for every draw in a draws object. Used primarily for its side effects and returns the input x invisibly.

Usage

```
for_each_draw(x, expr)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
expr	(expression) A bare expression that can contain references to variables in x by name. This expression will be executed once per draw of x, where references to variables in x resolve to the value of that variable in that draw. The expression supports quasiquotation.

Details

If x is not in the draws_rvars format, it is first converted to that format. This allows the variables in x to include their dimensions (i.e, to act as R vectors and arrays) when being referred to in expr.

Within expr, use .draw to refer to the draw index, which will be a value between 1 and ndraws(x). expr is executed in the calling environment of for_each_draw(), so it can use variables in that environment (however, due to the use of data masking, to modify variables in that environment, one must use <<-.)

Value

As for_each_draw() is used primarily for its side effects (the expression executed for each draw of x), it returns the input x invisibly.

Examples

```
eight_schools <- as_draws_rvars(example_draws())


# 1. A simple example --- looping over draws and printing each draw
# NOTE: You probably don't want to do this in practice! This example is
# just intended to show what for_each_draw() is doing. If you just want to
# print the draws of an rvar, it is probably better to use draws_of()
for_each_draw(eight_schools, {
  print(mu)
})


# 2. A more complex example --- building a parallel coordinates plot
# First, construct the plot bounds
plot(1, type = "n",
  xlim = c(1, length(eight_schools$theta)),
  ylim = range(range(eight_schools$theta)),
  xlab = "school", ylab = "theta"
)

# Then, use for_each_draw() to make a parallel coordinates plot of all draws
# of eight_schools$theta. Use resample_draws(eight_schools, n = ...)
# in place of eight_schools if a smaller sample is desired for the plot.
for_each_draw(eight_schools, {
  lines(seq_along(theta), theta, col = rgb(1, 0, 0, 0.05))
})

# Finally, add means and 90% intervals
lines(seq_along(eight_schools$theta), mean(eight_schools$theta))
with(summarise_draws(eight_schools$theta),
  segments(seq_along(eight_schools$theta), y0 = q5, y1 = q95)
)
```

is_rvar *Is* x *a random variable?*

Description

Test if x is an rvar.

Usage

```
is_rvar(x)
```

Arguments

x	(any object) An object to test.

Value

TRUE if x is an rvar, FALSE otherwise.

See Also

as_rvar() to convert objects to rvars.

is_rvar_factor *Is* x *a factor random variable?*

Description

Test if x is an rvar_factor or rvar_ordered.

Usage

```
is_rvar_factor(x)

is_rvar_ordered(x)
```

Arguments

x	(any object) An object to test.

Value

TRUE if x is an rvar_factor or rvar_ordered, FALSE otherwise.

See Also

as_rvar_factor() and as_rvar_ordered() to convert objects to rvar_factors and rvar_ordereds.

match *Value Matching*

Description

Generic version of base::match(). For base vectors, returns a vector of the positions of (first) matches of its first argument in its second. For rvars, returns an rvar of the matches.

Usage

```
match(x, table, ...)

## Default S3 method:
match(x, ...)

## S3 method for class 'rvar'
match(x, ...)

x %in% table
```

Arguments

x	(multiple options) the values to be matched. Can be: • A base vector: see base::match() • An rvar
table	(vector) the values to be matched against.
...	Arguments passed on to base::match nomatch the value to be returned in the case when no match is found. Note that it is coerced to integer. incomparables a vector of values that cannot be matched. Any value in x matching a value in this vector is assigned the nomatch value. For historical reasons, FALSE is equivalent to NULL.

Details

For more information on how match behaves with base vectors, see base::match().

When x is an rvar, the draws of x are matched against table using base::match(), and the result is returned as an rvar.

The implementation of %in% here is identical to base::%in%, except it uses the generic version of match() so that non-base vectors (such as rvars) are supported.

Value

When x is a base vector, a vector of the same length as x.

When x is an rvar, an rvar the same shape as x.

Examples

```
x <- rvar(c("a","b","b","c","d"))
x %in% c("b","d")

# for additional examples, see base::match()
```

`mcse_mean` *Monte Carlo standard error for the mean*

Description

Compute the Monte Carlo standard error for the mean (expectation) of a single variable.

Usage

```
mcse_mean(x, ...)

## Default S3 method:
mcse_mean(x, ...)

## S3 method for class 'rvar'
mcse_mean(x, ...)
```

Arguments

`x`	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
`...`	Arguments passed to individual methods (if applicable).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Andrew Gelman, John B. Carlin, Hal S. Stern, David B. Dunson, Aki Vehtari and Donald B. Rubin (2013). *Bayesian Data Analysis, Third Edition*. Chapman and Hall/CRC.

See Also

Other diagnostics: ess_basic(), ess_bulk(), ess_quantile(), ess_sd(), ess_tail(), mcse_quantile(), mcse_sd(), pareto_diags(), pareto_khat(), rhat(), rhat_basic(), rhat_nested(), rstar()

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
mcse_mean(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
mcse_mean(d$Sigma)
```

---

mcse_quantile *Monte Carlo standard error for quantiles*

Description

Compute Monte Carlo standard errors for quantile estimates of a single variable.

Usage

```
mcse_quantile(x, probs = c(0.05, 0.95), ...)

## Default S3 method:
mcse_quantile(x, probs = c(0.05, 0.95), names = TRUE, ...)

## S3 method for class 'rvar'
mcse_quantile(x, probs = c(0.05, 0.95), names = TRUE, ...)

mcse_median(x, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See extract_variable_matrix(). • An rvar.
probs	(numeric vector) Probabilities in [0, 1].
...	Arguments passed to individual methods (if applicable).
names	(logical) Should the result have a names attribute? The default is TRUE, but use FALSE for improved speed if there are many values in probs.

Value

If the input is an array, returns a numeric vector with one element per quantile. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be a vector of (numeric) `NA` values. Also, if all draws of a variable are the same (constant), the returned output will be a vector of (numeric) `NA` values as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar` and `length(probs) == 1`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function. If `length(probs) > 1`, the first dimension of the result indexes the input probabilities; i.e. the result has dimension `c(length(probs), dim(x))`.

References

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

See Also

Other diagnostics: `ess_basic()`, `ess_bulk()`, `ess_quantile()`, `ess_sd()`, `ess_tail()`, `mcse_mean()`, `mcse_sd()`, `pareto_diags()`, `pareto_khat()`, `rhat()`, `rhat_basic()`, `rhat_nested()`, `rstar()`

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
mcse_quantile(mu, probs = c(0.1, 0.9))

d <- as_draws_rvars(example_draws("multi_normal"))
mcse_quantile(d$mu)
```

`mcse_sd` *Monte Carlo standard error for the standard deviation*

Description

Compute the Monte Carlo standard error for the standard deviation (SD) of a single variable without assuming normality using moments of moments and first order Taylor series approximation (Kenney and Keeping, 1951, p. 141).

Usage

```
mcse_sd(x, ...)

## Default S3 method:
mcse_sd(x, ...)

## S3 method for class 'rvar'
mcse_sd(x, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
...	Arguments passed to individual methods (if applicable).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

J. F. Kenney & E. S. Keeping (1951). *Mathematics of Statistics, Vol. II.*

See Also

Other diagnostics: `ess_basic()`, `ess_bulk()`, `ess_quantile()`, `ess_sd()`, `ess_tail()`, `mcse_mean()`, `mcse_quantile()`, `pareto_diags()`, `pareto_khat()`, `rhat()`, `rhat_basic()`, `rhat_nested()`, `rstar()`

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
mcse_sd(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
```

```
mcse_sd(d$Sigma)
```

`merge_chains` *Merge chains of* `draws` *objects*

Description

Merge chains of `draws` objects into a single chain. Some operations will trigger an automatic merging of chains, for example, because chains do not match between two objects involved in a binary operation. By default, no warning will be issued when this happens but you can activate one via `options(posterior.warn_on_merge_chains = TRUE)`.

Usage

```
merge_chains(x, ...)

## S3 method for class 'draws_matrix'
merge_chains(x, ...)

## S3 method for class 'draws_array'
merge_chains(x, ...)

## S3 method for class 'draws_df'
merge_chains(x, ...)

## S3 method for class 'draws_list'
merge_chains(x, ...)

## S3 method for class 'rvar'
merge_chains(x, ...)

## S3 method for class 'draws_rvars'
merge_chains(x, ...)
```

Arguments

`x`	(draws) A `draws` object or another R object for which the method is defined.
`...`	Arguments passed to individual methods (if applicable).

Value

A `draws` object of the same class as `x`.

Examples

```
x <- example_draws()

# draws_array with 4 chains, 100 iters each
str(x)

# draws_array with 1 chain of 400 iterations
str(merge_chains(x))
```

`modal_category` *Modal category*

Description

Modal category of a vector.

Usage

```
modal_category(x)

## Default S3 method:
modal_category(x)

## S3 method for class 'rvar'
modal_category(x)
```

Arguments

`x`	(multiple options) A vector to be interpreted as draws from a categorical distribution, such as: • A factor • A numeric (should be integer or integer-like) • An rvar, rvar_factor, or rvar_ordered

Details

Finds the modal category (i.e., most frequent value) in `x`. In the case of ties, returns the first tie.

Value

If `x` is a factor or numeric, returns a length-1 vector containing the modal value.

If `x` is an rvar, returns an array of the same shape as `x`, where each cell is the modal value of the draws in the corresponding cell of `x`.

Examples

```
x <- factor(c("a","b","b","c","d"))
modal_category(x)

# in the case of ties, the first tie is returned
y <- factor(c("a","c","c","d","d"))
modal_category(y)

# both together, as an rvar
xy <- c(rvar(x), rvar(y))
xy
modal_category(xy)
```

`mutate_variables` *Mutate variables in* `draws` *objects*

Description

Mutate variables in a `draws` object.

Usage

```
mutate_variables(.x, ...)

## S3 method for class 'draws_matrix'
mutate_variables(.x, ...)

## S3 method for class 'draws_array'
mutate_variables(.x, ...)

## S3 method for class 'draws_df'
mutate_variables(.x, ...)

## S3 method for class 'draws_list'
mutate_variables(.x, ...)

## S3 method for class 'draws_rvars'
mutate_variables(.x, ...)
```

Arguments

`.x`	(draws) A `draws` object.
`...`	Name-value pairs of expressions, each with either length 1 or the same length as in the entire input (i.e., number of iterations or draws). The name of each argument will be the name of a new variable, and the value will be its corresponding value. Use a `NULL` value in `mutate_variables` to drop a variable. New variables overwrite existing variables of the same name.

Details

In order to mutate variables in draws_matrix and draws_array objects, they are transformed to draws_df objects first and then transformed back after mutation. As those transformations are quite expensive for larger number of draws, we recommend using mutate_variables on draws_df and draws_list objects if speed is an issue.

In draws_rvars objects, the output of each expression in ... is coerced to an rvar object if it is not already one using as_rvar().

Value

Returns a draws object of the same format as .x, with variables mutated according to the expressions provided in ....

See Also

variables, rename_variables

Examples

```
x <- as_draws_df(example_draws())
x <- subset(x, variable = c("mu", "tau"))

mutate_variables(x, tau2 = tau^2)
mutate_variables(x, scale = 1.96 * tau, lower = mu - scale)
```

---

order_draws *Order* draws *objects*

Description

Order draws objects according to iteration and chain number. By default, draws objects are ordered but subsetting or extracting parts of them may leave them in an unordered state.

Usage

```
order_draws(x, ...)

## S3 method for class 'draws_matrix'
order_draws(x, ...)

## S3 method for class 'draws_array'
order_draws(x, ...)

## S3 method for class 'draws_df'
order_draws(x, ...)
```

```
## S3 method for class 'draws_list'
order_draws(x, ...)

## S3 method for class 'draws_rvars'
order_draws(x, ...)

## S3 method for class 'rvar'
order_draws(x, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).

Value

A draws object of the same class as x.

See Also

repair_draws()

Examples

```
x <- as_draws_array(example_draws())
dimnames(x[10:5, 4:3, ])
dimnames(order_draws(x[10:5, 4:3, ]))
```

---

pareto_diags *Pareto smoothing diagnostics*

Description

Compute diagnostics for Pareto smoothing the tail draws of x by replacing tail draws by order statistics of a generalized Pareto distribution fit to the tail(s).

Usage

```
pareto_diags(x, ...)

## Default S3 method:
pareto_diags(
  x,
  tail = c("both", "right", "left"),
  r_eff = NULL,
  ndraws_tail = NULL,
  verbose = FALSE,
```

```
  are_log_weights = FALSE,
  ...
)

## S3 method for class 'rvar'
pareto_diags(x, ...)

pareto_khat_threshold(x, ...)

## Default S3 method:
pareto_khat_threshold(x, ...)

## S3 method for class 'rvar'
pareto_khat_threshold(x, ...)

pareto_min_ss(x, ...)

## Default S3 method:
pareto_min_ss(x, ...)

## S3 method for class 'rvar'
pareto_min_ss(x, ...)

pareto_convergence_rate(x, ...)

## Default S3 method:
pareto_convergence_rate(x, ...)

## S3 method for class 'rvar'
pareto_convergence_rate(x, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
...	Arguments passed to individual methods (if applicable).
tail	(string) The tail to diagnose/smooth: • `"right"`: diagnose/smooth only the right (upper) tail • `"left"`: diagnose/smooth only the left (lower) tail • `"both"`: diagnose/smooth both tails and return the maximum k-hat value The default is `"both"`.
r_eff	(numeric) relative effective sample size estimate. If `r_eff` is NULL, it will be calculated assuming the draws are from MCMC. Default is NULL.
ndraws_tail	(numeric) number of draws for the tail. If `ndraws_tail` is not specified, it will be calculated as ceiling(3 * sqrt(length(x) / r_eff)) if length(x) > 225 and length(x) / 5 otherwise (see Appendix H in Vehtari et al. (2024)).

verbose (logical) Should diagnostic messages be printed? If TRUE, messages related to Pareto diagnostics will be printed. Default is FALSE.

are_log_weights (logical) Are the draws log weights? Default is FALSE. If TRUE computation will take into account that the draws are log weights, and only right tail will be smoothed.

Details

When the fitted Generalized Pareto Distribution is used to smooth the tail values and these smoothed values are used to compute expectations, the following diagnostics can give further information about the reliability of these estimates.

- min_ss: Minimum sample size for reliable Pareto smoothed estimate. If the actual sample size is greater than min_ss, then Pareto smoothed estimates can be considered reliable. If the actual sample size is lower than min_ss, increasing the sample size might result in more reliable estimates. For further details, see Section 3.2.3, Equation 11 in Vehtari et al. (2024).
- khat_threshold: Threshold below which k-hat values result in reliable Pareto smoothed estimates. The threshold is lower for smaller effective sample sizes. If k-hat is larger than the threshold, increasing the total sample size may improve reliability of estimates. For further details, see Section 3.2.4, Equation 13 in Vehtari et al. (2024).
- convergence_rate: Relative convergence rate compared to the central limit theorem. Applicable only if the actual sample size is sufficiently large (greater than min_ss). The convergence rate tells the rate at which the variance of an estimate reduces when the sample size is increased, compared to the central limit theorem convergence rate. See Appendix B in Vehtari et al. (2024).

Value

List of Pareto smoothing diagnostics:

- khat: estimated Pareto k shape parameter,
- min_ss: minimum sample size for reliable Pareto smoothed estimate,
- khat_threshold: khat-threshold for reliable Pareto smoothed estimate,
- convergence_rate: Pareto smoothed estimate RMSE convergence rate.

References

Aki Vehtari, Daniel Simpson, Andrew Gelman, Yuling Yao and Jonah Gabry (2024). Pareto Smoothed Importance Sampling. *Journal of Machine Learning Research*, 25(72):1-58. PDF

See Also

pareto_khat, pareto_min_ss, pareto_khat_threshold, and pareto_convergence_rate for individual diagnostics; and pareto_smooth for Pareto smoothing draws.

Other diagnostics: ess_basic(), ess_bulk(), ess_quantile(), ess_sd(), ess_tail(), mcse_mean(), mcse_quantile(), mcse_sd(), pareto_khat(), rhat(), rhat_basic(), rhat_nested(), rstar()

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
pareto_diags(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
pareto_diags(d$Sigma)
```

`pareto_khat` *Pareto khat diagnostic*

Description

Estimate Pareto k value by fitting a Generalized Pareto Distribution to one or two tails of x. This can be used to estimate the number of fractional moments that is useful for convergence diagnostics. For further details see Vehtari et al. (2024).

Usage

```
pareto_khat(x, ...)

## Default S3 method:
pareto_khat(
  x,
  tail = c("both", "right", "left"),
  r_eff = NULL,
  ndraws_tail = NULL,
  verbose = FALSE,
  are_log_weights = FALSE,
  ...
)

## S3 method for class 'rvar'
pareto_khat(x, ...)
```

Arguments

`x`	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
`...`	Arguments passed to individual methods (if applicable).
`tail`	(string) The tail to diagnose/smooth: • `"right"`: diagnose/smooth only the right (upper) tail • `"left"`: diagnose/smooth only the left (lower) tail • `"both"`: diagnose/smooth both tails and return the maximum k-hat value The default is `"both"`.

`r_eff` (numeric) relative effective sample size estimate. If `r_eff` is NULL, it will be calculated assuming the draws are from MCMC. Default is NULL.

`ndraws_tail` (numeric) number of draws for the tail. If `ndraws_tail` is not specified, it will be calculated as ceiling(3 * sqrt(length(x) / r_eff)) if length(x) > 225 and length(x) / 5 otherwise (see Appendix H in Vehtari et al. (2024)).

`verbose` (logical) Should diagnostic messages be printed? If `TRUE`, messages related to Pareto diagnostics will be printed. Default is `FALSE`.

`are_log_weights` (logical) Are the draws log weights? Default is `FALSE`. If `TRUE` computation will take into account that the draws are log weights, and only right tail will be smoothed.

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Aki Vehtari, Daniel Simpson, Andrew Gelman, Yuling Yao and Jonah Gabry (2024). Pareto Smoothed Importance Sampling. *Journal of Machine Learning Research*, 25(72):1-58. PDF

See Also

`pareto_diags` for additional related diagnostics, and `pareto_smooth` for Pareto smoothed draws.

Other diagnostics: `ess_basic()`, `ess_bulk()`, `ess_quantile()`, `ess_sd()`, `ess_tail()`, `mcse_mean()`, `mcse_quantile()`, `mcse_sd()`, `pareto_diags()`, `rhat()`, `rhat_basic()`, `rhat_nested()`, `rstar()`

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
pareto_khat(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
pareto_khat(d$Sigma)
```

pareto_pit *Pareto-smoothed probability integral transform*

Description

Compute PIT values using the empirical CDF, then refine values in the tails by fitting a generalized Pareto distribution (GPD) to the tail draws. This gives smoother, more accurate PIT values in the tails where the ECDF is coarse, and avoids PIT values of 0 and 1. Due to use of generalized Pareto distribution CDF in tails, the PIT values are not anymore rank based and continuous uniformity test is appropriate.

Usage

```
pareto_pit(x, y, ...)

## Default S3 method:
pareto_pit(x, y, weights = NULL, log = FALSE, ndraws_tail = NULL, ...)

## S3 method for class 'draws_matrix'
pareto_pit(x, y, weights = NULL, log = FALSE, ndraws_tail = NULL, ...)

## S3 method for class 'rvar'
pareto_pit(x, y, weights = NULL, log = FALSE, ndraws_tail = NULL, ...)
```

Arguments

x	(draws) A draws_matrix object or one coercible to a draws_matrix object, or an rvar object.
y	(observations) A 1D vector, or an array of dim(x), if x is rvar. Each element of y corresponds to a variable in x.
...	Arguments passed to individual methods (if applicable).
weights	A matrix of weights for each draw and variable. weights should have one column per variable in x, and ndraws(x) rows.
log	(logical) Are the weights passed already on the log scale? The default is FALSE, that is, expecting weights to be on the standard (non-log) scale.
ndraws_tail	(integer) Number of tail draws to use for GPD fitting. If NULL (the default), computed using ps_tail_length().

Details

The function first computes raw PIT values identically to pit() (including support for weighted draws). It then fits a GPD to both tails of the draws (using the same approach as pareto_smooth()) and replaces PIT values for observations falling in the tail regions:

For a right-tail observation $y_i > c_R$ (where c_R is the right-tail cutoff):

$$PIT(y_i) = 1 - p_{tail}(1 - F_{GPD}(y_i; c_R, \sigma_R, k_R))$$

For a left-tail observation $y_i < c_L$:

$$PIT(y_i) = p_{tail}(1 - F_{GPD}(-y_i; -c_L, \sigma_L, k_L))$$

where p_{tail} is the proportion of (weighted) mass in the tail.

When (log-)weights in `weights` are provided, they are used for the raw PIT computation (as in `pit()`) and for GPD fit.

Value

A numeric vector of length `length(y)` containing the PIT values, or an array of shape `dim(y)`, if `x` is an `rvar`.

See Also

`pit()` for the unsmoothed version, `pareto_smooth()` for Pareto smoothing of draws.

Examples

```
x <- example_draws()
y <- rnorm(nvariables(x), 5, 5)
pareto_pit(x, y)
```

---

pareto_smooth *Pareto smoothing*

Description

Smooth the tail draws of x by replacing tail draws by order statistics of a generalized Pareto distribution fit to the tail(s). For further details see Vehtari et al. (2024).

Usage

```
pareto_smooth(x, ...)

## S3 method for class 'rvar'
pareto_smooth(x, return_k = FALSE, extra_diags = FALSE, ...)

## Default S3 method:
pareto_smooth(
  x,
  tail = c("both", "right", "left"),
  r_eff = NULL,
  ndraws_tail = NULL,
```

```
  return_k = FALSE,
  extra_diags = FALSE,
  verbose = TRUE,
  are_log_weights = FALSE,
  ...
)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
...	Arguments passed to individual methods (if applicable).
return_k	(logical) Should the Pareto khat be included in output? If `TRUE`, output will be a list containing smoothed draws and diagnostics, otherwise it will be a numeric of the smoothed draws. Default is `FALSE`.
extra_diags	(logical) Should extra Pareto khat diagnostics be included in output? If `TRUE`, `min_ss`, `khat_threshold` and `convergence_rate` for the estimated k value will be returned. Default is `FALSE`.
tail	(string) The tail to diagnose/smooth: • `"right"`: diagnose/smooth only the right (upper) tail • `"left"`: diagnose/smooth only the left (lower) tail • `"both"`: diagnose/smooth both tails and return the maximum k-hat value The default is `"both"`.
r_eff	(numeric) relative effective sample size estimate. If `r_eff` is NULL, it will be calculated assuming the draws are from MCMC. Default is NULL.
ndraws_tail	(numeric) number of draws for the tail. If `ndraws_tail` is not specified, it will be calculated as ceiling(3 * sqrt(length(x) / r_eff)) if length(x) > 225 and length(x) / 5 otherwise (see Appendix H in Vehtari et al. (2024)).
verbose	(logical) Should diagnostic messages be printed? If `TRUE`, messages related to Pareto diagnostics will be printed. Default is `FALSE`.
are_log_weights	(logical) Are the draws log weights? Default is `FALSE`. If `TRUE` computation will take into account that the draws are log weights, and only right tail will be smoothed.

Value

Either a vector `x` of smoothed values or a named list containing the vector `x` and a named list `diagnostics` containing numeric values:

- `khat`: estimated Pareto k shape parameter, and optionally
- `min_ss`: minimum sample size for reliable Pareto smoothed estimate
- `khat_threshold`: sample size specific khat threshold for reliable Pareto smoothed estimates
- `convergence_rate`: Relative convergence rate for Pareto smoothed estimates

If any of the draws is non-finite, that is, NA, NaN, Inf, or -Inf, Pareto smoothing will not be performed, and the original draws will be returned and and diagnostics will be NA (numeric).

References

Aki Vehtari, Daniel Simpson, Andrew Gelman, Yuling Yao and Jonah Gabry (2024). Pareto Smoothed Importance Sampling. *Journal of Machine Learning Research*, 25(72):1-58. PDF

See Also

pareto_khat for only calculating khat, and pareto_diags for additional diagnostics.

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
pareto_smooth(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
pareto_smooth(d$Sigma)
```

---

pit *Probability integral transform*

Description

Probability integral transform (PIT). LOO-PIT is given by a weighted sample.

Usage

```
pit(x, y, ...)

## Default S3 method:
pit(x, y, weights = NULL, log = FALSE, ...)

## S3 method for class 'draws_matrix'
pit(x, y, weights = NULL, log = FALSE, ...)

## S3 method for class 'rvar'
pit(x, y, weights = NULL, log = FALSE, ...)
```

Arguments

x	(draws) A draws_matrix object or one coercible to a draws_matrix object, or an rvar object.
y	(observations) A 1D vector, or an array of dim(x), if x is rvar. Each element of y corresponds to a variable in x.
...	Arguments passed to individual methods (if applicable).

weights	A matrix of weights for each draw and variable. weights should have one column per variable in x, and ndraws(x) rows.
log	(logical) Are the weights passed already on the log scale? The default is FALSE, that is, expecting weights to be on the standard (non-log) scale.

Details

The pit() function computes the probability integral transform of y using the empirical cumulative distribution computed from the draws in x. For continuous valued y and x, the PIT for the elements of y is computed as the empirical cumulative distribution value:

```
PIT(y_i) = Pr(x_i < y_i),
```

where x_i, is the corresponding set of draws in x. For draws objects, this corresponds to the draws of the *i*th variable, and for rvar the elements of y and x are matched.

The draws in x can further be provided (log-)weights in

If y and x are discrete, randomization is used to obtain continuous PIT values. (see, e.g., Czado, C., Gneiting, T., Held, L.: Predictive model assessment for count data. Biometrics 65(4), 1254–1261 (2009).)

Value

A numeric vector of length length(y) containing the PIT values, or an array of shape dim(y), if x is an rvar.

Examples

```
# PIT for a draws object
x <- example_draws()
# Create a vector of observations
y <- rnorm(nvariables(x), 5, 5)
pit(x, y)

# Compute weighted PIT (for example LOO-PIT)
weights <- matrix(runif(length(x)), ncol = nvariables(x))

pit(x, y, weights)

# PIT for an rvar
x <- rvar(example_draws())
# Create an array of observations with the same dimensions as x.
y_arr <- array(rnorm(length(x), 5, 5), dim = dim(x))
pit(x, y_arr)
```

print.draws_array *Print* draws_array *objects*

Description

Pretty printing for draws_array objects.

Usage

```
## S3 method for class 'draws_array'
print(
  x,
  digits = 2,
  max_iterations = getOption("posterior.max_iterations", 5),
  max_chains = getOption("posterior.max_chains", 8),
  max_variables = getOption("posterior.max_variables", 4),
  reserved = FALSE,
  ...
)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
digits	(nonnegative integer) The minimum number of significant digits to print. If NULL, defaults to getOption("posterior.digits", 2).
max_iterations	(positive integer) The maximum number of iterations to print. Can be controlled globally via the "posterior.max_iterations" option.
max_chains	(positive integer) The maximum number of chains to print. Can be controlled globally via the "posterior.max_chains" option.
max_variables	(positive integer) The maximum number of variables to print. Can be controlled globally via the "posterior.max_variables" option.
reserved	(logical) Should reserved variables be included in the output? Defaults to FALSE. See reserved_variables for an overview of currently reserved variable names.
...	Further arguments passed to the underlying print() methods.

Value

A draws object of the same class as x.

Examples

```
x <- as_draws_array(example_draws())
print(x)
```

print.draws_df *Print* draws_df *objects*

Description

Pretty printing for draws_df objects.

Usage

```
## S3 method for class 'draws_df'
print(
  x,
  digits = 2,
  max_draws = getOption("posterior.max_draws", 10),
  max_variables = getOption("posterior.max_variables", 8),
  reserved = FALSE,
  ...
)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
digits	(nonnegative integer) The minimum number of significant digits to print. If NULL, defaults to getOption("posterior.digits", 2).
max_draws	(positive integer) The maximum number of draws to print. Can be controlled globally via the "posterior.max_draws" option.
max_variables	(positive integer) The maximum number of variables to print. Can be controlled globally via the "posterior.max_variables" option.
reserved	(logical) Should reserved variables be included in the output? Defaults to FALSE. See reserved_variables for an overview of currently reserved variable names.
...	Further arguments passed to the underlying print() methods.

Value

A draws object of the same class as x.

Examples

```
x <- as_draws_df(example_draws())
print(x)
```

print.draws_list *Print* draws_list *objects*

Description

Pretty printing for draws_list objects.

Usage

```
## S3 method for class 'draws_list'
print(
  x,
  digits = 2,
  max_iterations = getOption("posterior.max_iterations", 10),
  max_chains = getOption("posterior.max_chains", 2),
  max_variables = getOption("posterior.max_variables", 4),
  reserved = FALSE,
  ...
)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
digits	(nonnegative integer) The minimum number of significant digits to print. If NULL, defaults to getOption("posterior.digits", 2).
max_iterations	(positive integer) The maximum number of iterations to print. Can be controlled globally via the "posterior.max_iterations" option.
max_chains	(positive integer) The maximum number of chains to print. Can be controlled globally via the "posterior.max_chains" option.
max_variables	(positive integer) The maximum number of variables to print. Can be controlled globally via the "posterior.max_variables" option.
reserved	(logical) Should reserved variables be included in the output? Defaults to FALSE. See reserved_variables for an overview of currently reserved variable names.
...	Further arguments passed to the underlying print() methods.

Value

A draws object of the same class as x.

Examples

```
x <- as_draws_list(example_draws())
print(x)
```

print.draws_matrix *Print* draws_matrix *objects*

Description

Pretty printing for draws_matrix objects.

Usage

```
## S3 method for class 'draws_matrix'
print(
  x,
  digits = 2,
  max_draws = getOption("posterior.max_draws", 10),
  max_variables = getOption("posterior.max_variables", 8),
  reserved = FALSE,
  ...
)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
digits	(nonnegative integer) The minimum number of significant digits to print. If NULL, defaults to getOption("posterior.digits", 2).
max_draws	(positive integer) The maximum number of draws to print. Can be controlled globally via the "posterior.max_draws" option.
max_variables	(positive integer) The maximum number of variables to print. Can be controlled globally via the "posterior.max_variables" option.
reserved	(logical) Should reserved variables be included in the output? Defaults to FALSE. See reserved_variables for an overview of currently reserved variable names.
...	Further arguments passed to the underlying print() methods.

Value

A draws object of the same class as x.

Examples

```
x <- as_draws_matrix(example_draws())
print(x)
```

print.draws_rvars *Print* draws_rvars *objects*

Description

Pretty printing for draws_rvars objects.

Usage

```
## S3 method for class 'draws_rvars'
print(
  x,
  digits = 2,
  max_variables = getOption("posterior.max_variables", 8),
  summary = getOption("posterior.rvar_summary", "mean_sd"),
  reserved = FALSE,
  ...
)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
digits	(nonnegative integer) The minimum number of significant digits to print. If NULL, defaults to getOption("posterior.digits", 2).
max_variables	(positive integer) The maximum number of variables to print. Can be controlled globally via the "posterior.max_variables" option.
summary	(string) The style of summary to display: • "mean_sd" displays mean ± sd • "median_mad" displays median ± mad • "mode_entropy" displays mode <entropy>, and is used automatically for rvar_factors. It shows normalized entropy, which ranges from 0 (all probability in one category) to 1 (uniform). See entropy(). • "mode_dissent" displays mode <dissent>, and is used automatically for rvar_ordereds. It shows Tastle and Wierman's (2007) *dissention* measure, which ranges from 0 (all probability in one category) through 0.5 (uniform) to 1 (bimodal: all probability split equally between the first and last category). See dissent(). • NULL uses getOption("posterior.rvar_summary") (default "mean_sd")
reserved	(logical) Should reserved variables be included in the output? Defaults to FALSE. See reserved_variables for an overview of currently reserved variable names.
...	Further arguments passed to the underlying print() methods.

Value

A draws object of the same class as x.

Examples

```
x <- as_draws_rvars(example_draws())
print(x)
```

`print.draws_summary` *Print summaries of* `draws` *objects*

Description

Print output from `summarise_draws()`.

Usage

```
## S3 method for class 'draws_summary'
print(x, ..., num_args = NULL)
```

Arguments

`x`	(draws_summary) A `"draws_summary"` object as output by `summarise_draws()`.
`...`	Additional arguments passed to `tibble::print.tbl_df()`
`num_args`	(named list) Optional arguments passed to num() for pretty printing of summaries. If `NULL` (the default), uses the arguments stored in the `"num_args"` attribute of `x`, as set by the `.num_args` argument of `summarise_draws()`, which itself can be controlled globally via the `posterior.num_args` option.

Value

An invisible version of the input object.

Examples

```
x <- example_draws("eight_schools")

# adjust how summaries are printed when calling summarise_draws()...
summarise_draws(x, .num_args = list(sigfig = 2, notation = "dec"))

# ... or when printing
s <- summarise_draws(x)
print(s, num_args = list(sigfig = 2, notation = "dec"))
print(s, num_args = list(digits = 3))
```

print.rvar *Print or format a random variable*

Description

Printing and formatting methods for rvars.

Usage

```
## S3 method for class 'rvar'
print(
  x,
  ...,
  summary = NULL,
  digits = NULL,
  color = TRUE,
  width = getOption("width")
)

## S3 method for class 'rvar'
format(x, ..., summary = NULL, digits = NULL, color = FALSE)

## S3 method for class 'rvar'
str(
  object,
  ...,
  summary = NULL,
  vec.len = NULL,
  indent.str = paste(rep.int(" ", max(0, nest.lev + 1)), collapse = ".."),
  nest.lev = 0,
  give.attr = TRUE
)
```

Arguments

x, object	(rvar) The rvar to print.
...	Further arguments passed to the underlying print() methods.
summary	(string) The style of summary to display: • "mean_sd" displays mean ± sd • "median_mad" displays median ± mad • "mode_entropy" displays mode <entropy>, and is used automatically for rvar_factors. It shows normalized entropy, which ranges from 0 (all probability in one category) to 1 (uniform). See entropy(). • "mode_dissent" displays mode <dissent>, and is used automatically for rvar_ordereds. It shows Tastle and Wierman's (2007) *dissention* measure, which ranges from 0 (all probability in one category) through 0.5 (uniform)

to 1 (bimodal: all probability split equally between the first and last category). See dissent().
- NULL uses getOption("posterior.rvar_summary") (default "mean_sd")

digits — (nonnegative integer) The minimum number of significant digits to print. If NULL, defaults to getOption("posterior.digits", 2).

color — (logical) Whether or not to use color when formatting the output. If TRUE, the pillar::style_num() functions may be used to produce strings containing control sequences to produce colored output on the terminal.

width — The maximum width used to print out lists of factor levels for rvar_factors. See format().

vec.len — (nonnegative integer) How many 'first few' elements are displayed of each vector. If NULL, defaults to getOption("str")$vec.len, which defaults to 4.

indent.str — (string) The indentation string to use.

nest.lev — (nonnegative integer) Current nesting level in the recursive calls to str().

give.attr — (logical) If TRUE (default), show attributes as sub structures.

Details

print() and str() print out rvar objects by summarizing each element in the random variable with either its mean±sd or median±mad, depending on the value of summary. Both functions use the format() implementation for rvar objects under the hood, which returns a character vector in the mean±sd or median±mad form.

Value

For print(), an invisible version of the input object.

For str(), nothing; i.e. invisible(NULL).

For format(), a character vector of the same dimensions as x where each entry is of the form "mean±sd" or "median±mad", depending on the value of summary.

References

William J. Tastle, Mark J. Wierman (2007). Consensus and dissention: A measure of ordinal dispersion. *International Journal of Approximate Reasoning*. 45(3), 531–545. doi:10.1016/j.ijar.2006.06.024.

Examples

```
set.seed(5678)
x = rbind(
  cbind(rvar(rnorm(1000, 1)), rvar(rnorm(1000, 2))),
  cbind(rvar(rnorm(1000, 3)), rvar(rnorm(1000, 4)))
)

print(x)
print(x, summary = "median_mad")

str(x)
```

```
format(x)
```

`ps_convergence_rate` *Pareto convergence rate*

Description

Given number of draws and scalar or array of k's, compute the relative convergence rate of PSIS estimate RMSE. See Appendix B of Vehtari et al. (2024). This function is exported to be usable by other packages. For user-facing diagnostic functions, see `pareto_convergence_rate` and `pareto_diags`.

Usage

```
ps_convergence_rate(k, ndraws, ...)
```

Arguments

`k`	pareto-k values
`ndraws`	number of draws
`...`	unused

Value

convergence rate

See Also

Other helper-functions: `gpdfit()`, `ps_khat_threshold()`, `ps_min_ss()`, `ps_tail_length()`

`ps_khat_threshold` *Pareto k-hat threshold*

Description

Given number of draws, computes khat threshold for reliable Pareto smoothed estimate (to have small probability of large error). See section 3.2.4, equation (13) of Vehtari et al. (2024). This function is exported to be usable by other packages. For user-facing diagnostic functions, see `pareto_khat_threshold` and `pareto_diags`.

Usage

```
ps_khat_threshold(ndraws, ...)
```

Arguments

ndraws	number of draws
...	unused

Value

threshold

See Also

Other helper-functions: gpdfit(), ps_convergence_rate(), ps_min_ss(), ps_tail_length()

---

ps_min_ss *Pareto-smoothing minimum sample-size*

Description

Given Pareto-k computes the minimum sample size for reliable Pareto smoothed estimate (to have small probability of large error). See section 3.2.3 in Vehtari et al. (2024). This function is exported to be usable by other packages. For user-facing diagnostic functions, see pareto_min_ss and pareto_diags.

Usage

```
ps_min_ss(k, ...)
```

Arguments

k	pareto k value
...	unused

Value

minimum sample size

See Also

Other helper-functions: gpdfit(), ps_convergence_rate(), ps_khat_threshold(), ps_tail_length()

`ps_tail` *Pareto smooth tail function to Pareto smooth the tail of a vector. Exported for usage in other packages, not by users.*

Description

Pareto smooth tail function to Pareto smooth the tail of a vector. Exported for usage in other packages, not by users.

Usage

```
ps_tail(
  x,
  ndraws_tail,
  smooth_draws = TRUE,
  tail = c("right", "left"),
  are_log_weights = FALSE,
  ...
)
```

Arguments

`x`	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
`ndraws_tail`	(numeric) number of draws for the tail. If `ndraws_tail` is not specified, it will be set to `length(x)`.
`smooth_draws`	(logical) Should the tails be smoothed? Default is `TRUE`. If `FALSE`, `k` will be calculated but `x` will remain untouched.
`tail`	(string) The tail to diagnose/smooth: • `"right"`: diagnose/smooth only the right (upper) tail • `"left"`: diagnose/smooth only the left (lower) tail
`are_log_weights`	(logical) Are the draws log weights? Default is `FALSE`. If `TRUE` computation will take into account that the draws are log weights, and only right tail will be smoothed.
`...`	Arguments passed to individual methods (if applicable).

References

Aki Vehtari, Daniel Simpson, Andrew Gelman, Yuling Yao and Jonah Gabry (2024). Pareto Smoothed Importance Sampling. *Journal of Machine Learning Research*, 25(72):1-58. PDF

See Also

`pareto_smooth` for the user-facing function.

ps_tail_length *Pareto tail length*

Description

Calculate the tail length from number of draws and relative efficiency r_eff. See Appendix H in Vehtari et al. (2024). This function is used internally and is exported to be available for other packages.

Usage

```
ps_tail_length(ndraws, r_eff, ...)
```

Arguments

ndraws	number of draws
r_eff	relative efficiency
...	unused

Value

tail length

See Also

Other helper-functions: gpdfit(), ps_convergence_rate(), ps_khat_threshold(), ps_min_ss()

quantile2 *Compute Quantiles*

Description

Compute quantiles of a sample and return them in a format consistent with other summary functions in the **posterior** package.

Usage

```
quantile2(x, probs = c(0.05, 0.95), na.rm = FALSE, ...)

## Default S3 method:
quantile2(x, probs = c(0.05, 0.95), na.rm = FALSE, names = TRUE, ...)

## S3 method for class 'rvar'
quantile2(x, probs = c(0.05, 0.95), na.rm = FALSE, names = TRUE, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See extract_variable_matrix(). • An rvar.
probs	(numeric vector) Probabilities in [0, 1].
na.rm	(logical) Should NA and NaN values be removed from x prior to computing quantiles? The default is FALSE.
...	Arguments passed to individual methods (if applicable) and then on to stats::quantile().
names	(logical) Should the result have a names attribute? The default is TRUE, but use FALSE for improved speed if there are many values in probs.

Value

A numeric vector of length length(probs). If names = TRUE, it has a names attribute with names like "q5", "q95", etc, based on the values of probs.

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
quantile2(mu)
```

---

rdo *Execute expressions of random variables*

Description

Execute (nearly) arbitrary R expressions that may include rvars, producing a new rvar.

Usage

```
rdo(expr, dim = NULL, ndraws = NULL)
```

Arguments

expr	(expression) A bare expression that can (optionally) contain rvars. The expression supports quasiquotation.
dim	(integer vector) One or more integers giving the maximal indices in each dimension to override the dimensions of the rvar to be created (see dim()). If NULL (the default), dim is determined by the input. **NOTE:** This argument controls the dimensions of the rvar, not the underlying array, so you cannot change the number of draws using this argument.
ndraws	(positive integer) The number of draws used to construct new random variables if no rvars are supplied in expr. If NULL, getOption("posterior.rvar_ndraws") is used (default 4000). If expr contains rvars, the number of draws in the provided rvars is used instead of the value of this argument.

Details

This function evaluates `expr` possibly multiple times, once for each draw of the `rvar`s it contains, then returns a new `rvar` representing the output of those expressions. To identify `rvar`s, `rdo()` searches the calling environment for any variables named in `expr` for which `is_rvar()` evaluates to `TRUE`. If `expr` contains no `rvar`s, then it will be executed `ndraws` times and an `rvar` with that many draws returned.

`rdo()` is not necessarily *fast* (in fact in some cases it may be very slow), but it has the advantage of allowing a nearly arbitrary R expression to be executed against `rvar`s simply by wrapping it with `rdo( ... )`. This makes it especially useful as a prototyping tool. If you create code with `rdo()` and it is unacceptably slow for your application, consider rewriting it using math operations directly on `rvar`s (which should be fast), using `rvar_rng()`, and/or using operations directly on the arrays that back the `rvar`s (via `draws_of()`).

Value

An `rvar`.

See Also

Other rfun: `rfun()`, `rvar_rng()`

Examples

```
mu <- rdo(rnorm(10, mean = 1:10, sd = 1))
sigma <- rdo(rgamma(1, shape = 1, rate = 1))
x <- rdo(rnorm(10, mu, sigma))
x
```

---

rename_variables *Rename variables in* draws *objects*

Description

Rename variables in a `draws` object.

Usage

```
rename_variables(.x, ...)

## S3 method for class 'draws'
rename_variables(.x, ...)
```

Arguments

.x	(draws) A draws object.
...	One or more expressions, separated by commas, indicating the variables to rename. The variable names can be unquoted (new_name = old_name) or quoted ("new_name" = "old_name"). For non-scalar variables, all elements can be renamed together ("new_name" = "old_name") or they can be renamed individually ("new_name[1]" = "old_name[1]").

Value

Returns a draws object of the same format as .x, with variables renamed according to the expressions provided in ....

See Also

variables, variables<-, mutate_variables

Examples

```
x <- as_draws_df(example_draws())
variables(x)

x <- rename_variables(x, mean = mu, sigma = tau)
variables(x)

x <- rename_variables(x, b = `theta[1]`) # or b  = "theta[1]"
variables(x)

# rename all elements of 'theta' at once
x <- rename_variables(x, alpha = theta)
variables(x)
```

---

repair_draws *Repair indices of* draws *objects*

Description

Repair indices of draws objects so that iterations, chains, and draws are continuously and consistently numbered.

Usage

```
repair_draws(x, order = TRUE, ...)

## S3 method for class 'draws_matrix'
repair_draws(x, order = TRUE, ...)
```

```
## S3 method for class 'draws_array'
repair_draws(x, order = TRUE, ...)

## S3 method for class 'draws_df'
repair_draws(x, order = TRUE, ...)

## S3 method for class 'draws_list'
repair_draws(x, order = TRUE, ...)

## S3 method for class 'draws_rvars'
repair_draws(x, order = TRUE, ...)

## S3 method for class 'rvar'
repair_draws(x, order = TRUE, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
order	(logical) Should draws be ordered (via order_draws()) before repairing indices? Defaults to TRUE.
...	Arguments passed to individual methods (if applicable).

Value

A draws object of the same class as x.

See Also

order_draws()

Examples

```
x <- as_draws_array(example_draws())
(x <- x[10:5, 3:4, ])
repair_draws(x)
```

---

resample_draws *Resample* draws *objects*

Description

Resample draws objects according to provided weights, for example weights obtained through importance sampling.

Usage

```
resample_draws(x, ...)

## S3 method for class 'draws'
resample_draws(x, weights = NULL, method = "stratified", ndraws = NULL, ...)

## S3 method for class 'rvar'
resample_draws(x, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).
weights	(numeric vector) A vector of positive weights of length ndraws(x). The weights will be internally normalized. If weights is not specified, an attempt will be made to extract any weights already stored in the draws object (via weight_draws()). If no weights are stored in the draws object, equal weight is supplied to each draw. How exactly the weights influence the resampling depends on the method argument.
method	(string) The resampling method to use: • "simple": simple random resampling with replacement • "simple_no_replace": simple random resampling without replacement • "stratified": stratified resampling with replacement • "deterministic": deterministic resampling with replacement Currently, "stratified" is the default as it has comparably low variance and bias with respect to ideal resampling. The latter would sample perfectly proportional to the weights, but this is not possible in practice due to the finite number of draws available. For more details about resampling methods, see Kitagawa (1996).
ndraws	(positive integer) The number of draws to be returned. By default ndraws is set internally to the total number of draws in x if sensible.

Details

Upon usage of resample_draws(), chains will automatically be merged due to subsetting of individual draws (see subset_draws for details). Also, weights stored in the draws object will be removed in the process, as resampling invalidates existing weights.

Value

A draws object of the same class as x.

References

Kitagawa, G., Monte Carlo Filter and Smoother for Non-Gaussian Nonlinear ' State Space Models, *Journal of Computational and Graphical Statistics*, 5(1):1-25, 1996.

See Also

resample_draws()

Examples

```
x <- as_draws_df(example_draws())

# random weights for justr for demonstration
w <- runif(ndraws(x), 0, 10)

# use default stratified sampling
x_rs <- resample_draws(x, weights = w)
summarise_draws(x_rs, default_summary_measures())

# use simple random sampling
x_rs <- resample_draws(x, weights = w, method = "simple")
summarise_draws(x_rs, default_summary_measures())
```

---

reserved_variables *Reserved variables*

Description

Get names of reserved variables from objects in the **posterior** package.

Usage

```
reserved_variables(x, ...)

## Default S3 method:
reserved_variables(x, ...)

## S3 method for class 'draws_matrix'
reserved_variables(x, ...)

## S3 method for class 'draws_array'
reserved_variables(x, ...)

## S3 method for class 'draws_df'
reserved_variables(x, ...)

## S3 method for class 'draws_list'
reserved_variables(x, ...)

## S3 method for class 'draws_rvars'
reserved_variables(x, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).

Details

`reserved_variables()` returns the names of reserved variables in use by an object.

The following variables names are currently reserved for special use cases in all draws formats:

- `.log_weight`: Log weights per draw (see weight_draws).

Further, specific for the draws_df format, there are three additional reserved variables:

- `.chain`: Chain index per draw
- `.iteration`: Iteration index within each chain
- `.draw`: Draw index across chains

More reserved variables may be added in the future.

Value

A character vector of reserved variables used in x.

Examples

```
x <- example_draws()
reserved_variables(x)

# if we add weights, the `.log_weight` reserved variable is used
x <- weight_draws(x, rexp(ndraws(x)))
reserved_variables(x)
```

---

rfun *Create functions of random variables*

Description

Function that create functions that can accept and/or produce rvars.

Usage

```
rfun(.f, rvar_args = NULL, rvar_dots = TRUE, ndraws = NULL)
```

Arguments

`.f`	(multiple options) A function to turn into a function that accepts and/or produces random variables: • A function • A one-sided formula that can be parsed by `rlang::as_function()`
`rvar_args`	(character vector) The names of the arguments of `.f` that should be allowed to accept `rvar`s as arguments. If `NULL` (the default), all arguments to `.f` are turned into arguments that accept `rvar`s, including arguments passed via `...` (if `rvar_dots` is `TRUE`).
`rvar_dots`	(logical) Should dots (`...`) arguments also be converted? Only applies if `rvar_args` is `NULL` (i.e., all arguments are allowed to be `rvar`s).
`ndraws`	(positive integer). The number of draws used to construct new random variables if no `rvar`s are supplied as arguments to the returned function. If `NULL`, `getOption("posterior.rvar_ndraws")` is used (default `4000`). If any arguments to the returned function contain `rvar`s, the number of draws in the provided `rvar`s is used instead of the value of this argument.

Details

This function wraps an existing function (`.f`) such that it returns `rvar`s containing whatever type of data `.f` would normally return.

The returned function, when called, executes `.f` possibly multiple times, once for each draw of the `rvar`s passed to it, then returns a new `rvar` representing the output of those function evaluations. If the arguments contain no `rvar`s, then `.f` will be executed `ndraws` times and an `rvar` with that many draws returned.

Functions created by `rfun()` are not necessarily *fast* (in fact in some cases they may be very slow), but they have the advantage of allowing a nearly arbitrary R functions to be executed against `rvar`s simply by wrapping them with `rfun()`. This makes it especially useful as a prototyping tool. If you create code with `rfun()` and it is unacceptably slow for your application, consider rewriting it using math operations directly on `rvar`s (which should be fast), using `rvar_rng()`, and/or using operations directly on the arrays that back the `rvar`s (via `draws_of()`).

Value

A function with the same argument specification as `.f`, but which can accept and return `rvar`s.

See Also

Other rfun: `rdo()`, `rvar_rng()`

Examples

```
rvar_norm <- rfun(rnorm)
rvar_gamma <- rfun(rgamma)

mu <- rvar_norm(10, mean = 1:10, sd = 1)
sigma <- rvar_gamma(1, shape = 1, rate = 1)
```

```
x <- rvar_norm(10, mu, sigma)
x
```

`rhat` *Rhat convergence diagnostic*

Description

Compute the Rhat convergence diagnostic for a single variable as the maximum of rank normalized split-Rhat and rank normalized folded-split-Rhat as proposed in Vehtari et al. (2021).

Usage

```
rhat(x, ...)

## Default S3 method:
rhat(x, ...)

## S3 method for class 'rvar'
rhat(x, ...)
```

Arguments

`x`	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
`...`	Arguments passed to individual methods (if applicable).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

See Also

Other diagnostics: ess_basic(), ess_bulk(), ess_quantile(), ess_sd(), ess_tail(), mcse_mean(), mcse_quantile(), mcse_sd(), pareto_diags(), pareto_khat(), rhat_basic(), rhat_nested(), rstar()

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
rhat(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
rhat(d$Sigma)
```

---

rhat_basic *Basic version of the Rhat convergence diagnostic*

Description

Compute the basic Rhat convergence diagnostic for a single variable as described in Gelman et al. (2013) with some changes according to Vehtari et al. (2021). For practical applications, we strongly recommend the improved Rhat convergence diagnostic implemented in rhat().

Usage

```
rhat_basic(x, ...)

## Default S3 method:
rhat_basic(x, split = TRUE, ...)

## S3 method for class 'rvar'
rhat_basic(x, split = TRUE, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See extract_variable_matrix(). • An rvar.
...	Arguments passed to individual methods (if applicable).
split	(logical) Should the estimate be computed on split chains? The default is TRUE.

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Andrew Gelman, John B. Carlin, Hal S. Stern, David B. Dunson, Aki Vehtari and Donald B. Rubin (2013). *Bayesian Data Analysis, Third Edition*. Chapman and Hall/CRC.

Aki Vehtari, Andrew Gelman, Daniel Simpson, Bob Carpenter, and Paul-Christian Bürkner (2021). Rank-normalization, folding, and localization: An improved R-hat for assessing convergence of MCMC (with discussion). *Bayesian Analysis*. 16(2), 667-–718. doi:10.1214/20-BA1221

See Also

Other diagnostics: `ess_basic()`, `ess_bulk()`, `ess_quantile()`, `ess_sd()`, `ess_tail()`, `mcse_mean()`, `mcse_quantile()`, `mcse_sd()`, `pareto_diags()`, `pareto_khat()`, `rhat()`, `rhat_nested()`, `rstar()`

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
rhat_basic(mu)

d <- as_draws_rvars(example_draws("multi_normal"))
rhat_basic(d$Sigma)
```

---

rhat_nested *Nested Rhat convergence diagnostic*

Description

Compute the nested Rhat convergence diagnostic for a single variable as proposed in Margossian et al. (2024).

Usage

```
rhat_nested(x, ...)

## Default S3 method:
rhat_nested(x, superchain_ids, ...)

## S3 method for class 'rvar'
rhat_nested(x, superchain_ids, ...)
```

Arguments

x	(multiple options) One of: • A matrix of draws for a single variable (iterations x chains). See `extract_variable_matrix()`. • An `rvar`.
...	Arguments passed to individual methods (if applicable).
superchain_ids	(numeric) Vector of length nchains specifying which superchain each chain belongs to. There should be equal numbers of chains in each superchain. All chains within the same superchain are assumed to have been initialized at the same point.

Details

Nested Rhat is a convergence diagnostic useful when running many short chains. It is calculated on superchains, which are groups of chains that have been initialized at the same point.

Note that there is a slight difference in the calculation of Rhat and nested Rhat, as nested Rhat is lower bounded by 1. This means that nested Rhat with one chain per superchain will not be exactly equal to basic Rhat (see Footnote 3 in Margossian et al. (2024)).

Value

If the input is an array, returns a single numeric value. If any of the draws is non-finite, that is, `NA`, `NaN`, `Inf`, or `-Inf`, the returned output will be (numeric) `NA`. Also, if all draws within any of the chains of a variable are the same (constant), the returned output will be (numeric) `NA` as well. The reason for the latter is that, for constant draws, we cannot distinguish between variables that are supposed to be constant (e.g., a diagonal element of a correlation matrix is always 1) or variables that just happened to be constant because of a failure of convergence or other problems in the sampling process.

If the input is an `rvar`, returns an array of the same dimensions as the `rvar`, where each element is equal to the value that would be returned by passing the draws array for that element of the `rvar` to this function.

References

Charles C. Margossian, Matthew D. Hoffman, Pavel Sountsov, Lionel Riou-Durand, Aki Vehtari and Andrew Gelman (2024). Nested R-hat: Assessing the convergence of Markov chain Monte Carlo when running many short chains. *Bayesian Analysis*. doi:10.1214/24-BA1453

See Also

Other diagnostics: ess_basic(), ess_bulk(), ess_quantile(), ess_sd(), ess_tail(), mcse_mean(), mcse_quantile(), mcse_sd(), pareto_diags(), pareto_khat(), rhat(), rhat_basic(), rstar()

Examples

```
mu <- extract_variable_matrix(example_draws(), "mu")
rhat_nested(mu, superchain_ids = c(1, 1, 2, 2))

d <- as_draws_rvars(example_draws("multi_normal"))
rhat_nested(d$Sigma, superchain_ids = c(1, 1, 2, 2))
```

---

rstar *Calculate R* convergence diagnostic*

Description

The rstar() function generates a measure of convergence for MCMC draws based on whether it is possible to determine the Markov chain that generated a draw with probability greater than chance. To do so, it fits a machine learning classifier to a training set of MCMC draws and evaluates its predictive accuracy on a testing set: giving the ratio of accuracy to predicting a chain uniformly at random.

Usage

```
rstar(
  x,
  split = TRUE,
  uncertainty = FALSE,
  method = "rf",
  hyperparameters = NULL,
  training_proportion = 0.7,
  nsimulations = 1000,
  ...
)
```

Arguments

x	(draws) A draws_df object or one coercible to a draws_df object.
split	(logical) Should the estimate be computed on split chains? The default is TRUE.
uncertainty	(logical). Indicates whether to provide a vector of R* values representing uncertainty in the calculated value (if TRUE) or a single value (if FALSE). The default is TRUE.
method	(string) The machine learning classifier to use (must be available in the **caret** package). The default is "rf", which calls the random forest classifier.

`hyperparameters`
: (named list) Hyperparameter settings passed to the classifier. The default for the random forest classifier (`method = "rf"`) is `list(mtry = floor(sqrt(nvariables(x))))`. The default for the gradient-based model (`method = "gbm"`) is `list(interaction.depth = 3, n.trees = 50, shrinkage = 0.1, n.minobsinnode = 10)`.

`training_proportion`
: (positive real) The proportion (in `(0,1)`) of iterations in used to train the classifier. The default is `0.7`.

`nsimulations`
: (positive integer) The number of R* values in the returned vector if `uncertainty` is `TRUE`. The default is `1000`.

`...`
: Other arguments passed to `caret::train()`.

Details

The `rstar()` function provides a measure of MCMC convergence based on whether it is possible to determine the chain that generated a particular draw with a probability greater than chance. To do so, it fits a machine learning classifier to a subset of the original MCMC draws (the training set) and evaluates its predictive accuracy on the remaining draws (the testing set). If predictive accuracy exceeds chance (i.e. predicting the chain that generated a draw uniformly at random), the diagnostic measure R* will be above 1, indicating that convergence has yet to occur. This statistic is recently developed, and it is currently unclear what is a reasonable threshold for diagnosing convergence.

The statistic, R*, is stochastic, meaning that each time the test is run, unless the random seed is fixed, it will generally produce a different result. To minimize the implications of this stochasticity, it is recommended to repeatedly run this function to calculate a distribution of R*; alternatively, an approximation to this distribution can be obtained by setting `uncertainty = TRUE`, although this approximation of uncertainty will generally have a lower mean.

By default, a random forest classifier is used (`method = "rf"`), which tends to perform best for target distributions of around 4 dimensions and above. For lower dimensional targets, gradient boosted models (called via `method = "gbm"`) tend to have a higher classification accuracy. On a given MCMC sample, it is recommended to try both of these classifiers.

Value

A numeric vector of length 1 (by default) or length `nsimulations` (if `uncertainty = TRUE`).

References

Ben Lambert, Aki Vehtari (2020) R*: A robust MCMC convergence diagnostic with uncertainty using gradient-boosted machines. *arXiv preprint* `arXiv:2003.07900`.

See Also

Other diagnostics: `ess_basic()`, `ess_bulk()`, `ess_quantile()`, `ess_sd()`, `ess_tail()`, `mcse_mean()`, `mcse_quantile()`, `mcse_sd()`, `pareto_diags()`, `pareto_khat()`, `rhat()`, `rhat_basic()`, `rhat_nested()`

Examples

```
if (require("caret", quietly = TRUE) && require("randomForest", quietly = TRUE)) {
  x <- example_draws("eight_schools")
  print(rstar(x))
  print(rstar(x, split = FALSE))
  print(rstar(x, method = "gbm"))
  # can pass additional arguments to methods
  print(rstar(x, method = "gbm", verbose = FALSE))

  # with uncertainty, returns a vector of R* values
  hist(rstar(x, uncertainty = TRUE))
  hist(rstar(x, uncertainty = TRUE, nsimulations = 100))

  # can use other classification methods in caret library
  print(rstar(x, method = "knn"))
}
```

---

rvar *Random variables of arbitrary dimension*

Description

Random variables backed by arrays of arbitrary dimension

Usage

```
rvar(
  x = double(),
  dim = NULL,
  dimnames = NULL,
  nchains = NULL,
  with_chains = FALSE
)
```

Arguments

x	(multiple options) The object to convert to an rvar: • A vector of draws from a distribution. • An array where the first dimension represents draws from a distribution. The resulting rvar will have dimension dim(x)[-1]; that is, everything except the first dimension is used for the shape of the variable, and the first dimension is used to index draws from the distribution (see **Examples**). Optionally, if with_chains == TRUE, the first dimension indexes the iteration and the second dimension indexes the chain (see with_chains). • An rvar.

dim	(integer vector) One or more integers giving the maximal indices in each dimension to override the dimensions of the rvar to be created (see dim()). If NULL (the default), dim is determined by the input. **NOTE:** This argument controls the dimensions of the rvar, not the underlying array, so you cannot change the number of draws using this argument.
dimnames	(list) Character vectors giving the names in each dimension to override the names of the dimensions of the rvar to be created (see dimnames()). If NULL (the default), this is determined by the input. **NOTE:** This argument controls the names of the dimensions of the rvar, not the underlying array.
nchains	(positive integer) The number of chains. The if NULL (the default), 1 is used unless x is already an rvar, in which case the number of chains it has is used.
with_chains	(logical) Does x include a dimension for chains? If FALSE (the default), chains are not included, the first dimension of the input array should index draws, and the nchains argument can be used to determine the number of chains. If TRUE, the nchains argument is ignored and the second dimension of x is used to index chains. Internally, the array will be converted to a format without the chain index. Ignored when x is already an rvar.

Details

The "rvar" class internally represents random variables as arrays of arbitrary dimension, where the first dimension is used to index draws from the distribution. Most mathematical operators and functions are supported, including efficient matrix multiplication and vector and array-style indexing. The intent is that an rvar works as closely as possible to how a base vector/matrix/array does, with a few differences:

- The default behavior when subsetting is not to drop extra dimensions (i.e. the default drop argument for [is FALSE, not TRUE).
- Rather than base R-style recycling, rvars use a limited form of broadcasting: if an operation is being performed on two vectors with different size of the same dimension, the smaller vector will be recycled up to the size of the larger one along that dimension so long as it has size 1.

For functions that expect base numeric arrays and for which rvars cannot be used directly as arguments, you can use rfun() or rdo() to translate your code into code that executes across draws from one or more random variables and returns a random variable as output. Typically rdo() offers the most straightforward translation.

As rfun() and rdo() incur some performance cost, you can also operate directly on the underlying array using the draws_of() function. To re-use existing random number generator functions to efficiently create rvars, use rvar_rng().

Value

An object of class "rvar" representing a random variable.

See Also

as_rvar() to convert objects to rvars. See rdo(), rfun(), and rvar_rng() for higher-level interfaces for creating rvars.

Examples

```
set.seed(1234)

# To create a "scalar" `rvar`, pass a one-dimensional array or a vector
# whose length (here `4000`) is the desired number of draws:
x <- rvar(rnorm(4000, mean = 1, sd = 1))
x

# Create random vectors by adding an additional dimension:
n <- 4   # length of output vector
x <- rvar(array(rnorm(4000 * n, mean = rep(1:n, each = 4000), sd = 1), dim = c(4000, n)))
x

# Create a random matrix:
rows <- 4
cols <- 3
x <- rvar(array(rnorm(4000 * rows * cols, mean = 1, sd = 1), dim = c(4000, rows, cols)))
x

# If the input sample comes from multiple chains, we can indicate that using the
# nchains argument (here, 1000 draws each from 4 chains):
x <- rvar(rnorm(4000, mean = 1, sd = 1), nchains = 4)
x

# Or if the input sample has chain information as its second dimension, we can
# use with_chains to create the rvar
x <- rvar(array(rnorm(4000, mean = 1, sd = 1), dim = c(1000, 4)), with_chains = TRUE)
x
```

`rvar-dist` *Density, CDF, and quantile functions of random variables*

Description

The probability density function (`density()`), cumulative distribution function (`cdf()`), and quantile function / inverse CDF (`quantile()`) of an `rvar`.

Usage

```
## S3 method for class 'rvar'
density(x, at, ...)

## S3 method for class 'rvar_factor'
density(x, at, ...)

## S3 method for class 'rvar'
cdf(x, q, ...)
```

```
## S3 method for class 'rvar_factor'
cdf(x, q, ...)

## S3 method for class 'rvar_ordered'
cdf(x, q, ...)

## S3 method for class 'rvar'
quantile(x, probs, ...)

## S3 method for class 'rvar_factor'
quantile(x, probs, ...)

## S3 method for class 'rvar_ordered'
quantile(x, probs, ...)
```

Arguments

x	(rvar) An rvar object.
...	Additional arguments passed onto underlying methods: • For density(), these are passed to stats::density(). • For cdf(), these are ignored. • For quantile(), these are passed to stats::quantile().
q, at	(numeric vector) One or more quantiles.
probs	(numeric vector) One or more probabilities in [0,1].

Value

If x is a scalar rvar, returns a vector of the same length as the input (q, at, or probs) containing values from the corresponding function of the given rvar.

If x has length greater than 1, returns an array with dimensions c(length(y), dim(x)) where y is q, at, or probs, where each result[i,...] is the value of the corresponding function,f(y[i]), for the corresponding cell in the input array, x[...].

Examples

```
set.seed(1234)
x = rvar(rnorm(100))

density(x, seq(-2, 2, length.out = 10))
cdf(x, seq(-2, 2, length.out = 10))
quantile(x, ppoints(10))

x2 = c(rvar(rnorm(100, mean = -0.5)), rvar(rnorm(100, mean = 0.5)))
density(x2, seq(-2, 2, length.out = 10))
cdf(x2, seq(-2, 2, length.out = 10))
quantile(x2, ppoints(10))
```

`rvar-matmult` *Matrix multiplication of random variables*

Description

Matrix multiplication of random variables.

Usage

```
x %**% y

## S3 method for class 'rvar'
matrixOps(x, y)
```

Arguments

`x`	(multiple options) The object to be postmultiplied by `y`: • An `rvar` • A `numeric` vector or matrix • A `logical` vector or matrix If a vector is used, it is treated as a *row* vector.
`y`	(multiple options) The object to be premultiplied by `x`: • An `rvar` • A `numeric` vector or matrix • A `logical` vector or matrix If a vector is used, it is treated as a *column* vector.

Details

If `x` or `y` are vectors, they are converted into matrices prior to multiplication, with `x` converted to a row vector and `y` to a column vector. Numerics and logicals can be multiplied by `rvar`s and are broadcasted across all draws of the `rvar` argument. Tensor multiplication is used to efficiently multiply matrices across draws, so if either `x` or `y` is an `rvar`, `x %**% y` will be much faster than `rdo(x %*% y)`.

In R >= 4.3, you can also use `%*%` in place of `%**%` for matrix multiplication of `rvar`s. In R < 4.3, S3 classes cannot properly override `%*%`, so you must use `%**%` for matrix multiplication of `rvar`s.

Value

An `rvar` representing the matrix product of `x` and `y`.

Examples

```
# d has mu (mean vector of length 3) and Sigma (3x3 covariance matrix)
d <- as_draws_rvars(example_draws("multi_normal"))
d$Sigma

# trivial example: multiplication by a non-random matrix
d$Sigma %**% diag(1:3)

# Decompose Sigma into R s.t. R'R = Sigma ...
R <- chol(d$Sigma)
# ... and recreate Sigma using matrix multiplication
t(R) %**% R
```

`rvar-slice` *Random variable slicing*

Description

Operations for slicing `rvar`s and replacing parts of `rvar`s.

Usage

```
## S3 method for class 'rvar'
x[[i, ...]]

## S3 replacement method for class 'rvar'
x[[i, ...]] <- value

## S3 method for class 'rvar'
x[..., drop = FALSE]

## S3 replacement method for class 'rvar'
x[i, ...] <- value
```

Arguments

`x`	an `rvar`.
`i, ...`	indices; see *Details*.
`value`	(`rvar` or coercible to `rvar`) Value to insert into `x` at the location determined by the indices.
`drop`	(logical) Should singular dimensions be dropped when slicing array `rvar`s? Unlike base array slicing operations, defaults to `FALSE`.

Details

The rvar slicing operators ([and [[) attempt to implement the same semantics as the base array slicing operators. There are some exceptions; most notably, rvar slicing defaults to drop = FALSE instead of drop = TRUE.

Extracting or replacing single elements with [[

The [[operator extracts (or replaces) single elements. It always returns (or replaces) a scalar (length-1) rvar.

The x[[i, ...]] operator can be used as follows:

- x[[<numeric>]] for scalar numeric i: gives the ith element of x. If x is multidimensional (i.e. length(dim(x)) > 1), extra dimensions are ignored when indexing. For example, if x is a 6×2 rvar array, the 7th element, x[[7]], will be the first element of the second column, x[1, 2].
- x[[<numeric rvar>]] for scalar numeric rvar i: a generalization of indexing when i is a scalar numeric. Within each draw of x, selects the element corresponding to the value of i within that same draw.
- x[[<character>]] for scalar character i: gives the element of x with name equal to i. **Unlike with base arrays**, does not work with multidimensional rvars.
- x[[i_1, i_2, ..., i_n]] for scalar numeric or character i_1, i_2, etc. Must provide exactly the same number of indices as dimensions in x. Selects the element at the corresponding position in the rvar by number and/or dimname (as a string).

Extracting or replacing multiple elements with [

The [operator extracts (or replaces) multiple elements. It always returns (or replaces) a possibly-multidimensional rvar.

The x[i, ...] operator can be used as follows:

- x[<logical>] for vector logical i: i is recycled to the same length as x, ignoring multiple dimensions in x, then an rvar vector is returned containing the elements in x where i is TRUE.
- x[<logical rvar>] for scalar logical rvar i: returns an rvar the same shape as x containing only those draws where i is TRUE.
- x[<numeric>] for vector numeric i: an rvar vector is returned containing the ith elements of x, ignoring dimensions.
- x[<matrix>] for numeric matrix i, where ncol(i) == length(dim(x)): each row of i should give the multidimensional index for a single element in x. The result is an rvar vector of length nrow(i) containing elements of x selected by each row of i.
- x[i_1, i_2, ..., i_n] for vector numeric, character, or logical i_1, i_2, etc. Returns a slice of x containing all elements from the dimensions specified in i_1, i_2, etc. If an argument is left empty, all elements from that dimension are included. Unlike base arrays, trailing dimensions can be omitted entirely and will still be selected; for example, if x has three dimensions, both x[1, ,] and x[1,] can be used to create a slice that includes all elements from the last two dimensions. Unlike base arrays, [defaults to drop = FALSE, so results retain the same number of dimensions as x.

Examples

```
x <- rvar(array(1:24, dim = c(4,2,3)))
dimnames(x) <- list(c("a","b"), c("d","e","f"))
x

## Slicing single elements
# x[[<numeric>]]
x[[2]]

# x[[<numeric rvar>]]
# notice the draws of x[1:4]...
draws_of(x[1:4])
x[[rvar(c(1,3,4,4))]]
# ... x[[rvar(c(1,3,4,4))]] creates a mixures of those draws
draws_of(x[[rvar(c(1,3,4,4))]])

# x[[i_1,i_2,...]]
x[[2,"e"]]


## Slicing multiple elements
# x[<logical>]
x[c(TRUE,TRUE,FALSE)]

# x[<logical rvar>]
# select every other draw
x[rvar(c(TRUE,FALSE,TRUE,FALSE))]

# x[<numeric>]
x[1:3]

# x[<matrix>]
x[rbind(
  c(1,2),
  c(1,3),
  c(2,2)
)]

# x[i_1,i_2,...,i_n]
x[1,]
x[1,2:3]
x[,2:3]
```

`rvar-summaries-over-draws`

Summaries of random variables within array elements, over draws

---

Description

Compute summaries within elements of an `rvar` and over draws of each element, producing an array of the same shape as the input random variable (except in the case of `range()`, see **Details**).

Usage

```
E(x, ...)

## S3 method for class 'rvar'
mean(x, ...)

Pr(x, ...)

## Default S3 method:
Pr(x, ...)

## S3 method for class 'logical'
Pr(x, ...)

## S3 method for class 'rvar'
Pr(x, ...)

## S3 method for class 'rvar'
median(x, ...)

## S3 method for class 'rvar'
min(x, ...)

## S3 method for class 'rvar'
max(x, ...)

## S3 method for class 'rvar'
sum(x, ...)

## S3 method for class 'rvar'
prod(x, ...)

## S3 method for class 'rvar'
all(x, ...)

## S3 method for class 'rvar'
any(x, ...)

## S3 method for class 'rvar'
Summary(...)

## S3 method for class 'rvar'
variance(x, ...)

var(x, ...)

## Default S3 method:
var(x, ...)
```

```
## S3 method for class 'rvar'
var(x, ...)

sd(x, ...)

## Default S3 method:
sd(x, ...)

## S3 method for class 'rvar'
sd(x, ...)

mad(x, ...)

## Default S3 method:
mad(x, ...)

## S3 method for class 'rvar'
mad(x, ...)

## S3 method for class 'rvar_ordered'
mad(x, ...)

## S3 method for class 'rvar'
range(x, ...)

## S3 method for class 'rvar'
is.finite(x)

## S3 method for class 'rvar'
is.infinite(x)

## S3 method for class 'rvar'
is.nan(x)

## S3 method for class 'rvar'
is.na(x)
```

Arguments

x	(rvar) An rvar.
...	Further arguments passed to underlying functions (e.g., base::mean() or base::median()), such as na.rm.

Details

Summaries include expectations (E() or mean()), probabilities (Pr()), medians (median()), spread (var(), variance(), sd(), mad()), sums and products (sum(), prod()), extrema and ranges

(`min()`, `max()`, `range()`), logical summaries (`all()`, `any()`), and special value predicates (`is.finite()`, `is.infinite()`, `is.nan()`, `is.na()`).

Unless otherwise stated, these functions return a numeric array with the same shape (same dimensions) as the input `rvar`, `x`.

`range(x)` returns an array with dimensions `c(2, dim(x))`, where the last dimension contains the minimum and maximum values.

`is.infinite(x)`, `is.nan(x)`, and `is.na(x)` return logical arrays, where each element is `TRUE` if **any** draws in its corresponding element in `x` match the predicate. Each elements in the result of `is.finite(x)` is `TRUE` if **all** draws in the corresponding element in `x` are finite.

Both `E()`, `mean()`, and `Pr()` return the means of each element in the input. `Pr()` additionally checks that the provided `rvar` is a logical variable (hence, taking its expectation results in a probability).

For consistency, `E()` and `Pr()` are also defined for base arrays so that they can be used as summary functions in `summarise_draws()`.

Value

A numeric or logical vector with the same dimensions as the given random variable, where each entry in the vector is the mean, median, or variance of the corresponding entry in `x`.

See Also

rvar-summaries-within-draws for summary functions within draws. rvar-dist for density, CDF, and quantile functions of random variables.

Other rvar-summaries: `rvar-summaries-within-draws`, `rvar_is_finite()`

Examples

```
set.seed(5678)
x = rvar_rng(rnorm, 4, mean = 1:4, sd = 2)

# These should all be ~= c(1, 2, 3, 4)
E(x)
mean(x)
median(x)

# This ...
Pr(x < 1.5)
# ... should be about the same as this:
pnorm(1.5, mean = 1:4, sd = 2)
```

`rvar-summaries-within-draws` *Summaries of random variables over array elements, within draws*

Description

Compute summaries of random variables over array elements and within draws, producing a new random variable of length 1 (except in the case of `rvar_range()`, see **Details**).

Usage

```
rvar_mean(..., na.rm = FALSE)

rvar_median(..., na.rm = FALSE)

rvar_sum(..., na.rm = FALSE)

rvar_prod(..., na.rm = FALSE)

rvar_min(..., na.rm = FALSE)

rvar_max(..., na.rm = FALSE)

rvar_sd(..., na.rm = FALSE)

rvar_var(..., na.rm = FALSE)

rvar_mad(..., constant = 1.4826, na.rm = FALSE)

rvar_range(..., na.rm = FALSE)

rvar_quantile(..., probs, names = FALSE, na.rm = FALSE)

rvar_all(..., na.rm = FALSE)

rvar_any(..., na.rm = FALSE)
```

Arguments

`...`	(rvar) One or more rvars.
`na.rm`	(logical) Should NAs be removed from the input before summaries are computed? The default is FALSE.
`constant`	(scalar real) For `rvar_mad()`, a scale factor for computing the median absolute deviation. See the details of `stats::mad()` for the justification for the default value.
`probs`	(numeric vector) For `rvar_quantile()`, probabilities in `[0, 1]`.
`names`	(logical) For `rvar_quantile()`, if TRUE, the result has a `names` attribute.

Details

These functions compute statistics within each draw of the random variable. For summaries over draws (such as expectations), see rvar-summaries-over-draws.

Each function defined here corresponds to the base function of the same name without the `rvar_` prefix (e.g., `rvar_mean()` calls `mean()` under the hood, etc).

Value

An `rvar` of length 1 (for `range()`, length 2; for `quantile()`, length equal to `length(probs)`) with the same number of draws as the input rvar(s) containing the summary statistic computed within each draw of the input rvar(s).

See Also

rvar-summaries-over-draws for summary functions across draws (e.g. expectations). rvar-dist for density, CDF, and quantile functions of random variables.

Other rvar-summaries: `rvar-summaries-over-draws`, `rvar_is_finite()`

Examples

```
set.seed(5678)
x = rvar_rng(rnorm, 4, mean = 1:4, sd = 2)

# These will give similar results to mean(1:4),
# median(1:4), sum(1:4), prod(1:4), etc
rvar_mean(x)
rvar_median(x)
rvar_sum(x)
rvar_prod(x)
rvar_range(x)
rvar_quantile(x, probs = c(0.25, 0.5, 0.75), names = TRUE)
```

`rvar_apply` *Random variable resulting from a function applied over margins of an array or random variable*

Description

Returns an `rvar` obtained by applying a function to margins of an array or `rvar`. Acts like `apply()`, except that the function supplied (`.f`) should return an `rvar`, and the final result is always an `rvar`.

Usage

```
rvar_apply(.x, .margin, .f, ...)
```

Arguments

`.x`	An array or an `rvar`.
`.margin`	(multiple options) The subscripts which the function will be applied over: • An integer vector. E.g., for a matrix `1` indicates rows, `2` indicates columns, `c(1, 2)` indicates rows and columns. • A character vector of dimension names if `.x` has named dimensions.
`.f`	(function) The function to be applied. The function `.f` must return an `rvar` and the dimensions of the result of `.f` applied to each margin of `.x` must be able to be broadcasted to a common shape (otherwise the resulting `rvar` cannot be simplified). See **Details**.
`...`	Optional arguments passed to `.f`.

Details

This function acts much like `apply()`, except that the function passed to it (`.f`) must return `rvar`s, and the result is simplified into an `rvar`. Unlike `apply()`, it also keeps the dimensions of the returned values along each margin, rather than simplifying each margin to a vector, and if the results of `.f` do not all have the same dimensions, it applies the `rvar` broadcasting rules to bind results together rather than using vector recycling.

If you wish to apply functions over `rvar`s where the result is not intended to be simplified into an `rvar`, you can use the standard `apply()`, `lapply()`, `sapply()`, or `vapply()` functions.

Value

An `rvar`.

If the result of each call to `.f` returns an `rvar` of dimension `d` after being broadcast to a common shape, then `rvar_apply()` returns an `rvar` of dimension `c(d, dim(.x)[.margin])`. If the last dimension of the result would be `1`, it is dropped (other dimensions equal to `1` are retained). If `d` is `0`, the result has length `0` but not necessarily the 'correct' dimension.

See Also

`as_rvar()` to convert objects to `rvar`s. See `rdo()`, `rfun()`, and `rvar_rng()` for higher-level interfaces for creating `rvar`s.

Examples

```
set.seed(3456)
x <- rvar_rng(rnorm, 24, mean = 1:24)
dim(x) <- c(2,3,4)

# we can find the distributions of marginal means of the above array
# using rvar_mean along with rvar_apply
rvar_apply(x, 1, rvar_mean)
rvar_apply(x, 2:3, rvar_mean)
```

`rvar_factor` *Factor random variables of arbitrary dimension*

Description

Random variables backed by factor-like arrays of arbitrary dimension.

Usage

```
rvar_factor(
  x = factor(),
  dim = NULL,
  dimnames = NULL,
  nchains = NULL,
  with_chains = FALSE,
  ...
)

rvar_ordered(
  x = ordered(NULL),
  dim = NULL,
  dimnames = NULL,
  nchains = NULL,
  with_chains = FALSE,
  ...
)
```

Arguments

`x`	(multiple options) The object to convert to an `rvar`: • A vector of draws from a distribution. • An array where the first dimension represents draws from a distribution. The resulting `rvar` will have dimension `dim(x)[-1]`; that is, everything except the first dimension is used for the shape of the variable, and the first dimension is used to index draws from the distribution (see **Examples**). Optionally, if `with_chains == TRUE`, the first dimension indexes the iteration and the second dimension indexes the chain (see `with_chains`). • An `rvar`.
`dim`	(integer vector) One or more integers giving the maximal indices in each dimension to override the dimensions of the `rvar` to be created (see `dim()`). If `NULL` (the default), `dim` is determined by the input. **NOTE:** This argument controls the dimensions of the `rvar`, not the underlying array, so you cannot change the number of draws using this argument.
`dimnames`	(list) Character vectors giving the names in each dimension to override the names of the dimensions of the `rvar` to be created (see `dimnames()`). If `NULL` (the default), this is determined by the input. **NOTE:** This argument controls the names of the dimensions of the `rvar`, not the underlying array.

nchains (positive integer) The number of chains. The if NULL (the default), 1 is used unless x is already an rvar, in which case the number of chains it has is used.

with_chains (logical) Does x include a dimension for chains? If FALSE (the default), chains are not included, the first dimension of the input array should index draws, and the nchains argument can be used to determine the number of chains. If TRUE, the nchains argument is ignored and the second dimension of x is used to index chains. Internally, the array will be converted to a format without the chain index. Ignored when x is already an rvar.

... Arguments passed on to base::factor

levels an optional vector of the unique values (as character strings) that x might have taken. The default is the unique set of values taken by as.character(x), sorted into increasing order *of* x. Note that this set can be specified as smaller than sort(unique(x)).

labels *either* an optional character vector of labels for the levels (in the same order as levels after removing those in exclude), *or* a character string of length 1. Duplicated values in labels can be used to map different values of x to the same factor level.

exclude a vector of values to be excluded when forming the set of levels. This may be factor with the same level set as x or should be a character.

ordered logical flag to determine if the levels should be regarded as ordered (in the order given).

nmax an upper bound on the number of levels; see 'Details'.

Details

A subtype of rvar() that represents a (possibly multidimensional) sample of a factor or an ordered factor. It is otherwise very similar to the basic rvar(): it is backed by a multidimensional array with draws as the first dimension. The primary difference is that the backing array has class "factor" (for rvar_factor()) or c("ordered", "factor") (for rvar_ordered()). If you pass a factor or ordered factor to rvar() it will automatically return an object with the classes "rvar_factor" or c("rvar_ordered", "rvar_factor").

See rvar() for more details on the internals of the random variable datatype.

Value

An object of class "rvar_factor" representing a factor-like random variable.

See Also

as_rvar_factor() to convert objects to rvar_factors. See rdo(), rfun(), and rvar_rng() for higher-level interfaces for creating rvars.

Examples

```
set.seed(1234)

# To create a "scalar" `rvar_factor`, pass a one-dimensional array or a vector
# whose length (here `4000`) is the desired number of draws:
```

```
x <- rvar(sample(c("a","a","a","b","c"), 4000, replace = TRUE))
x

# Create random vectors by adding an additional dimension:
x_array <- array(c(
    sample(c("a","a","a","b","c"), 4000, replace = TRUE),
    sample(c("a","a","b","c","c"), 4000, replace = TRUE),
    sample(c("b","b","b","b","c"), 4000, replace = TRUE),
    sample(c("d","d","b","b","c"), 4000, replace = TRUE)
  ), dim = c(4000, 4))
rvar_factor(x_array)

# You can also create ordered factors
rvar_ordered(x_array)

# arguments of factor() and ordered() are passed down by the constructor
# e.g. we can reorder levels of an ordered factor:
rvar_ordered(x_array, levels = c("d","c","b","a"))

# Unlike base factors, rvar factors can be matrices or arrays:
rvar_factor(x_array, dim = c(2, 2))

# If the input to rvar_factor() is an array with a `"levels"` attribute, it
# will use those as the levels of the factor
y_array <- t(array(rbinom(3000, 1, c(0.1, 0.5, 0.9)) + 1, dim = c(3, 1000)))
rvar(y_array)
# with levels
attr(y_array, "levels") = c("a", "b")
rvar_factor(y_array)
```

`rvar_ifelse` *Random variable ifelse*

Description

A version of `ifelse()` that returns an `rvar`.

Usage

```
rvar_ifelse(test, yes, no)
```

Arguments

`test`	(logical `rvar`, or castable to one) logical test determining whether the value in `yes` or `no` is assigned in the corresponding position of the result.
`yes`	(`rvar`, or castable to one) corresponding values assigned for entries in `test` that are `TRUE`.
`no`	(`rvar`, or castable to one) corresponding values assigned for entries in `test` that are `FALSE`.

Value

An rvar with the common type of yes and no (as determined by vctrs::vec_cast_common()) and a shape determined by broadcasting test, yes, and no to a common shape (see the section on broadcasting rules in vignette("rvar")). For every element of draws_of(test), the corresponding element of draws_of(yes) or draws_of(no) is placed into the result, depending on whether the element of test is TRUE or FALSE.

Examples

```
x <- rvar(1:4)
y <- rvar(5:8)

i <- rvar(c(TRUE,FALSE,TRUE,FALSE))
z <- rvar_ifelse(i, x, y)
z
draws_of(z)
```

---

rvar_is_finite *Special value predicates for random variables*

Description

Compute special value predicates (checking for finite / infinite values, NaN, and NA) on all draws within a random variable, returning a random variable.

Usage

```
rvar_is_finite(x)

rvar_is_infinite(x)

rvar_is_nan(x)

rvar_is_na(x)
```

Arguments

x	(rvar) An rvar.

Details

These functions return a new rvar that is the result of applying is.finite(), is.infinite(), is.nan(), or is.na() to every draw in the input random variable.

Value

A logical rvar of the same length as the input.

See Also

rvar-summaries-over-draws for summary functions across draws, including implementations of `is.finite()`, `is.infinite()`, `is.nan()`, and `is.na()` for rvars.

Other rvar-summaries: `rvar-summaries-over-draws`, `rvar-summaries-within-draws`

Examples

```
x <- rvar(c(1, Inf, -Inf, NaN, NA))
x

rvar_is_finite(x)
rvar_is_infinite(x)
rvar_is_nan(x)
rvar_is_na(x)
```

`rvar_rng` *Create random variables from existing random number generators*

Description

Specialized alternative to `rdo()` or `rfun()` for creating rvars from existing random-number generator functions (such as `rnorm()`, `rbinom()`, etc).

Usage

```
rvar_rng(.f, n, ..., ndraws = NULL)
```

Arguments

`.f`	(function) A function (or string naming a function) representing a random-number generating function that follows the pattern of base random number generators (like `rnorm()`, `rbinom()`, etc). It must: • Have a first argument, `n`, giving the number of draws to take from the distribution • Have vectorized parameter arguments • Return a single vector of length `n`
`n`	(positive integer) The length of the output rvar vector (**not** the number of draws).
`...`	Arguments passed to `.f`. These arguments may include rvars, so long as they are vectors only (no multidimensional rvars are allowed).
`ndraws`	(positive integer) The number of draws used to construct the returned random variable if no rvars are supplied in `...`. If `NULL`, `getOption("posterior.rvar_ndraws")` is used (default 4000). If `...` contains rvars, the number of draws in the provided rvars is used instead of the value of this argument.

Details

This function unwraps the arrays underlying the input rvars in ... and then passes them to .f, relying on the vectorization of .f to evaluate it across draws from the input rvars. This is why the arguments of .f **must** be vectorized. It asks for n times the number of draws in the input rvars (or ndraws if none are given) draws from the random number generator .f, then reshapes the output from .f into an rvar with length n.

rvar_rng() is a fast alternative to rdo() or rfun(), but you **must** ensure that .f satisfies the preconditions described above for the result to be correct. Most base random number generators satisfy these conditions. It is advisable to test against rdo() or rfun() (which should be correct, but slower) if you are uncertain.

Value

A single-dimensional rvar of length n.

See Also

Other rfun: rdo(), rfun()

Examples

```
mu <- rvar_rng(rnorm, 10, mean = 1:10, sd = 1)
sigma <- rvar_rng(rgamma, 1, shape = 1, rate = 1)
x <- rvar_rng(rnorm, 10, mu, sigma)
x
```

---

split_chains *Split Chains*

Description

Split chains by halving the number of iterations per chain and doubling the number of chains.

Usage

```
split_chains(x, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).

Value

A draws object of the same class as x.

Examples

```
x <- example_draws()
niterations(x)
nchains(x)

x <- split_chains(x)
niterations(x)
nchains(x)
```

---

subset_draws *Subset* draws *objects*

Description

Subset draws objects by variables, iterations, chains, and draws indices.

Usage

```
subset_draws(x, ...)

## S3 method for class 'draws_matrix'
subset_draws(
  x,
  variable = NULL,
  iteration = NULL,
  chain = NULL,
  draw = NULL,
  regex = FALSE,
  unique = TRUE,
  exclude = FALSE,
  scalar = FALSE,
  ...
)

## S3 method for class 'draws_array'
subset_draws(
  x,
  variable = NULL,
  iteration = NULL,
  chain = NULL,
  draw = NULL,
  regex = FALSE,
  unique = TRUE,
  exclude = FALSE,
  scalar = FALSE,
  ...
```

```
)

## S3 method for class 'draws_df'
subset_draws(
  x,
  variable = NULL,
  iteration = NULL,
  chain = NULL,
  draw = NULL,
  regex = FALSE,
  unique = TRUE,
  exclude = FALSE,
  scalar = FALSE,
  ...
)

## S3 method for class 'draws_list'
subset_draws(
  x,
  variable = NULL,
  iteration = NULL,
  chain = NULL,
  draw = NULL,
  regex = FALSE,
  unique = TRUE,
  exclude = FALSE,
  scalar = FALSE,
  ...
)

## S3 method for class 'draws_rvars'
subset_draws(
  x,
  variable = NULL,
  iteration = NULL,
  chain = NULL,
  draw = NULL,
  regex = FALSE,
  unique = TRUE,
  exclude = FALSE,
  scalar = FALSE,
  ...
)

## S3 method for class 'rvar'
subset_draws(x, variable = NULL, ...)

## S3 method for class 'draws'
```

```
subset(x, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).
variable	(character vector) The variables to select. All elements of non-scalar variables can be selected at once.
iteration	(integer vector) The iteration indices to select.
chain	(integer vector) The chain indices to select.
draw	(integer vector) The draw indices to be select. Subsetting draw indices will lead to an automatic merging of chains via merge_chains. By default, a message will be issued when this happens but you can deactivate it via options(posterior.warn_on_merge_chai = FALSE).
regex	(logical) Should variable should be treated as a (vector of) regular expressions? Any variable in x matching at least one of the regular expressions will be selected. Defaults to FALSE.
unique	(logical) Should duplicated selection of chains, iterations, or draws be allowed? If TRUE (the default) only unique chains, iterations, and draws are selected regardless of how often they appear in the respective selecting arguments.
exclude	(logical) Should the selected subset be excluded? If FALSE (the default) only the selected subset will be returned. If TRUE everything but the selected subset will be returned.
scalar	(logical) Should only scalar variables be selected? If FALSE (the default), all variables with matching names and *arbitrary* indices will be selected (see examples).

Details

To ensure that multiple consecutive subsetting operations work correctly, subset() *repairs* the draws object before and after subsetting.

Value

A draws object of the same class as x.

Examples

```
x <- example_draws()
subset_draws(x, variable = c("mu", "tau"))
subset_draws(x, chain = 2)
subset_draws(x, iteration = 5:10, chain = 3:4)

# extract the first chain twice
subset_draws(x, chain = c(1, 1), unique = FALSE)

# extract all elements of 'theta'
```

```
subset_draws(x, variable = "theta")

# trying to extract only a scalar 'theta' will fail
# subset_draws(x, variable = "theta", scalar = TRUE)
```

---

thin_draws *Thin* draws *objects*

Description

Thin draws objects to reduce their size and autocorrelation in the chains.

Usage

```
thin_draws(x, thin = NULL, ...)

## S3 method for class 'draws'
thin_draws(x, thin = NULL, ...)

## S3 method for class 'rvar'
thin_draws(x, thin = NULL, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
thin	(positive numeric) The period for selecting draws. Must be between 1 and the number of iterations. If the value is not an integer, the draws will be selected such that the number of draws returned is equal to round(ndraws(x) / thin). Intervals between selected draws will be either ceiling(thin) or floor(thin), such that the average interval will be close to the thin value. If NULL, it will be automatically calculated based on bulk and tail effective sample size as suggested by Säilynoja et al. (2022).
...	Arguments passed to individual methods (if applicable).

Value

A draws object of the same class as x.

References

Teemu Säilynoja, Paul-Christian Bürkner, and Aki Vehtari (2022). Graphical test for discrete uniformity and its applications in goodness-of-fit evaluation and multiple sample comparison. *Statistics and Computing*. 32, 32. doi:10.1007/s11222-022-10090-6

Examples

```
x <- example_draws()
niterations(x)

x <- thin_draws(x, thin = 5)
niterations(x)
```

`uniformity_test` *Uniformity test for PIT values*

Description

Tests whether PIT (probability integral transform) values deviate from uniformity while accounting for dependence among them. LOO-PIT values are not independent because each LOO predictive distribution is conditioned on nearly the same data (all but one observation), which induces correlation. Standard uniformity tests that assume independence can have inflated Type I error or reduced power. This function implements dependence-aware procedures from Tesso and Vehtari (2026).

Usage

```
uniformity_test(pit, test)
```

Arguments

`pit`	Numeric vector of PIT values in `[0, 1]`.
`test`	Character string. One of `"POT"`, `"PIET"`, or `"PRIT"`. See details above.

Details

Three test variants are available:

- **POT** (Pointwise Order Tests): Uses beta distributions for order statistics. Best for continuous PIT values; recommended as default for LOO-PIT model checking. Good power against diverse departures from uniformity.
- **PIET** (Pointwise Inverse-CDF Evaluation Tests): Uses a continuous reference distribution (exponential) via inverse-CDF transformation. Best for detecting tail deviations; maintains valid Type I error control under strong dependence.
- **PRIT** (Pointwise Rank-based Individual Tests): Uses binomial distributions on scaled ECDF (ranks). Intended for discrete or rank-based PIT values.

All procedures compute pointwise p-values, aggregate them via the (truncated) Cauchy combination test (Liu and Xie, 2020; Chen et al., 2025), and derive Shapley values to quantify each point's contribution to the overall test.

Value

A list with components:

- `pvalue`: Global p-value from the Cauchy combination test.
- `pointwise`: Shapley values (contributions of each PIT value to the test statistic). Non-negative values indicate points contributing to evidence against uniformity.

References

Tesso, H., and Vehtari, A. (2026). LOO-PIT predictive model checking. arXiv preprint arXiv:2603.02928.

Liu, Y., and Xie, J. (2020). Cauchy combination test: a powerful test with analytic p-value calculation under arbitrary dependency structures. *Journal of the American Statistical Association*, 115(529), 393-402.

Chen, B., Xu, W., and Gao, X. (2025). Truncated Cauchy combination test: a robust and powerful p-value combination method with arbitrary correlations. arXiv preprint arXiv:2506.12489.

`variables` *Get variable names from* `draws` *objects*

Description

Get variable names from `draws` objects.

Usage

```
variables(x, ...)

## S3 method for class 'draws_matrix'
variables(x, reserved = FALSE, with_indices = TRUE, ...)

## S3 method for class 'draws_array'
variables(x, reserved = FALSE, with_indices = TRUE, ...)

## S3 method for class 'draws_df'
variables(x, reserved = FALSE, with_indices = TRUE, ...)

## S3 method for class 'draws_list'
variables(x, reserved = FALSE, with_indices = TRUE, ...)

## S3 method for class 'draws_rvars'
variables(x, reserved = FALSE, with_indices = FALSE, ...)

nvariables(x, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).
reserved	(logical) Should reserved variables be included in the output? Defaults to FALSE. See reserved_variables for an overview of currently reserved variable names.
with_indices	(logical) Should indices be included in variable names? For example, if the object includes variables named "x[1]" and "x[2]", if TRUE, c("x[1]", "x[2]") is returned; if FALSE, only "x" is returned. Defaults to TRUE for all formats except draws_rvars().

Details

variables() returns a vector of all variable names, and nvariables() returns the number of variables.

Value

For variables(), a character vector.

For nvariables(), a scalar integer.

See Also

variables<-, rename_variables, draws-index

Examples

```
x <- example_draws()

variables(x)
nvariables(x)
variables(x) <- letters[1:nvariables(x)]
```

---

variables<- *Set variable names in* draws *objects*

Description

Set variable names for all variables in a draws object. The set_variables() form is useful when using pipe operators.

Usage

```
variables(x, ...) <- value

## S3 replacement method for class 'draws_matrix'
variables(x, with_indices = TRUE, ...) <- value

## S3 replacement method for class 'draws_array'
variables(x, with_indices = TRUE, ...) <- value

## S3 replacement method for class 'draws_df'
variables(x, with_indices = TRUE, ...) <- value

## S3 replacement method for class 'draws_list'
variables(x, with_indices = TRUE, ...) <- value

## S3 replacement method for class 'draws_rvars'
variables(x, with_indices = FALSE, ...) <- value

set_variables(x, variables, ...)
```

Arguments

x	(draws) A draws object or another R object for which the method is defined.
...	Arguments passed to individual methods (if applicable).
value, variables	(character vector) new variable names.
with_indices	(logical) Should indices be included in variable names? For example, if the object includes variables named "x[1]" and "x[2]", if TRUE, c("x[1]", "x[2]") is returned; if FALSE, only "x" is returned. Defaults to TRUE for all formats except draws_rvars().

Details

variables(x) <- value allows you to modify the vector of variable names, similar to how names(x) <- value works for vectors and lists. For renaming specific variables, set_variables(x, value) works equivalently, but is more intuitive when using the pipe operator.

For renaming specific variables, rename_variables() may offer a more convenient approach.

Value

Returns a draws object of the same format as x, with variables named as specified.

See Also

variables, rename_variables, draws-index

Examples

```
x <- example_draws()

variables(x)
nvariables(x)
variables(x) <- letters[1:nvariables(x)]

# or equivalently...
x <- set_variables(x, letters[1:nvariables(x)])
```

`weight_draws` *Weight* `draws` *objects*

Description

Add weights to `draws` objects, with one weight per draw, for use in subsequent weighting operations. For reasons of numerical accuracy, weights are stored in the form of unnormalized log-weights (in a variable called `.log_weight`). See `weights.draws()` for details how to extract weights from `draws` objects.

Usage

```
weight_draws(x, weights, ...)

## S3 method for class 'draws_matrix'
weight_draws(x, weights, log = FALSE, pareto_smooth = FALSE, ...)

## S3 method for class 'draws_array'
weight_draws(x, weights, log = FALSE, pareto_smooth = FALSE, ...)

## S3 method for class 'draws_df'
weight_draws(x, weights, log = FALSE, pareto_smooth = FALSE, ...)

## S3 method for class 'draws_list'
weight_draws(x, weights, log = FALSE, pareto_smooth = FALSE, ...)

## S3 method for class 'draws_rvars'
weight_draws(x, weights, log = FALSE, pareto_smooth = FALSE, ...)
```

Arguments

`x`	(draws) A `draws` object or another R object for which the method is defined.
`weights`	(numeric vector) A vector of weights of length `ndraws(x)`. Weights will be internally stored on the log scale (in a variable called `.log_weight`) and will not be normalized, but normalized (non-log) weights can be returned via the `weights.draws()` method later.

...	Arguments passed to individual methods (if applicable).
log	(logical) Are the weights passed already on the log scale? The default is FALSE, that is, expecting weights to be on the standard (non-log) scale.
pareto_smooth	(logical) Should the weights be Pareto-smoothed? The default is FALSE.

Value

A draws object of the same class as x.

See Also

weights.draws(), resample_draws()

Examples

```
x <- example_draws()

# sample some random weights for illustration
wts <- rexp(ndraws(x))
head(wts)

# add weights
x <- weight_draws(x, weights = wts)

# extract weights
head(weights(x)) # defaults to normalized weights
head(weights(x, normalize=FALSE)) # recover original weights
head(weights(x, log=TRUE)) # get normalized log-weights

# add weights which are already on the log scale
log_wts <- log(wts)
head(log_wts)

x <- weight_draws(x, weights = log_wts, log = TRUE)
# extract weights
head(weights(x))
head(weights(x, log=TRUE, normalize = FALSE)) # recover original log_wts

# add weights on log scale and Pareto smooth them
x <- weight_draws(x, weights = log_wts, log = TRUE, pareto_smooth = TRUE)
```

---

weights.draws *Extract Weights from Draws Objects*

Description

Extract weights from draws objects, with one weight per draw. See weight_draws for details how to add weights to draws objects.

Usage

```
## S3 method for class 'draws'
weights(object, log = FALSE, normalize = TRUE, ...)
```

Arguments

object	(draws) A draws object.
log	(logical) Should the weights be returned on the log scale? Defaults to FALSE.
normalize	(logical) Should the weights be normalized to sum to 1 on the standard scale? Defaults to TRUE.
...	Arguments passed to individual methods (if applicable).

Value

A vector of weights, with one weight per draw.

See Also

weight_draws, resample_draws

Examples

```
x <- example_draws()

# sample some random weights for illustration
wts <- rexp(ndraws(x))
head(wts)

# add weights
x <- weight_draws(x, weights = wts)

# extract weights
head(weights(x)) # defaults to normalized weights
head(weights(x, normalize=FALSE)) # recover original weights
head(weights(x, log=TRUE)) # get normalized log-weights

# add weights which are already on the log scale
log_wts <- log(wts)
head(log_wts)

x <- weight_draws(x, weights = log_wts, log = TRUE)
# extract weights
head(weights(x))
head(weights(x, log=TRUE, normalize = FALSE)) # recover original log_wts

# add weights on log scale and Pareto smooth them
x <- weight_draws(x, weights = log_wts, log = TRUE, pareto_smooth = TRUE)
```

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