

# Package ‘BayesTools’

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**Title** Tools for Bayesian Analyses

**Version** 0.2.11

**Description** Provides tools for conducting Bayesian analyses. The package contains functions for creating a wide range of prior distribution objects, mixing posterior samples from 'JAGS' and 'Stan' models, plotting posterior distributions, and etc... The tools for working with prior distribution span from visualization, generating 'JAGS' and 'bridgesampling' syntax to basic functions such as `rng`, `quantile`, and distribution functions.

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**URL** <https://fbartos.github.io/BayesTools/>

**BugReports** <https://github.com/FBartos/BayesTools/issues>

**License** GPL-3

**Encoding** UTF-8

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rmarkdown

**RdMacros** Rdpack

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<code>add_column</code>	<i>Adds column to BayesTools table</i>
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### Description

Adds column to a BayesTools table while not breaking formatting, attributes, etc...

### Usage

```
add_column(  
  table,  
  column_title,  
  column_values,  
  column_position = NULL,  
  column_type = NULL  
)
```

### Arguments

<code>table</code>	BayesTools table
<code>column_title</code>	title of the new column
<code>column_values</code>	values of the new column
<code>column_position</code>	position of the new column (defaults to NULL which appends the column to the end)
<code>column_type</code>	type of values of the new column table (important for formatting, defaults to NULL = the function tries to guess numeric / character based on the <code>column_values</code> but many more specific types are available)

### Value

returns an object of 'BayesTools\_table' class.

BayesTools

*BayesTools*

---

**Description**

BayesTools: Provides tools for conducting Bayesian analyses. The package contains functions for creating a wide range of prior distribution objects, mixing posterior samples from JAGS and Stan models, plotting posterior distributions, and etc... The tools for working with prior distribution span from visualization, generating JAGS and bridgesampling syntax to basic functions such as rng, quantile, and distribution functions.

**Author(s)**

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**See Also**

Useful links:

- <https://fbartos.github.io/BayesTools/>
- Report bugs at <https://github.com/FBartos/BayesTools/issues>

---

BayesTools\_ensemble\_tables*Create BayesTools ensemble summary tables*

---

**Description**

Creates estimate summaries based on posterior distributions created by [mix\\_posteriors](#), inference summaries based on inference created by [ensemble\\_inference](#), or ensemble summary/diagnostics based on a list of [models\\_inference](#) models.

**Usage**

```
ensemble_estimates_table(  
  samples,  
  parameters,  
  probs = c(0.025, 0.95),  
  title = NULL,  
  footnotes = NULL,  
  warnings = NULL,  
  transform_orthonormal = FALSE,  
  formula_prefix = TRUE  
)
```

```

ensemble_inference_table(
  inference,
  parameters,
  logBF = FALSE,
  BF01 = FALSE,
  title = NULL,
  footnotes = NULL,
  warnings = NULL
)

ensemble_summary_table(
  models,
  parameters,
  logBF = FALSE,
  BF01 = FALSE,
  title = NULL,
  footnotes = NULL,
  warnings = NULL,
  remove_spike_0 = TRUE,
  short_name = FALSE
)

ensemble_diagnostics_table(
  models,
  parameters,
  title = NULL,
  footnotes = NULL,
  warnings = NULL,
  remove_spike_0 = TRUE,
  short_name = FALSE
)

```

### Arguments

samples	posterior samples created by <a href="#">mix_posteriors</a>
parameters	character vector of parameters (or a named list with of character vectors for summary and diagnostics tables) specifying the parameters (and their grouping) for the summary table
probs	quantiles for parameter estimates
title	title to be added to the table
footnotes	footnotes to be added to the table
warnings	warnings to be added to the table
transform_orthonormal	whether factors with orthonormal prior distributions should be transformed to differences from the grand mean
formula_prefix	whether the parameter prefix from formula should be printed. Defaults to TRUE.
inference	model inference created by <a href="#">ensemble_inference</a>

logBF	whether the Bayes factor should be on log scale
BF01	whether the Bayes factor should be inverted
models	list of <a href="#">models_inference</a> model objects, each of which containing a list of priors and inference object, The inference must be a named list with information about the model: model number <code>m_number</code> , marginal likelihood <code>marglik</code> , prior and posterior probability <code>prior_prob</code> and <code>post_prob</code> , inclusion Bayes factor <code>inclusion_BF</code> , and fit summary generated by <a href="#">runjags_estimates_table</a> for the diagnostics table
remove_spike_0	whether prior distributions equal to spike at 0 should be removed from the <code>prior_list</code>
short_name	whether the prior distribution names should be shortened. Defaults to FALSE.

**Value**

`ensemble_estimates_table` returns a table with the model-averaged estimates, `ensemble_inference_table` returns a table with the prior and posterior probabilities and inclusion Bayes factors, `ensemble_summary_table` returns a table with overview of the models included in the ensemble, and `ensemble_diagnostics_table` returns an overview of the MCMC diagnostics for the models included in the ensemble. All of the tables are objects of class 'BayesTools\_table'.

**See Also**

[ensemble\\_inference](#) [mix\\_posteriors](#) [BayesTools\\_model\\_tables](#)

---

BayesTools\_model\_tables

*Create BayesTools model tables*

---

**Description**

Creates model summary based on a model objects or provides estimates table for a runjags fit.

**Usage**

```
model_summary_table(
  model,
  model_description = NULL,
  title = NULL,
  footnotes = NULL,
  warnings = NULL,
  remove_spike_0 = TRUE,
  short_name = FALSE,
  formula_prefix = TRUE
)

runjags_estimates_table(
```

```
    fit,
    transformations = NULL,
    title = NULL,
    footnotes = NULL,
    warnings = NULL,
    conditional = FALSE,
    remove_spike_0 = TRUE,
    transform_orthonormal = FALSE,
    formula_prefix = TRUE,
    remove_inclusion = FALSE
  )

runjags_inference_table(
  fit,
  title = NULL,
  footnotes = NULL,
  warnings = NULL,
  formula_prefix = TRUE
)

JAGS_estimates_table(
  fit,
  transformations = NULL,
  title = NULL,
  footnotes = NULL,
  warnings = NULL,
  conditional = FALSE,
  remove_spike_0 = TRUE,
  transform_orthonormal = FALSE,
  formula_prefix = TRUE,
  remove_inclusion = FALSE
)

JAGS_inference_table(
  fit,
  title = NULL,
  footnotes = NULL,
  warnings = NULL,
  formula_prefix = TRUE
)

JAGS_summary_table(
  model,
  model_description = NULL,
  title = NULL,
  footnotes = NULL,
  warnings = NULL,
  remove_spike_0 = TRUE,
```

```

    short_name = FALSE,
    formula_prefix = TRUE
  )

runjags_estimates_empty_table(title = NULL, footnotes = NULL, warnings = NULL)

runjags_inference_empty_table(title = NULL, footnotes = NULL, warnings = NULL)

```

### Arguments

<code>model</code>	model object containing a list of priors and inference object, The inference must be a named list with information about the model: model number <code>m_number</code> , marginal likelihood <code>marglik</code> , prior and posterior probability <code>prior_prob</code> and <code>post_prob</code> , and model inclusion Bayes factor <code>inclusion_BF</code>
<code>model_description</code>	named list with additional description to be added to the table
<code>title</code>	title to be added to the table
<code>footnotes</code>	footnotes to be added to the table
<code>warnings</code>	warnings to be added to the table
<code>remove_spike_0</code>	whether prior distributions equal to spike at 0 should be removed from the <code>prior_list</code>
<code>short_name</code>	whether the prior distribution names should be shortened. Defaults to FALSE.
<code>formula_prefix</code>	whether the parameter prefix from formula should be printed. Defaults to TRUE.
<code>fit</code>	runjags model fit
<code>transformations</code>	named list of transformations to be applied to specific parameters
<code>conditional</code>	summarizes estimates conditional on being included in the model for spike and slab priors. Defaults to FALSE.
<code>transform_orthonormal</code>	whether factors with orthonormal prior distributions should be transformed to differences from the grand mean
<code>remove_inclusion</code>	whether estimates of the inclusion probabilities should be excluded from the summary table. Defaults to FALSE.

### Value

`model_summary_table` returns a table with overview of the fitted model, `runjags_estimates_table` returns a table with MCMC estimates, and `runjags_estimates_empty_table` returns an empty estimates table. All of the tables are objects of class 'BayesTools\_table'.

### See Also

[BayesTools\\_ensemble\\_tables](#)

---

check_input	<i>Check input</i>
-------------	--------------------

---

**Description**

A set of convenience functions for checking objects/arguments to a function passed by a user.

**Usage**

```
check_bool(x, name, check_length = 1, allow_NULL = FALSE, call = "")
```

```
check_char(  
  x,  
  name,  
  check_length = 1,  
  allow_values = NULL,  
  allow_NULL = FALSE,  
  call = ""  
)
```

```
check_real(  
  x,  
  name,  
  lower = -Inf,  
  upper = Inf,  
  allow_bound = TRUE,  
  check_length = 1,  
  allow_NULL = FALSE,  
  call = ""  
)
```

```
check_int(  
  x,  
  name,  
  lower = -Inf,  
  upper = Inf,  
  allow_bound = TRUE,  
  check_length = 1,  
  allow_NULL = FALSE,  
  call = ""  
)
```

```
check_list(  
  x,  
  name,  
  check_length = 0,  
  check_names = NULL,
```

```

    all_objects = FALSE,
    allow_other = FALSE,
    allow_NULL = FALSE,
    call = ""
)

```

### Arguments

x	object to be checked
name	name of the object that will be print in the error message.
check_length	length of the object to be checked. Defaults to 1. Set to 0 in order to not check object length.
allow_NULL	whether the object can be NULL. If so, no checks are executed.
call	string to be placed as a prefix to the error call.
allow_values	names of values allowed in a character vector. Defaults to NULL (do not check).
lower	lower bound of allowed values. Defaults to -Inf (do not check).
upper	upper bound of allowed values. Defaults to Inf (do not check).
allow_bound	whether the values at the boundary are allowed. Defaults to TRUE.
check_names	names of entries allowed in a list. Defaults to NULL (do not check).
all_objects	whether all entries in check_names must be present. Defaults to FALSE.
allow_other	whether additional entries then the specified in check_names might be present

### Value

returns NULL, called for the input check.

### Examples

```

# check whether the object is logical
check_bool(TRUE, name = "input")

# will throw an error on any other type
## Not run:
  check_bool("TRUE", name = "input")

## End(Not run)

```

---

contr.orthonormal	<i>Orthornomal contrast matrix</i>
-------------------	------------------------------------

---

**Description**

Return a matrix of orthornomal contrasts. Code is based on `stanova::contr.bayes` and corresponding to description by Rouder et al. (2012)

**Usage**

```
contr.orthonormal(n, contrasts = TRUE)
```

**Arguments**

n	a vector of levels for a factor, or the number of levels
contrasts	logical indicating whether contrasts should be computed

**Value**

A matrix with n rows and k columns, with  $k = n - 1$  if `contrasts = TRUE` and  $k = n$  if `contrasts = FALSE`.

**References**

Rouder JN, Morey RD, Speckman PL, Province JM (2012). “Default Bayes factors for ANOVA designs.” *Journal of Mathematical Psychology*, **56**(5), 356–374. doi:10.1016/j.jmp.2012.08.001.

**Examples**

```
contr.orthonormal(c(1, 2))  
contr.orthonormal(c(1, 2, 3))
```

---

density.prior	<i>Prior density</i>
---------------	----------------------

---

**Description**

Computes density of a prior distribution across a range of values.

**Usage**

```
## S3 method for class 'prior'
density(
  x,
  x_seq = NULL,
  x_range = NULL,
  x_range_quant = NULL,
  n_points = 1000,
  n_samples = 10000,
  force_samples = FALSE,
  individual = FALSE,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  truncate_end = TRUE,
  ...
)
```

**Arguments**

<code>x</code>	a prior
<code>x_seq</code>	sequence of x coordinates
<code>x_range</code>	vector of length two with lower and upper range for the support (used if <code>x_seq</code> is unspecified)
<code>x_range_quant</code>	quantile used for automatically obtaining <code>x_range</code> if both <code>x_range</code> and <code>x_seq</code> are unspecified. Defaults to 0.005 for all but Cauchy, Student-t, Gamma, and Inverse-gamma distributions that use 0.010.
<code>n_points</code>	number of equally spaced points in the <code>x_range</code> if <code>x_seq</code> is unspecified
<code>n_samples</code>	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with <code>force_samples = TRUE</code> )
<code>force_samples</code>	should prior be sampled instead of obtaining analytic solution whenever possible
<code>individual</code>	should individual densities be returned (e.g., in case of weightfunction)
<code>transformation</code>	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations: <b>lin</b> linear transformation in form of $a + b \cdot x$ <b>tanh</b> also known as Fisher's z transformation <b>exp</b> exponential transformation , or a list containing the transformation function <code>fun</code> , inverse transformation function <code>inv</code> , and the Jacobian of the transformation <code>jac</code> . See examples for details.
<code>transformation_arguments</code>	a list with named arguments for the transformation
<code>transformation_settings</code>	boolean indicating whether the settings the <code>x_seq</code> or <code>x_range</code> was specified on the transformed support

truncate_end	whether the density should be set to zero in for the endpoints of truncated distributions
...	additional arguments

**Value**

density.prior returns an object of class 'density'.

**See Also**

[prior\(\)](#)

---

ensemble_inference	<i>Compute posterior probabilities and inclusion Bayes factors</i>
--------------------	--

---

**Description**

Computes prior probabilities, posterior probabilities, and inclusion Bayes factors based either on (1) a list of models, vector of parameters, and a list of indicators the models represent the null or alternative hypothesis for each parameter, (2) on prior model odds, marginal likelihoods, and indicator whether the models represent the null or alternative hypothesis, or (3) list of models for each model.

**Usage**

```
compute_inference(prior_weights, margliks, is_null = NULL, conditional = FALSE)
ensemble_inference(model_list, parameters, is_null_list, conditional = FALSE)
models_inference(model_list)
```

**Arguments**

prior_weights	vector of prior model odds
margliks	vector of marginal likelihoods
is_null	logical vector of indicators specifying whether the model corresponds to the null or alternative hypothesis (or an integer vector indexing models corresponding to the null hypothesis)
conditional	whether prior and posterior model probabilities should be returned only for the conditional model. Defaults to FALSE
model_list	list of models, each of which contains marginal likelihood estimated with bridge sampling marglik and prior model odds prior_weights
parameters	vector of parameters names for which inference should be drawn
is_null_list	list with entries for each parameter carrying either logical vector of indicators specifying whether the model corresponds to the null or alternative hypothesis (or an integer vector indexing models corresponding to the null hypothesis)

**Value**

compute\_inference returns a named list of prior probabilities, posterior probabilities, and Bayes factors, ppoint gives the distribution function, ensemble\_inference gives a list of named lists of inferences for each parameter, and models\_inference returns a list of models, each expanded by the inference list.

**See Also**

[mix\\_posteriors](#) [BayesTools\\_ensemble\\_tables](#)

---

format_BF	<i>Format Bayes factor</i>
-----------	----------------------------

---

**Description**

Formats Bayes factor

**Usage**

```
format_BF(BF, logBF = FALSE, BF01 = FALSE, inclusion = FALSE)
```

**Arguments**

BF	Bayes factor(s)
logBF	log(BF)
BF01	1/BF
inclusion	whether the Bayes factor is an inclusion BF (for naming purposes)

**Value**

format\_BF returns a formatted Bayes factor.

---

geom_prior	<i>Add prior object to a ggplot</i>
------------	-------------------------------------

---

**Description**

Add prior object to a ggplot

**Usage**

```
geom_prior(
  x,
  xlim = NULL,
  x_seq = NULL,
  x_range_quant = NULL,
  n_points = 1000,
  n_samples = 10000,
  force_samples = FALSE,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  show_parameter = if (individual) 1 else NULL,
  individual = FALSE,
  rescale_x = FALSE,
  scale_y2 = 1,
  ...
)
```

**Arguments**

<code>x</code>	a prior
<code>xlim</code>	plotting range of the prior
<code>x_seq</code>	sequence of x coordinates
<code>x_range_quant</code>	quantile used for automatically obtaining <code>x_range</code> if both <code>x_range</code> and <code>x_seq</code> are unspecified. Defaults to 0.005 for all but Cauchy, Student-t, Gamma, and Inverse-gamma distributions that use 0.010.
<code>n_points</code>	number of equally spaced points in the <code>x_range</code> if <code>x_seq</code> is unspecified
<code>n_samples</code>	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with <code>force_samples = TRUE</code> )
<code>force_samples</code>	should prior be sampled instead of obtaining analytic solution whenever possible
<code>transformation</code>	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations: <b>lin</b> linear transformation in form of $a + b \cdot x$ <b>tanh</b> also known as Fisher's z transformation <b>exp</b> exponential transformation , or a list containing the transformation function <code>fun</code> , inverse transformation function <code>inv</code> , and the Jacobian of the transformation <code>jac</code> . See examples for details.
<code>transformation_arguments</code>	a list with named arguments for the transformation
<code>transformation_settings</code>	boolean indicating whether the settings the <code>x_seq</code> or <code>x_range</code> was specified on the transformed support

show_parameter	which parameter should be returned in case of multiple parameters per prior. Useful when priors for the omega parameter are plotted and individual = TRUE.
individual	should individual densities be returned (e.g., in case of weightfunction)
rescale_x	allows to rescale x-axis in case a weightfunction is plotted.
scale_y2	scaling factor for a secondary axis
...	additional arguments

**Value**

geom\_prior\_list returns an object of class 'ggplot'.

**See Also**

[plot.prior\(\)](#) [lines.prior\(\)](#)

---

geom_prior_list	<i>Add list of prior objects to a plot</i>
-----------------	--

---

**Description**

Add list of prior objects to a plot

**Usage**

```
geom_prior_list(
  prior_list,
  xlim = NULL,
  x_seq = NULL,
  x_range_quant = NULL,
  n_points = 500,
  n_samples = 10000,
  force_samples = FALSE,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  rescale_x = FALSE,
  scale_y2 = NULL,
  prior_list_mu = NULL,
  ...
)
```

**Arguments**

prior_list	list of prior distributions
xlim	x plotting range
x_seq	sequence of x coordinates
x_range_quant	quantile used for automatically obtaining x_range if both x_range and x_seq are unspecified. Defaults to 0.005 for all but Cauchy, Student-t, Gamma, and Inverse-gamme distributions that use 0.010.
n_points	number of equally spaced points in the x_range if x_seq is unspecified
n_samples	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with force_samples = TRUE)
force_samples	should prior be sampled instead of obtaining analytic solution whenever possible
transformation	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations:  <b>lin</b> linear transformation in form of $a + b*x$ <b>tanh</b> also known as Fisher's z transformation <b>exp</b> exponential transformation  , or a list containing the transformation function fun, inverse transformation function inv, and the Jacobian of the transformation jac. See examples for details.
transformation_arguments	a list with named arguments for the transformation
transformation_settings	boolean indicating whether the settings the x_seq or x_range was specified on the transformed support
rescale_x	allows to rescale x-axis in case a weightfunction is plotted.
scale_y2	scaling factor for a secondary axis
prior_list_mu	list of priors for the mu parameter required when plotting PET-PEESE
...	additional arguments

**Value**

geom\_prior\_list returns an object of class 'ggplot'.

**See Also**

[plot\\_prior\\_list\(\)](#) [lines\\_prior\\_list\(\)](#)

---

inclusion_BF	<i>Compute inclusion Bayes factors</i>
--------------	--

---

**Description**

Computes inclusion Bayes factors based on prior model probabilities, posterior model probabilities (or marginal likelihoods), and indicator whether the models represent the null or alternative hypothesis.

**Usage**

```
inclusion_BF(prior_probs, post_probs, margliks, is_null)
```

**Arguments**

prior_probs	vector of prior model probabilities
post_probs	vector of posterior model probabilities
margliks	vector of marginal likelihoods.
is_null	logical vector of indicators whether the model corresponds to the null or alternative hypothesis (or an integer vector indexing models corresponding to the null hypothesis)

**Details**

Supplying margliks as the input is preferred since it is better at dealing with under/overflow (posterior probabilities are very close to either 0 or 1). In case that both the post\_probs and margliks are supplied, the results are based on margliks.

**Value**

inclusion\_BF returns a Bayes factor.

---

interpret	<i>Interpret ensemble inference and estimates</i>
-----------	---

---

**Description**

Provides textual summary for posterior distributions created by [mix\\_posteriors](#) and ensemble inference created by [ensemble\\_inference](#).

**Usage**

```
interpret(inference, samples, specification, method)
```

**Arguments**

inference	model inference created by <a href="#">ensemble_inference</a>
samples	posterior samples created by <a href="#">mix_posteriors</a>
specification	list of lists specifying the generated text. Each inner list carries: (1) inference specifying the name of in the inference entry and optionally inference_name as a name to use in the text and inference_BF_name as a symbol to be used instead of the default "BF", (2) samples specifying the name of in the samples entry and optionally samples_name as a name to use in the text, samples_units as a unit text to be appended after the estimate, and samples_conditional specifying whether the estimate is conditional or model-averaged.
method	character specifying name of the method to be appended at the beginning of each sentence.

**Value**

interpret returns character.

**See Also**

[ensemble\\_inference](#) [mix\\_posteriors](#) [BayesTools\\_model\\_tables](#) [BayesTools\\_ensemble\\_tables](#)

---

is.prior	<i>Reports whether x is a a prior object</i>
----------	--

---

**Description**

Reports whether x is a a prior object. Note that point priors inherit the prior.simple property

**Usage**

```
is.prior(x)
is.prior.point(x)
is.prior.none(x)
is.prior.simple(x)
is.prior.discrete(x)
is.prior.vector(x)
is.prior.PET(x)
is.prior.PEERE(x)
```

```
is.prior.weightfunction(x)
is.prior.factor(x)
is.prior.orthonormal(x)
is.prior.dummy(x)
is.prior.spike_and_slab(x)
```

### Arguments

x                    an object of test

### Value

returns a boolean indicating whether the test object is a prior (of specific type).

### Examples

```
# create some prior distributions
p0 <- prior(distribution = "point", parameters = list(location = 0))
p1 <- prior_PET(distribution = "normal", parameters = list(mean = 0, sd = 1))

is.prior(p0)
is.prior.simple(p0)
is.prior.point(p0)
is.prior.PET(p0)

is.prior(p1)
is.prior.simple(p1)
is.prior.point(p1)
is.prior.PET(p1)
```

---

JAGS\_add\_priors            *Add 'JAGS' prior*

---

### Description

Adds priors to a 'JAGS' syntax.

### Usage

```
JAGS_add_priors(syntax, prior_list)
```

### Arguments

syntax                JAGS model syntax  
prior\_list            named list of prior distribution (names correspond to the parameter names)

**Value**

JAGS\_add\_priors returns a JAGS syntax.

---

JAGS\_bridgesampling     *Compute marginal likelihood of a 'JAGS' model*

---

**Description**

A wrapper around [bridge\\_sampler](#) that automatically computes likelihood part dependent on the prior distribution and prepares parameter samples. `log_posterior` must specify a function that takes two arguments - a named list of samples from the prior distributions and the data, and returns log likelihood of the model part.

**Usage**

```
JAGS_bridgesampling(
  fit,
  log_posterior,
  data = NULL,
  prior_list = NULL,
  formula_list = NULL,
  formula_data_list = NULL,
  formula_prior_list = NULL,
  add_parameters = NULL,
  add_bounds = NULL,
  maxiter = 10000,
  silent = TRUE,
  ...
)
```

**Arguments**

<code>fit</code>	model fitted with either <a href="#">runjags</a> posterior samples obtained with <a href="#">rjags-package</a>
<code>log_posterior</code>	function that takes a named list of samples, the data, and additional list of parameters passed as <code>...</code> as input and returns the log of the unnormalized posterior density of the model part
<code>data</code>	list containing data to fit the model (not including data for the formulas)
<code>prior_list</code>	named list of prior distribution (names correspond to the parameter names) of parameters not specified within the <code>formula_list</code>
<code>formula_list</code>	named list of formulas to be added to the model (names correspond to the parameter name created by each of the formula)
<code>formula_data_list</code>	named list of data frames containing data for each formula (names of the lists correspond to the parameter name created by each of the formula)

`formula_prior_list`      named list of named lists of prior distributions (names of the lists correspond to the parameter name created by each of the formula and the names of the prior distribution correspond to the parameter names) of parameters specified within the formula

`add_parameters`      vector of additional parameter names that should be used in bridgesampling but were not specified in the `prior_list`

`add_bounds`      list with two name vectors ("lb" and "up") containing lower and upper bounds of the additional parameters that were not specified in the `prior_list`

`maxiter`      maximum number of iterations for the [bridge\\_sampler](#)

`silent`      whether the progress should be printed, defaults to TRUE

`...`      additional argument to the [bridge\\_sampler](#) and `log_posterior` function

**Value**

JAGS\_bridgesampling returns an object of class 'bridge'.

**Examples**

```
## Not run:
# simulate data
set.seed(1)
data <- list(
  x = rnorm(10),
  N = 10
)
data$x

# define priors
priors_list <- list(mu = prior("normal", list(0, 1)))

# define likelihood for the data
model_syntax <-
"model{
  for(i in 1:N){
    x[i] ~ dnorm(mu, 1)
  }
}"

# fit the models
fit <- JAGS_fit(model_syntax, data, priors_list)

# define log posterior for bridge sampling
log_posterior <- function(parameters, data){
  sum(dnorm(data$x, parameters$mu, 1, log = TRUE))
}

# get marginal likelihoods
marglik <- JAGS_bridgesampling(fit, log_posterior, data, priors_list)

## End(Not run)
```

---

```
JAGS_bridgesampling_posterior
      Prepare 'JAGS' posterior for 'bridgesampling'
```

---

### Description

prepares posterior distribution for 'bridgesampling' by removing unnecessary parameters and attaching lower and upper bounds of parameters based on a list of prior distributions.

### Usage

```
JAGS_bridgesampling_posterior(
  posterior,
  prior_list,
  add_parameters = NULL,
  add_bounds = NULL
)
```

### Arguments

posterior	matrix of mcmc samples from the posterior distribution
prior_list	named list of prior distribution (names correspond to the parameter names) of parameters not specified within the formula_list
add_parameters	vector of additional parameter names that should be used in bridgesampling but were not specified in the prior_list
add_bounds	list with two name vectors ("lb" and "up") containing lower and upper bounds of the additional parameters that were not specified in the prior_list

### Value

JAGS\_bridgesampling\_posterior returns a matrix of posterior samples with 'lb' and 'ub' attributes carrying the lower and upper boundaries.

---

```
JAGS_check_and_list      Check and list 'JAGS' fitting settings
```

---

### Description

Checks and lists settings for the [JAGS\\_fit](#) function.

**Usage**

```
JAGS_check_and_list_fit_settings(
  chains,
  adapt,
  burnin,
  sample,
  thin,
  autofit,
  parallel,
  cores,
  silent,
  seed,
  check_mins = list(chains = 1, adapt = 50, burnin = 50, sample = 100, thin = 1),
  call = ""
)

JAGS_check_and_list_autofit_settings(
  autofit_control,
  skip_sample_extend = FALSE,
  call = ""
)
```

**Arguments**

chains	number of chains to be run, defaults to 4
adapt	number of samples used for adapting the MCMC chains, defaults to 500
burnin	number of burnin iterations of the MCMC chains, defaults to 1000
sample	number of sampling iterations of the MCMC chains, defaults to 4000
thin	thinning interval for the MCMC samples, defaults to 1
autofit	whether the models should be refitted until convergence criteria specified in <code>autofit_control</code> . Defaults to FALSE.
parallel	whether the chains should be run in parallel FALSE
cores	number of cores used for multithreading if <code>parallel = TRUE</code> , defaults to <code>chains</code>
silent	whether the function should proceed silently, defaults to TRUE
seed	seed for random number generation
check_mins	named list of minimal values for which should some input be checked. Defaults to: <b>chains</b> 1 <b>adapt</b> 50 <b>burnin</b> 50 <b>sample</b> 100 <b>thin</b> 1
call	string to be placed as a prefix to the error call.

autofit\_control

a list of arguments controlling the autofit function. Possible options are:

**max\_Rhat** maximum R-hat error for the autofit function. Defaults to 1.05.

**min\_ESS** minimum effective sample size. Defaults to 500.

**max\_error** maximum MCMC error. Defaults to 1.01.

**max\_SD\_error** maximum MCMC error as the proportion of standard deviation of the parameters. Defaults to 0.05.

**max\_time** list specifying the time `time` and units after which the automatic fitting function is stopped. The units arguments need to correspond to units passed to `difftime` function.

**sample\_extend** number of samples between each convergence check. Defaults to 1000.

skip\_sample\_extend

whether `sample_extend` is allowed to be NULL and skipped in the check

## Value

`JAGS_check_and_list_fit_settings` invisibly returns a list of checked fit settings. `JAGS_check_and_list_autofit_set` invisibly returns a list of checked autofit settings. parameter names.

---

JAGS\_check\_convergence

*Assess convergence of a runjags model*

---

## Description

Checks whether the supplied [runjags-package](#) model satisfied convergence criteria.

## Usage

```
JAGS_check_convergence(
  fit,
  prior_list,
  max_Rhat = 1.05,
  min_ESS = 500,
  max_error = 0.01,
  max_SD_error = 0.05
)
```

## Arguments

<code>fit</code>	a runjags model
<code>prior_list</code>	named list of prior distribution (names correspond to the parameter names)
<code>max_Rhat</code>	maximum R-hat error for the autofit function. Defaults to 1.05.
<code>min_ESS</code>	minimum effective sample size. Defaults to 500.
<code>max_error</code>	maximum MCMC error. Defaults to 1.01.
<code>max_SD_error</code>	maximum MCMC error as the proportion of standard deviation of the parameters. Defaults to 0.05.

**Value**

JAGS\_check\_convergence returns a boolean indicating whether the model converged or not, with an attribute 'errors' carrying the failed convergence checks (if any).

**See Also**

[JAGS\\_fit\(\)](#)

**Examples**

```
## Not run:
# simulate data
set.seed(1)
data <- list(
  x = rnorm(10),
  N = 10
)
data$x

# define priors
priors_list <- list(mu = prior("normal", list(0, 1)))

# define likelihood for the data
model_syntax <-
  "model{
    for(i in 1:N){
      x[i] ~ dnorm(mu, 1)
    }
  }"

# fit the models
fit <- JAGS_fit(model_syntax, data, priors_list)
JAGS_check_convergence(fit, priors_list)

## End(Not run)
```

---

JAGS\_diagnostics

*Plot diagnostics of a 'JAGS' model*

---

**Description**

Creates density plots, trace plots, and autocorrelation plots for a given parameter of a JAGS model.

**Usage**

```
JAGS_diagnostics(
  fit,
  parameter,
  type,
```

```
    plot_type = "base",
    xlim = NULL,
    ylim = NULL,
    lags = 30,
    n_points = 1000,
    transformations = NULL,
    transform_orthonormal = FALSE,
    short_name = FALSE,
    parameter_names = FALSE,
    formula_prefix = TRUE,
    ...
)

JAGS_diagnostics_density(
  fit,
  parameter,
  plot_type = "base",
  xlim = NULL,
  n_points = 1000,
  transformations = NULL,
  transform_orthonormal = FALSE,
  short_name = FALSE,
  parameter_names = FALSE,
  formula_prefix = TRUE,
  ...
)

JAGS_diagnostics_trace(
  fit,
  parameter,
  plot_type = "base",
  ylim = NULL,
  transformations = NULL,
  transform_orthonormal = FALSE,
  short_name = FALSE,
  parameter_names = FALSE,
  formula_prefix = TRUE,
  ...
)

JAGS_diagnostics_autocorrelation(
  fit,
  parameter,
  plot_type = "base",
  lags = 30,
  transformations = NULL,
  transform_orthonormal = FALSE,
  short_name = FALSE,
```

```

    parameter_names = FALSE,
    formula_prefix = TRUE,
    ...
)

```

### Arguments

<code>fit</code>	a JAGS model fitted via <a href="#">JAGS_fit()</a>
<code>parameter</code>	parameter to be plotted
<code>type</code>	what type of model diagnostic should be plotted. The available options are "density", "trace", and "autocorrelation"
<code>plot_type</code>	whether to use a base plot "base" or ggplot2 "ggplot" for plotting.
<code>xlim</code>	x plotting range
<code>ylim</code>	y plotting range
<code>lags</code>	number of lags to be shown for the autocorrelation plot. Defaults to 30.
<code>n_points</code>	number of equally spaced points in the <code>x_range</code> if <code>x_seq</code> is unspecified
<code>transformations</code>	named list of transformations to be applied to specific parameters
<code>transform_orthonormal</code>	whether factors with orthonormal prior distributions should be transformed to differences from the grand mean
<code>short_name</code>	whether prior distribution names should be shorted
<code>parameter_names</code>	whether parameter names should be printed
<code>formula_prefix</code>	whether the parameter prefix from formula should be printed. Defaults to TRUE.
<code>...</code>	additional arguments

### Value

`diagnostics` returns either NULL if `plot_type = "base"` or an object/list of objects (depending on the number of parameters to be plotted) of class 'ggplot2' if `plot_type = "ggplot2"`.

### See Also

[JAGS\\_fit\(\)](#) [JAGS\\_check\\_convergence\(\)](#)

---

JAGS\_evaluate\_formula *Evaluate JAGS formula using posterior samples*

---

**Description**

Evaluates a JAGS formula on a posterior distribution obtained from a fitted model.

**Usage**

```
JAGS_evaluate_formula(fit, formula, parameter, data, prior_list)
```

**Arguments**

fit	model fitted with either <a href="#">runjags</a> posterior samples obtained with <a href="#">rjags-package</a>
formula	formula specifying the right hand side of the assignment (the left hand side is ignored)
parameter	name of the parameter created with the formula
data	data.frame containing predictors included in the formula
prior_list	named list of prior distribution of parameters specified within the formula

**Value**

JAGS\_evaluate\_formula returns a matrix of the evaluated posterior samples on the supplied data.

**See Also**

[JAGS\\_fit\(\)](#) [JAGS\\_formula\(\)](#)

---

JAGS\_fit

*Fits a 'JAGS' model*

---

**Description**

A wrapper around [run.jags](#) that simplifies fitting 'JAGS' models with usage with pre-specified model part of the 'JAGS' syntax, data and list of prior distributions.

**Usage**

```
JAGS_fit(
  model_syntax,
  data = NULL,
  prior_list = NULL,
  formula_list = NULL,
  formula_data_list = NULL,
  formula_prior_list = NULL,
  chains = 4,
  adapt = 500,
  burnin = 1000,
  sample = 4000,
  thin = 1,
  autofit = FALSE,
  autofit_control = list(max_Rhat = 1.05, min_ESS = 500, max_error = 0.01, max_SD_error
    = 0.05, max_time = list(time = 60, unit = "mins"), sample_extend = 1000),
  parallel = FALSE,
  cores = chains,
  silent = TRUE,
  seed = NULL,
  add_parameters = NULL,
  required_packages = NULL
)
```

**Arguments**

<code>model_syntax</code>	jags syntax for the model part
<code>data</code>	list containing data to fit the model (not including data for the formulas)
<code>prior_list</code>	named list of prior distribution (names correspond to the parameter names) of parameters not specified within the <code>formula_list</code>
<code>formula_list</code>	named list of formulas to be added to the model (names correspond to the parameter name created by each of the formula)
<code>formula_data_list</code>	named list of data frames containing data for each formula (names of the lists correspond to the parameter name created by each of the formula)
<code>formula_prior_list</code>	named list of named lists of prior distributions (names of the lists correspond to the parameter name created by each of the formula and the names of the prior distribution correspond to the parameter names) of parameters specified within the formula
<code>chains</code>	number of chains to be run, defaults to 4
<code>adapt</code>	number of samples used for adapting the MCMC chains, defaults to 500
<code>burnin</code>	number of burnin iterations of the MCMC chains, defaults to 1000
<code>sample</code>	number of sampling iterations of the MCMC chains, defaults to 4000
<code>thin</code>	thinning interval for the MCMC samples, defaults to 1

<code>autofit</code>	whether the models should be refitted until convergence criteria specified in <code>autofit_control</code> . Defaults to FALSE.
<code>autofit_control</code>	a list of arguments controlling the autofit function. Possible options are: <b>max_Rhat</b> maximum R-hat error for the autofit function. Defaults to 1.05. <b>min_ESS</b> minimum effective sample size. Defaults to 500. <b>max_error</b> maximum MCMC error. Defaults to 1.01. <b>max_SD_error</b> maximum MCMC error as the proportion of standard deviation of the parameters. Defaults to 0.05. <b>max_time</b> list specifying the time <code>time</code> and units after which the automatic fitting function is stopped. The units arguments need to correspond to units passed to <code>difftime</code> function. <b>sample_extend</b> number of samples between each convergence check. Defaults to 1000.
<code>parallel</code>	whether the chains should be run in parallel FALSE
<code>cores</code>	number of cores used for multithreading if <code>parallel = TRUE</code> , defaults to <code>chains</code>
<code>silent</code>	whether the function should proceed silently, defaults to TRUE
<code>seed</code>	seed for random number generation
<code>add_parameters</code>	vector of additional parameter names that should be used monitored but were not specified in the <code>prior_list</code>
<code>required_packages</code>	character vector specifying list of packages containing JAGS models required for sampling (in case that the function is run in parallel or in detached R session). Defaults to NULL.

**Value**

JAGS\_fit returns an object of class 'runjags'.

**See Also**

[JAGS\\_check\\_convergence\(\)](#)

**Examples**

```
## Not run:
# simulate data
set.seed(1)
data <- list(
  x = rnorm(10),
  N = 10
)
data$x

# define priors
priors_list <- list(mu = prior("normal", list(0, 1)))
```

```
# define likelihood for the data
model_syntax <-
  "model{
    for(i in 1:N){
      x[i] ~ dnorm(mu, 1)
    }
  }"

# fit the models
fit <- JAGS_fit(model_syntax, data, priors_list)

## End(Not run)
```

---

JAGS\_formula

*Create JAGS formula syntax and data object*

---

### Description

Creates a JAGS formula syntax, prepares data input, and returns modified prior list for further processing in the JAGS\_fit function

### Usage

```
JAGS_formula(formula, parameter, data, prior_list)
```

### Arguments

formula	formula specifying the right hand side of the assignment (the left hand side is ignored)
parameter	name of the parameter to be created with the formula
data	data.frame containing predictors included in the formula
prior_list	named list of prior distribution of parameters specified within the formula

### Value

JAGS\_formula returns a list containing the formula JAGS syntax, JAGS data object, and modified prior\_list.

### See Also

[JAGS\\_fit\(\)](#)

**Examples**

```

# simulate data
set.seed(1)
df <- data.frame(
  y      = rnorm(60),
  x_cont = rnorm(60),
  x_bin  = rbinom(60, 1, .5),
  x_fac3 = factor(rep(c("A", "B", "C"), 20), levels = c("A", "B", "C")),
  x_fac4 = factor(rep(c("A", "B", "C", "D"), 15), levels = c("A", "B", "C", "D"))
)

# specify priors
prior_list <- list(
  "intercept" = prior("normal", list(0, 1)),
  "x_cont"     = prior("normal", list(0, .5)),
  "x_fac3"     = prior_factor("normal", list(0, 1), contrast = "treatment"),
  "x_fac4"     = prior_factor("mnormal", list(0, 1), contrast = "orthonormal"),
  "x_fac3:x_fac4" = prior_factor("mnormal", list(0, .5), contrast = "orthonormal")
)

# create the formula object
formula <- JAGS_formula(
  formula = ~ x_cont + x_fac3 * x_fac4,
  parameter = "mu", data = df, prior_list = prior_list)

```

---

JAGS\_get\_inits

*Create initial values for 'JAGS' model*


---

**Description**

Creates initial values for priors in a 'JAGS' model.

**Usage**

```
JAGS_get_inits(prior_list, chains, seed)
```

**Arguments**

prior_list	named list of prior distribution (names correspond to the parameter names)
chains	number of chains
seed	seed for random number generation

**Value**

JAGS\_add\_priors returns a list of JAGS initial values.

---

JAGS\_marglik\_parameters

*Extract parameters for 'JAGS' priors*

---

### Description

Extracts transformed parameters from the prior part of a 'JAGS' model inside of a 'bridgesampling' function (returns them as a named list)

### Usage

```
JAGS_marglik_parameters(samples, prior_list)
```

```
JAGS_marglik_parameters_formula(  
  samples,  
  formula_data_list,  
  formula_prior_list,  
  prior_list_parameters  
)
```

### Arguments

<code>samples</code>	samples provided by bridgesampling function
<code>prior_list</code>	named list of prior distribution (names correspond to the parameter names) of parameters not specified within the <code>formula_list</code>
<code>formula_data_list</code>	named list of data frames containing data for each formula (names of the lists correspond to the parameter name created by each of the formula)
<code>formula_prior_list</code>	named list of named lists of prior distributions (names of the lists correspond to the parameter name created by each of the formula and the names of the prior distribution correspond to the parameter names) of parameters specified within the formula
<code>prior_list_parameters</code>	named list of prior distributions on model parameters (not specified within the formula but that might scale the formula parameters)

### Value

JAGS\_marglik\_parameters returns a named list of (transformed) posterior samples.

---

JAGS\_marglik\_priors     *Compute marginal likelihood for 'JAGS' priors*

---

**Description**

Computes marginal likelihood for the prior part of a 'JAGS' model within 'bridgesampling' function

**Usage**

```
JAGS_marglik_priors(samples, prior_list)
```

```
JAGS_marglik_priors_formula(samples, formula_prior_list)
```

**Arguments**

samples	samples provided by bridgesampling function
prior_list	named list of prior distribution (names correspond to the parameter names) of parameters not specified within the formula_list
formula_prior_list	named list of named lists of prior distributions (names of the lists correspond to the parameter name created by each of the formula and the names of the prior distribution correspond to the parameter names) of parameters specified within the formula

**Value**

JAGS\_marglik\_priors returns a numeric value of likelihood evaluated at the current posterior sample.

---

JAGS\_to\_monitor     *Create list of monitored parameters for 'JAGS' model*

---

**Description**

Creates a vector of parameter names to be monitored in a 'JAGS' model.

**Usage**

```
JAGS_to_monitor(prior_list)
```

**Arguments**

prior_list	named list of prior distribution (names correspond to the parameter names)
------------	--

**Value**

JAGS\_to\_monitor returns a character vector of parameter names.

---

kitchen_rolls	<i>Kitchen Rolls data from Wagenmakers et al. (2015) replication study.</i>
---------------	---

---

**Description**

The data set contains mean NEO PI-R scores for two groups of students. Each of them filled a personality questionnaire while rotating a kitchen roll either clock or counter-clock wise. See Wagenmakers et al. (2015) for more details about the replication study and the <https://osf.io/uszvx/> for the original data.

**Usage**

```
kitchen_rolls
```

**Format**

A data.frame with 2 columns and 102 observations.

**Value**

a data.frame.

**References**

Wagenmakers E, Beek TF, Rotteveel M, Gierholz A, Matzke D, Steingroever H, Ly A, Verhagen J, Selker R, Sasiadek A, others (2015). “Turning the hands of time again: a purely confirmatory replication study and a Bayesian analysis.” *Frontiers in Psychology*, **6**, 494. [doi:10.3389/fpsyg.2015.00494](https://doi.org/10.3389/fpsyg.2015.00494).

---

lines.prior	<i>Add prior object to a plot</i>
-------------	-----------------------------------

---

**Description**

Add prior object to a plot

**Usage**

```
## S3 method for class 'prior'  
lines(  
  x,  
  xlim = NULL,  
  x_seq = NULL,  
  x_range_quant = NULL,  
  n_points = 1000,  
  n_samples = 10000,
```

```

    force_samples = FALSE,
    transformation = NULL,
    transformation_arguments = NULL,
    transformation_settings = FALSE,
    show_parameter = if (individual) 1 else NULL,
    individual = FALSE,
    rescale_x = FALSE,
    scale_y2 = 1,
    ...
)

```

### Arguments

<code>x</code>	a prior
<code>xlim</code>	plotting range of the prior
<code>x_seq</code>	sequence of x coordinates
<code>x_range_quant</code>	quantile used for automatically obtaining <code>x_range</code> if both <code>x_range</code> and <code>x_seq</code> are unspecified. Defaults to 0.005 for all but Cauchy, Student-t, Gamma, and Inverse-gamma distributions that use 0.010.
<code>n_points</code>	number of equally spaced points in the <code>x_range</code> if <code>x_seq</code> is unspecified
<code>n_samples</code>	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with <code>force_samples = TRUE</code> )
<code>force_samples</code>	should prior be sampled instead of obtaining analytic solution whenever possible
<code>transformation</code>	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations: <b>lin</b> linear transformation in form of $a + b \cdot x$ <b>tanh</b> also known as Fisher's z transformation <b>exp</b> exponential transformation , or a list containing the transformation function <code>fun</code> , inverse transformation function <code>inv</code> , and the Jacobian of the transformation <code>jac</code> . See examples for details.
<code>transformation_arguments</code>	a list with named arguments for the transformation
<code>transformation_settings</code>	boolean indicating whether the settings the <code>x_seq</code> or <code>x_range</code> was specified on the transformed support
<code>show_parameter</code>	which parameter should be returned in case of multiple parameters per prior. Useful when priors for the omega parameter are plotted and <code>individual = TRUE</code> .
<code>individual</code>	should individual densities be returned (e.g., in case of weightfunction)
<code>rescale_x</code>	allows to rescale x-axis in case a weightfunction is plotted.
<code>scale_y2</code>	scaling factor for a secondary axis
<code>...</code>	additional arguments

**Value**

lines.prior returns NULL.

**See Also**

[plot.prior\(\)](#) [geom\\_prior\(\)](#)

---

lines_prior_list	<i>Add list of prior objects to a plot</i>
------------------	--

---

**Description**

Add list of prior objects to a plot

**Usage**

```
lines_prior_list(
  prior_list,
  xlim = NULL,
  x_seq = NULL,
  x_range_quant = NULL,
  n_points = 500,
  n_samples = 10000,
  force_samples = FALSE,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  rescale_x = FALSE,
  scale_y2 = NULL,
  prior_list_mu = NULL,
  ...
)
```

**Arguments**

prior_list	list of prior distributions
xlim	x plotting range
x_seq	sequence of x coordinates
x_range_quant	quantile used for automatically obtaining x_range if both x_range and x_seq are unspecified. Defaults to 0.005 for all but Cauchy, Student-t, Gamma, and Inverse-gamme distributions that use 0.010.
n_points	number of equally spaced points in the x_range if x_seq is unspecified
n_samples	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with force_samples = TRUE)
force_samples	should prior be sampled instead of obtaining analytic solution whenever possible

**transformation** transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations:

- lin** linear transformation in form of  $a + b \cdot x$
- tanh** also known as Fisher's z transformation
- exp** exponential transformation

, or a list containing the transformation function `fun`, inverse transformation function `inv`, and the Jacobian of the transformation `jac`. See examples for details.

**transformation\_arguments** a list with named arguments for the transformation

**transformation\_settings** boolean indicating whether the settings the `x_seq` or `x_range` was specified on the transformed support

**rescale\_x** allows to rescale x-axis in case a weightfunction is plotted.

**scale\_y2** scaling factor for a secondary axis

**prior\_list\_mu** list of priors for the mu parameter required when plotting PET-PEESE

**...** additional arguments

**Value**

`lines_prior_list` returns NULL.

**See Also**

[plot\\_prior\\_list\(\)](#) [geom\\_prior\\_list\(\)](#)

---

mean.prior

*Prior mean*

---

**Description**

Computes mean of a prior distribution. (In case of orthonormal prior distributions for factors, the mean of for the deviations from intercept is returned.)

**Usage**

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

**Arguments**

`x` a prior

`...` unused

**Value**

a mean of an object of class 'prior'.

**See Also**

[prior\(\)](#)

**Examples**

```
# create a standard normal prior distribution
p1 <- prior(distribution = "normal", parameters = list(mean = 1, sd = 1))

# compute mean of the prior distribution
mean(p1)
```

---

mix\_posteriors

*Model-average posterior distributions*

---

**Description**

Model-averages posterior distributions based on a list of models, vector of parameters, and a list of indicators the models represent the null or alternative hypothesis for each parameter.

**Usage**

```
mix_posteriors(
  model_list,
  parameters,
  is_null_list,
  conditional = FALSE,
  seed = NULL,
  n_samples = 10000
)
```

**Arguments**

model_list	list of models, each of which contains marginal likelihood estimated with bridge sampling <code>marglik</code> and prior model odds <code>prior_weights</code>
parameters	vector of parameters names for which inference should be drawn
is_null_list	list with entries for each parameter carrying either logical vector of indicators specifying whether the model corresponds to the null or alternative hypothesis (or an integer vector indexing models corresponding to the null hypothesis)
conditional	whether prior and posterior model probabilities should be returned only for the conditional model. Defaults to FALSE
seed	integer specifying seed for sampling posteriors for model averaging. Defaults to 1.
n_samples	number of samples to be drawn for the model-averaged posterior distribution

**Value**

mix\_posteriors returns a named list of mixed posterior distributions (either a vector of matrix).

**See Also**

[ensemble\\_inference](#) [BayesTools\\_ensemble\\_tables](#)

---

parameter_names	<i>Clean parameter names from JAGS</i>
-----------------	--

---

**Description**

Removes additional formatting from parameter names outputted from JAGS.

**Usage**

```
format_parameter_names(
  parameters,
  formula_parameters = NULL,
  formula_prefix = TRUE
)

JAGS_parameter_names(parameters, formula_parameter = NULL)
```

**Arguments**

parameters      a vector of parameter names  
 formula\_parameters  
                   a vector of formula parameter prefix names  
 formula\_prefix   whether the formula\_parameters names should be kept. Defaults to TRUE.  
 formula\_parameter  
                   a formula parameter prefix name

**Value**

A character vector with reformatted parameter names.

**Examples**

```
format_parameter_names(c("mu_x_cont", "mu_x_fac3t", "mu_x_fac3t__xXx__x_cont"),
  formula_parameters = "mu")
```

plot.prior

*Plots a prior object***Description**

Plots a prior object

**Usage**

```
## S3 method for class 'prior'
plot(
  x,
  plot_type = "base",
  x_seq = NULL,
  xlim = NULL,
  x_range_quant = NULL,
  n_points = 1000,
  n_samples = 10000,
  force_samples = FALSE,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  show_figures = if (individual) -1 else NULL,
  individual = FALSE,
  rescale_x = FALSE,
  par_name = NULL,
  ...
)
```

**Arguments**

x	a prior
plot_type	whether to use a base plot "base" or ggplot2 "ggplot" for plotting.
x_seq	sequence of x coordinates
xlim	x plotting range
x_range_quant	quantile used for automatically obtaining x_range if both x_range and x_seq are unspecified. Defaults to 0.005 for all but Cauchy, Student-t, Gamma, and Inverse-gamme distributions that use 0.010.
n_points	number of equally spaced points in the x_range if x_seq is unspecified
n_samples	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with force_samples = TRUE)
force_samples	should prior be sampled instead of obtaining analytic solution whenever possible
transformation	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations:

	<b>lin</b> linear transformation in form of $a + b \cdot x$
	<b>tanh</b> also known as Fisher's z transformation
	<b>exp</b> exponential transformation
	, or a list containing the transformation function <code>fun</code> , inverse transformation function <code>inv</code> , and the Jacobian of the transformation <code>jac</code> . See examples for details.
<code>transformation_arguments</code>	a list with named arguments for the transformation
<code>transformation_settings</code>	boolean indicating whether the settings the <code>x_seq</code> or <code>x_range</code> was specified on the transformed support
<code>show_figures</code>	which figures should be returned in case of multiple plots are generated. Useful when priors for the omega parameter are plotted and <code>individual = TRUE</code> .
<code>individual</code>	should individual densities be returned (e.g., in case of weightfunction)
<code>rescale_x</code>	allows to rescale x-axis in case a weightfunction is plotted.
<code>par_name</code>	a type of parameter for which the prior is specified. Only relevant if the prior corresponds to a mu parameter that needs to be transformed.
<code>...</code>	additional arguments

**Value**

`plot.prior` returns either NULL or an object of class 'ggplot' if `plot_type` is `plot_type = "ggplot"`.

**See Also**

[prior\(\)](#) [lines.prior\(\)](#) [geom\\_prior\(\)](#)

**Examples**

```
# create some prior distributions
p0 <- prior(distribution = "point", parameters = list(location = 0))
p1 <- prior(distribution = "normal", parameters = list(mean = 0, sd = 1))
p2 <- prior(distribution = "normal", parameters = list(mean = 0, sd = 1), truncation = list(0, Inf))

# a default plot
plot(p0)

# manipulate line thickness and color, change the parameter name
plot(p1, lwd = 2, col = "blue", par_name = bquote(mu))

# use ggplot
plot(p2, plot_type = "ggplot")

# utilize the ggplot prior geom
plot(p2, plot_type = "ggplot", xlim = c(-2, 2)) + geom_prior(p1, col = "red", lty = 2)

# apply transformation
plot(p1, transformation = "exp")
```

plot\_models

*Plot estimates from models***Description**

Plot estimates from models

**Usage**

```
plot_models(
  model_list,
  samples,
  inference,
  parameter,
  plot_type = "base",
  prior = FALSE,
  conditional = FALSE,
  order = NULL,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  par_name = NULL,
  formula_prefix = TRUE,
  ...
)
```

**Arguments**

model_list	list of models, each of which contains marginal likelihood estimated with bridge sampling <code>marglik</code> and prior model odds <code>prior_weights</code>
samples	samples from a posterior distribution for a parameter generated by <a href="#">mix_posteriors</a> .
inference	object created by <a href="#">ensemble_inference</a> function
parameter	parameter name to be plotted. Does not support PET-PEESE and <code>weightfunction</code> .
plot_type	whether to use a base plot "base" or <code>ggplot2</code> "ggplot" for plotting.
prior	whether prior distribution should be added to the figure
conditional	whether conditional models should be displayed
order	list specifying ordering of the models. The first element describes whether the ordering should be "increasing" or "decreasing" and the second element describes whether the ordering should be based "model" order, "estimate" size, posterior "probability", or the inclusion "BF".
transformation	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations: <b>lin</b> linear transformation in form of $a + b \cdot x$

	<b>tanh</b> also known as Fisher's z transformation
	<b>exp</b> exponential transformation
	, or a list containing the transformation function <code>fun</code> , inverse transformation function <code>inv</code> , and the Jacobian of the transformation <code>jac</code> . See examples for details.
<code>transformation_arguments</code>	a list with named arguments for the transformation
<code>transformation_settings</code>	boolean indicating whether the settings the <code>x_seq</code> or <code>x_range</code> was specified on the transformed support
<code>par_name</code>	a type of parameter for which the prior is specified. Only relevant if the prior corresponds to a <code>mu</code> parameter that needs to be transformed.
<code>formula_prefix</code>	whether the <code>formula_parameters</code> names should be kept. Defaults to <code>TRUE</code> .
<code>...</code>	additional arguments. E.g.: <ul style="list-style-type: none"> <li>"<code>show_updating</code>" whether Bayes factors and change from prior to posterior odds should be shown on the secondary y-axis</li> <li>"<code>show_estimates</code>" whether posterior estimates and 95% CI should be shown on the secondary y-axis</li> <li>"<code>y_axis2</code>" whether the secondary y-axis should be shown</li> </ul>

### Details

Plots prior and posterior estimates of the same parameter across multiple models (prior distributions with orthonormal contrast) are always plotted as differences from the grand mean).

### Value

`plot_models` returns either `NULL` or an object of class `'ggplot'` if `plot_type` is `plot_type = "ggplot"`.

### See Also

[prior\(\)](#) [lines\\_prior\\_list\(\)](#) [geom\\_prior\\_list\(\)](#)

---

plot\_posterior

*Plot samples from the mixed posterior distributions*

---

### Description

Plot samples from the mixed posterior distributions

**Usage**

```
plot_posterior(
  samples,
  parameter,
  plot_type = "base",
  prior = FALSE,
  n_points = 1000,
  n_samples = 10000,
  force_samples = FALSE,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  rescale_x = FALSE,
  par_name = NULL,
  dots_prior = list(),
  ...
)
```

**Arguments**

samples	samples from a posterior distribution for a parameter generated by <a href="#">mix_posteriors</a> .
parameter	parameter name to be plotted. Use "PETPEESE" for PET-PEESE plot with parameters "PET" and "PEESE", and "weightfunction" for plotting a weightfunction with parameters "omega".
plot_type	whether to use a base plot "base" or ggplot2 "ggplot" for plotting.
prior	whether prior distribution should be added to the figure
n_points	number of equally spaced points in the x_range if x_seq is unspecified
n_samples	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with force_samples = TRUE)
force_samples	should prior be sampled instead of obtaining analytic solution whenever possible
transformation	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations: <b>lin</b> linear transformation in form of $a + b \cdot x$ <b>tanh</b> also known as Fisher's z transformation <b>exp</b> exponential transformation , or a list containing the transformation function fun, inverse transformation function inv, and the Jacobian of the transformation jac. See examples for details.
transformation_arguments	a list with named arguments for the transformation
transformation_settings	boolean indicating whether the settings the x_seq or x_range was specified on the transformed support
rescale_x	allows to rescale x-axis in case a weightfunction is plotted.

par_name	a type of parameter for which the prior is specified. Only relevant if the prior corresponds to a mu parameter that needs to be transformed.
dots_prior	additional arguments for the prior distribution plot
...	additional arguments

**Value**

plot\_posterior returns either NULL or an object of class 'ggplot' if plot\_type is plot\_type = "ggplot".

**See Also**

[prior\(\)](#) [lines\\_prior\\_list\(\)](#) [geom\\_prior\\_list\(\)](#)

---

plot_prior_list	<i>Plot a list of prior distributions</i>
-----------------	---

---

**Description**

Plot a list of prior distributions

**Usage**

```
plot_prior_list(
  prior_list,
  plot_type = "base",
  x_seq = NULL,
  xlim = NULL,
  x_range_quant = NULL,
  n_points = 500,
  n_samples = 10000,
  force_samples = FALSE,
  transformation = NULL,
  transformation_arguments = NULL,
  transformation_settings = FALSE,
  rescale_x = FALSE,
  par_name = NULL,
  prior_list_mu = NULL,
  ...
)
```

**Arguments**

prior_list	list of prior distributions
plot_type	whether to use a base plot "base" or ggplot2 "ggplot" for plotting.
x_seq	sequence of x coordinates

xlim	x plotting range
x_range_quant	quantile used for automatically obtaining x_range if both x_range and x_seq are unspecified. Defaults to 0.005 for all but Cauchy, Student-t, Gamma, and Inverse-gamme distributions that use 0.010.
n_points	number of equally spaced points in the x_range if x_seq is unspecified
n_samples	number of samples from the prior distribution if the density cannot be obtained analytically (or if samples are forced with force_samples = TRUE)
force_samples	should prior be sampled instead of obtaining analytic solution whenever possible
transformation	transformation to be applied to the prior distribution. Either a character specifying one of the prepared transformations:  <b>lin</b> linear transformation in form of $a + b \cdot x$ <b>tanh</b> also known as Fisher's z transformation <b>exp</b> exponential transformation  , or a list containing the transformation function fun, inverse transformation function inv, and the Jacobian of the transformation jac. See examples for details.
transformation_arguments	a list with named arguments for the transformation
transformation_settings	boolean indicating whether the settings the x_seq or x_range was specified on the transformed support
rescale_x	allows to rescale x-axis in case a weightfunction is plotted.
par_name	a type of parameter for which the prior is specified. Only relevant if the prior corresponds to a mu parameter that needs to be transformed.
prior_list_mu	list of priors for the mu parameter required when plotting PET-PEESE
...	additional arguments

**Value**

plot\_prior\_list returns either NULL or an object of class 'ggplot' if plot\_type is plot\_type = "ggplot".

**See Also**

[prior\(\)](#) [lines\\_prior\\_list\(\)](#) [geom\\_prior\\_list\(\)](#)

---

point	<i>Point mass distribution</i>
-------	--------------------------------

---

**Description**

Density, distribution function, quantile function and random generation for point distribution.

**Usage**

```
dpoint(x, location, log = FALSE)
```

```
rpoint(n, location)
```

```
ppoint(q, location, lower.tail = TRUE, log.p = FALSE)
```

```
qpoint(p, location, lower.tail = TRUE, log.p = FALSE)
```

**Arguments**

x, q	vector or matrix of quantiles.
location	vector of locations.
log, log.p	logical; if TRUE, probabilities p are given as log(p).
n	number of observations.
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise, $P[X \geq x]$ .
p	vector of probabilities.

**Value**

dpoint gives the density, ppoint gives the distribution function, qpoint gives the quantile function, and rpoint generates random deviates.

**Examples**

```
# draw samples from a point distribution  
rpoint(10, location = 1)
```

```
print.BayesTools_table  
Print a BayesTools table
```

---

**Description**

Print a BayesTools table

**Usage**

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

**Arguments**

x                    a BayesTools\_values\_tables  
...                  additional arguments.

**Value**

print.BayesTools\_table returns NULL.

---

```
print.prior            Prints a prior object
```

---

**Description**

Prints a prior object

**Usage**

```
## S3 method for class 'prior'  
print(  
  x,  
  short_name = FALSE,  
  parameter_names = FALSE,  
  plot = FALSE,  
  digits_estimates = 2,  
  silent = FALSE,  
  ...  
)
```

**Arguments**

x	a prior
short_name	whether prior distribution names should be shorted
parameter_names	whether parameter names should be printed
plot	to return <code>bquote</code> formatted prior name for plotting.
digits_estimates	number of decimals to be displayed for printed parameters.
silent	to silently return the print message.
...	additional arguments

**Value**

`print.prior` invisibly returns the print statement.

**See Also**

[prior\(\)](#)

**Examples**

```
# create some prior distributions
p0 <- prior(distribution = "point", parameters = list(location = 0))
p1 <- prior(distribution = "normal", parameters = list(mean = 0, sd = 1))

# print them
p0
p1

# use short names
print(p1, short_name = TRUE)

# print parameter names
print(p1, parameter_names = TRUE)

# generate bquote plotting syntax
plot(0, main = print(p1, plot = TRUE))
```

---

prior

*Creates a prior distribution*

---

**Description**

`prior` creates a prior distribution. The prior can be visualized by the `plot` function.

**Usage**

```
prior(
  distribution,
  parameters,
  truncation = list(lower = -Inf, upper = Inf),
  prior_weights = 1
)

prior_none(prior_weights = 1)
```

**Arguments**

distribution	name of the prior distribution. The possible options are "point" for a point density characterized by a location parameter. "normal" for a normal distribution characterized by a mean and sd parameters. "lognormal" for a lognormal distribution characterized by a meanlog and sdlog parameters. "cauchy" for a Cauchy distribution characterized by a location and scale parameters. Internally converted into a generalized t-distribution with df = 1. "t" for a generalized t-distribution characterized by a location, scale, and df parameters. "gamma" for a gamma distribution characterized by either shape and rate, or shape and scale parameters. The later is internally converted to the shape and rate parametrization "invgamma" for an inverse-gamma distribution characterized by a shape and scale parameters. The JAGS part uses a 1/gamma distribution with a shape and rate parameter. "beta" for a beta distribution characterized by an alpha and beta parameters. "exp" for an exponential distribution characterized by either rate or scale parameter. The later is internally converted to rate. "uniform" for a uniform distribution defined on a range from a to b
parameters	list of appropriate parameters for a given distribution.
truncation	list with two elements, lower and upper, that define the lower and upper truncation of the distribution. Defaults to <code>list(lower = -Inf, upper = Inf)</code> . The truncation is automatically set to the bounds of the support.
prior_weights	prior odds associated with a given distribution. The value is passed into the model fitting function, which creates models corresponding to all combinations of prior distributions for each of the model parameters and sets the model priors odds to the product of its prior distributions.

**Value**

`prior` and `prior_none` return an object of class 'prior'. A named list containing the distribution name, parameters, and prior weights.

**See Also**

[plot.prior\(\)](#), [Normal](#), [Lognormal](#), [Cauchy](#), [Beta](#), [Exponential](#), [LocationScaleT](#), [InvGamma](#).

**Examples**

```
# create a standard normal prior distribution
p1 <- prior(distribution = "normal", parameters = list(mean = 1, sd = 1))

# create a half-normal standard normal prior distribution
p2 <- prior(distribution = "normal", parameters = list(mean = 1, sd = 1),
truncation = list(lower = 0, upper = Inf))

# the prior distribution can be visualized using the plot function
# (see ?plot.prior for all options)
plot(p1)
```

---

prior\_factor

*Creates a prior distribution for factors*

---

**Description**

prior\_factor creates a prior distribution for fitting models with factor predictors. (Note that results across different operating systems might vary due to differences in JAGS numerical precision.)

**Usage**

```
prior_factor(
  distribution,
  parameters,
  truncation = list(lower = -Inf, upper = Inf),
  prior_weights = 1,
  contrast = "orthonormal"
)
```

**Arguments**

distribution    name of the prior distribution. The possible options are

- "point" for a point density characterized by a location parameter.
- "normal" for a normal distribution characterized by a mean and sd parameters.
- "lognormal" for a lognormal distribution characterized by a meanlog and sdlog parameters.
- "cauchy" for a Cauchy distribution characterized by a location and scale parameters. Internally converted into a generalized t-distribution with df = 1.
- "t" for a generalized t-distribution characterized by a location, scale, and df parameters.



---

prior_functions	<i>Elementary prior related functions</i>
-----------------	---

---

**Description**

Density (pdf / lpdf), distribution function (cdf / ccdf), quantile function (quant), random generation (rng), mean, standard deviation (sd), and marginal variants of the functions (mpdf, mlpf, mcdf, mccdf, mquant) for prior distributions.

**Usage**

```
## S3 method for class 'prior'  
rng(x, n, ...)
```

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

```
## S3 method for class 'prior'  
ccdf(x, q, ...)
```

```
## S3 method for class 'prior'  
lpdf(x, y, ...)
```

```
## S3 method for class 'prior'  
pdf(x, y, ...)
```

```
## S3 method for class 'prior'  
quant(x, p, ...)
```

```
## S3 method for class 'prior'  
mcdf(x, q, ...)
```

```
## S3 method for class 'prior'  
mccdf(x, q, ...)
```

```
## S3 method for class 'prior'  
mlpdf(x, y, ...)
```

```
## S3 method for class 'prior'  
mpdf(x, y, ...)
```

```
## S3 method for class 'prior'  
mquant(x, p, ...)
```

**Arguments**

x                    prior distribution

n	number of observations
...	unused arguments
q	vector or matrix of quantiles
y	vector of observations
p	vector of probabilities

**Value**

pdf (mpdf) and lpdf (mlpdf) give the (marginal) density and the log of (marginal) density, cdf (mcdf) and ccdf (mccdf) give the (marginal) distribution and the complement of (marginal) distribution function, quant (mquant) give the (marginal) quantile function, and rng generates random deviates for an object of class 'prior'.

**Examples**

```
# create a standard normal prior distribution
p1 <- prior(distribution = "normal", parameters = list(mean = 1, sd = 1))

# generate a random sample from the prior
rng(p1, 10)

# compute cumulative density function
cdf(p1, 0)

# obtain quantile
quant(p1, .5)

# compute probability density
pdf(p1, c(0, 1, 2))
```

---

prior\_functions\_methods

*Creates generics for common statistical functions*

---

**Description**

Density (pdf / lpdf), distribution function (cdf / ccdf), quantile function (quant), random generation (rng), mean, standard deviation (sd), and marginal variants of the functions (mpdf, mlpdf, mcdf, mccdf, mquant).

**Usage**

```
rng(x, ...)
```

```
cdf(x, ...)
```

```
ccdf(x, ...)  
quant(x, ...)  
lpdf(x, ...)  
pdf(x, ...)  
mcdf(x, ...)  
mccdf(x, ...)  
mquant(x, ...)  
mlpdf(x, ...)  
mpdf(x, ...)
```

### Arguments

x	main argument
...	unused arguments

### Value

pdf (mpdf) and lpdf (mlpdf) give the (marginal) density and the log of (marginal) density, cdf (mcdf) and ccdf (mccdf) give the (marginal) distribution and the complement of (marginal) distribution function, quant (mquant) give the (marginal) quantile function, and rng generates random deviates for an object of class 'prior'.

The pdf function proceeds to PDF graphics device if x is a character.

---

prior_informed	<i>Creates an informed prior distribution based on research</i>
----------------	---

---

### Description

prior\_informed creates an informed prior distribution based on past research. The prior can be visualized by the plot function.

### Usage

```
prior_informed(name, parameter = NULL, type = "smd")
```

**Arguments**

name	name of the prior distribution. There are many options based on prior psychological or medical research. For psychology, the possible options are "van Erp" for an informed prior distribution for the heterogeneity parameter tau of meta-analytic effect size estimates based on standardized mean differences (van Erp et al. 2017), "Oosterwijk" for an informed prior distribution for the effect sizes expected in social psychology based on prior elicitation with dr. Oosterwijk (Gronau et al. 2017).  For medicine, the possible options are based on Bartoš et al. (2021) who developed empirical prior distributions for the effect size and heterogeneity parameters of the continuous standardized outcomes based on the Cochrane database of systematic reviews. Use "Cochrane" for a prior distribution based on the whole database or call <code>print(prior_informed_medicine_names)</code> to inspect the names of all 46 subfields and set the appropriate parameter and type.
parameter	parameter name describing what prior distribution is supposed to be produced in cases where the name corresponds to multiple prior distributions. Relevant only for the empirical medical prior distributions.
type	prior type describing what prior distribution is supposed to be produced in cases where the name and parameter correspond to multiple prior distributions. Relevant only for the empirical medical prior distributions.

**Value**

`prior_informed` returns an object of class 'prior'.

**References**

- Bartoš F, Gronau QF, Timmers B, Otte WM, Ly A, Wagenmakers E (2021). "Bayesian model-averaged meta-analysis in medicine." *Statistics in Medicine*. doi:10.1002/sim.9170.
- Gronau QF, Van Erp S, Heck DW, Cesario J, Jonas KJ, Wagenmakers E (2017). "A Bayesian model-averaged meta-analysis of the power pose effect with informed and default priors: The case of felt power." *Comprehensive Results in Social Psychology*, 2(1), 123–138. doi:10.1080/23743603.2017.1326760.
- van Erp S, Verhagen J, Grasman RP, Wagenmakers E (2017). "Estimates of between-study heterogeneity for 705 meta-analyses reported in Psychological Bulletin from 1990–2013." *Journal of Open Psychology Data*, 5(1). doi:10.5334/jopd.33.

**See Also**

[prior\(\)](#), [prior\\_informed\\_medicine\\_names](#)

**Examples**

```
# prior distribution representing expected effect sizes in social psychology
# based on prior elicitation with dr. Oosterwijk
```

```
p1 <- prior_informed("Oosterwijk")

# the prior distribution can be visualized using the plot function
# (see ?plot.prior for all options)
plot(p1)

# empirical prior distribution for the standardized mean differences from the oral health
# medical subfield based on meta-analytic effect size estimates from the
# Cochrane database of systematic reviews
p2 <- prior_informed("Oral Health", parameter = "effect", type = "smd")
print(p2)
```

---

prior\_informed\_medicine\_names

*Names of medical subfields from the Cochrane database of systematic reviews*

---

### Description

Contain names identifying the individual subfields from the Cochrane database of systematic reviews. The individual elements correspond to valid name arguments for the `prior_informed()` function.

### Usage

```
prior_informed_medicine_names
```

### Format

An object of class character of length 47.

### Value

returns a character vector with names of medical subfields from Cochrane database of systematic reviews.

### See Also

`prior_informed()`

### Examples

```
print(prior_informed_medicine_names)
```

---

prior\_PP                      *Creates a prior distribution for PET or PEESE models*

---

### Description

prior creates a prior distribution for fitting a PET or PEESE style models in RoBMA. The prior distribution can be visualized by the plot function.

### Usage

```
prior_PET(
  distribution,
  parameters,
  truncation = list(lower = 0, upper = Inf),
  prior_weights = 1
)

prior_PEESE(
  distribution,
  parameters,
  truncation = list(lower = 0, upper = Inf),
  prior_weights = 1
)
```

### Arguments

**distribution**    name of the prior distribution. The possible options are

- "point" for a point density characterized by a location parameter.
- "normal" for a normal distribution characterized by a mean and sd parameters.
- "lognormal" for a lognormal distribution characterized by a meanlog and sdlog parameters.
- "cauchy" for a Cauchy distribution characterized by a location and scale parameters. Internally converted into a generalized t-distribution with df = 1.
- "t" for a generalized t-distribution characterized by a location, scale, and df parameters.
- "gamma" for a gamma distribution characterized by either shape and rate, or shape and scale parameters. The later is internally converted to the shape and rate parametrization
- "invgamma" for an inverse-gamma distribution characterized by a shape and scale parameters. The JAGS part uses a 1/gamma distribution with a shape and rate parameter.
- "beta" for a beta distribution characterized by an alpha and beta parameters.
- "exp" for an exponential distribution characterized by either rate or scale parameter. The later is internally converted to rate.

	"uniform" for a uniform distribution defined on a range from a to b
parameters	list of appropriate parameters for a given distribution.
truncation	list with two elements, lower and upper, that define the lower and upper truncation of the distribution. Defaults to <code>list(lower = -Inf, upper = Inf)</code> . The truncation is automatically set to the bounds of the support.
prior_weights	prior odds associated with a given distribution. The value is passed into the model fitting function, which creates models corresponding to all combinations of prior distributions for each of the model parameters and sets the model priors odds to the product of its prior distributions.

**Value**

`prior_PET` and `prior_PEESE` return an object of class 'prior'.

**See Also**

[plot.prior\(\)](#), [prior\(\)](#)

**Examples**

```
# create a half-Cauchy prior distribution
# (PET and PEESE specific functions automatically set lower truncation at 0)
p1 <- prior_PET(distribution = "Cauchy", parameters = list(location = 0, scale = 1))

plot(p1)
```

---

`prior_spike_and_slab` *Creates a spike and slab prior distribution*

---

**Description**

`prior_spike_and_slab` creates a spike and slab prior distribution corresponding to the specification in Kuo and Mallick (1998) (see O'Hara and Sillanpää (2009) for further details). I.e., a prior distribution is multiplied by an independent indicator with values either zero or one.

**Usage**

```
prior_spike_and_slab(
  prior_parameter,
  prior_inclusion = prior(distribution = "spike", parameters = list(location = 0.5)),
  prior_weights = 1
)
```

**Arguments**

- `prior_parameter` a prior distribution for the parameter
- `prior_inclusion` a prior distribution for the inclusion probability. The inclusion probability must be bounded within 0 and 1 range. Defaults to `prior("spike", parameters = list(location = 0.5))` which corresponds to 1/2 prior probability of including the slab prior distribution (but other prior distributions, like beta etc can be also specified).
- `prior_weights` prior odds associated with a given distribution. The value is passed into the model fitting function, which creates models corresponding to all combinations of prior distributions for each of the model parameters and sets the model priors odds to the product of its prior distributions.

**Value**

return an object of class 'prior'.

**See Also**

[prior\(\)](#)

**Examples**

```
# create a spike and slab prior distribution
p1 <- prior_spike_and_slab(
  prior(distribution = "normal", parameters = list(mean = 0, sd = 1)),
  prior_inclusion = prior(distribution = "beta", parameters = list(alpha = 1, beta = 1))
)
```

---

`prior_weightfunction` *Creates a prior distribution for a weight function*

---

**Description**

`prior_weightfunction` creates a prior distribution for fitting a RoBMA selection model. The prior can be visualized by the `plot` function.

**Usage**

```
prior_weightfunction(distribution, parameters, prior_weights = 1)
```

**Arguments**

- `distribution` name of the prior distribution. The possible options are  
 "two.sided" for a two-sided weight function characterized by a vector `steps` and vector `alpha` parameters. The `alpha` parameter determines an alpha parameter of Dirichlet distribution which cumulative sum is used for the weights `omega`.  
 "one.sided" for a one-sided weight function characterized by either a vector `steps` and vector `alpha` parameter, leading to a monotonic one-sided function, or by a vector `steps`, vector `alpha1`, and vector `alpha2` parameters leading non-monotonic one-sided weight function. The `alpha` / `alpha1` and `alpha2` parameters determine an alpha parameter of Dirichlet distribution which cumulative sum is used for the weights `omega`.
- `parameters` list of appropriate parameters for a given distribution.
- `prior_weights` prior odds associated with a given distribution. The model fitting function usually creates models corresponding to all combinations of prior distributions for each of the model parameters, and sets the model priors odds to the product of its prior distributions.

**Value**

`prior_weightfunction` returns an object of class 'prior'.

**See Also**

[plot.prior\(\)](#)

**Examples**

```
p1 <- prior_weightfunction("one-sided", parameters = list(steps = c(.05, .10), alpha = c(1, 1, 1)))

# the prior distribution can be visualized using the plot function
# (see ?plot.prior for all options)
plot(p1)
```

---

range.prior

*Prior range*

---

**Description**

Computes range of a prior distribution (if the prior distribution is unbounded range from quantiles to 1 -quantiles) is returned.

**Usage**

```
## S3 method for class 'prior'
range(x, quantiles = NULL, ..., na.rm = FALSE)
```

**Arguments**

x	a prior
quantiles	quantile to be returned in case of unbounded distribution.
...	additional arguments
na.rm	unused

**Value**

range.prior returns a numeric vector of length with a plotting range of a prior distribution.

**See Also**

[prior\(\)](#)

---

sd	<i>Creates generic for sd function</i>
----	--

---

**Description**

Creates generic for sd function

**Usage**

```
sd(x, ...)
```

**Arguments**

x	main argument
...	additional arguments

**Value**

sd returns a standard deviation of the supplied object (if it is either a numeric vector or an object of class 'prior').

**See Also**

[sd](#)

---

sd.prior	<i>Prior sd</i>
----------	-----------------

---

**Description**

Computes standard deviation of a prior distribution.

**Usage**

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

**Arguments**

x	a prior
...	unused arguments

**Value**

a standard deviation of an object of class 'prior'.

**See Also**

[prior\(\)](#)

**Examples**

```
# create a standard normal prior distribution
p1 <- prior(distribution = "normal", parameters = list(mean = 1, sd = 1))

# compute sd of the prior distribution
sd(p1)
```

---

transform_orthonormal_samples	<i>Transform orthonormal posterior samples into differences from the mean</i>
-------------------------------	---

---

**Description**

Transforms posterior samples from model-averaged posterior distributions based on orthonormal prior distributions into differences from the mean.

**Usage**

```
transform_orthonormal_samples(samples)
```

**Arguments**

`samples` (a list) of mixed posterior distributions created with `mix_posteriors` function

**Value**

`transform_orthonormal_samples` returns a named list of mixed posterior distributions (either a vector of matrix).

**See Also**

[mix\\_posteriors](#) [contr.orthonormal](#)

---

`var` *Creates generic for var function*

---

**Description**

Creates generic for var function

**Usage**

```
var(x, ...)
```

**Arguments**

`x` main argument  
`...` additional arguments

**Value**

`var` returns a variance of the supplied object (if it is either a numeric vector or an object of class 'prior').

**See Also**

[cor](#)

---

var.prior	<i>Prior var</i>
-----------	------------------

---

**Description**

Computes variance of a prior distribution.

**Usage**

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

**Arguments**

x	a prior
...	unused arguments

**Value**

a variance of an object of class 'prior'.

**See Also**

[prior\(\)](#)

**Examples**

```
# create a standard normal prior distribution  
p1 <- prior(distribution = "normal", parameters = list(mean = 1, sd = 1))  
  
# compute variance of the prior distribution  
var(p1)
```

---

weightfunctions	<i>Weight functions</i>
-----------------	-------------------------

---

**Description**

Marginal density, marginal distribution function, marginal quantile function and random generation for weight functions.

**Usage**

```
mdone.sided(x, alpha = NULL, alpha1 = NULL, alpha2 = NULL, log = FALSE)
mdtwo.sided(x, alpha, log = FALSE)
mdone.sided_fixed(x, omega, log = FALSE)
mdtwo.sided_fixed(x, omega, log = FALSE)
rone.sided(n, alpha = NULL, alpha1 = NULL, alpha2 = NULL)
rtwo.sided(n, alpha)
rone.sided_fixed(n, omega)
rtwo.sided_fixed(n, omega)

mpone.sided(
  q,
  alpha = NULL,
  alpha1 = NULL,
  alpha2 = NULL,
  lower.tail = TRUE,
  log.p = FALSE
)
mptwo.sided(q, alpha, lower.tail = TRUE, log.p = FALSE)
mpone.sided_fixed(q, omega, lower.tail = TRUE, log.p = FALSE)
mptwo.sided_fixed(q, omega, lower.tail = TRUE, log.p = FALSE)

mqone.sided(
  p,
  alpha = NULL,
  alpha1 = NULL,
  alpha2 = NULL,
  lower.tail = TRUE,
  log.p = FALSE
)
mqtwo.sided(p, alpha, lower.tail = TRUE, log.p = FALSE)
mqone.sided_fixed(p, omega, lower.tail = TRUE, log.p = FALSE)
mqtwo.sided_fixed(p, omega, lower.tail = TRUE, log.p = FALSE)
```

**Arguments**

x, q	vector or matrix of quantiles.
alpha	vector or matrix with concentration parameters for the Dirichlet distribution for a monotonic one.sided or a two.sided weight function.
alpha1	vector or matrix with concentration parameters for the Dirichlet distribution for the expected direction of non-monotonic one.sided of weight function.
alpha2	vector or matrix with concentration parameters for the Dirichlet distribution for the unexpected direction of non-monotonic one.sided of weight function.
log, log.p	logical; if TRUE, probabilities p are given as log(p).
omega	vector or matrix of fixed probabilities for a one.sided or a two.sided weight function.
n	number of observations.
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise, $P[X \geq x]$ .
p	vector of probabilities.

**Value**

mdone.sided, mdtwo.sided, mdone.sided\_fixed, and mdtwo.sided\_fixed give the marginal density, mpone.sided, mptwo.sided, mpone.sided\_fixed, and mptwo.sided\_fixed give the marginal distribution function, mqone.sided, mqtwo.sided, mqone.sided\_fixed, and mqtwo.sided\_fixed give the marginal quantile function, and rone.sided, rtwo.sided, rone.sided\_fixed, and rtwo.sided\_fixed generate random deviates.

**Examples**

```
# draw samples from a two-sided weight function
rtwo.sided(10, alpha = c(1, 1))

# draw samples from a monotone one-sided weight function
rone.sided(10, alpha = c(1, 1, 1))

# draw samples from a non-monotone one-sided weight function
rone.sided(10, alpha1 = c(1, 1), alpha2 = c(1, 1))
```

---

weightfunctions\_mapping

*Create coefficient mapping between multiple weightfunctions*

---

**Description**

Creates coefficients mapping between multiple weightfunctions.

**Usage**

```
weightfunctions_mapping(prior_list, cuts_only = FALSE)
```

**Arguments**

`prior_list` list of prior distributions  
`cuts_only` whether only p-value cuts should be returned

**Value**

`weightfunctions_mapping` returns a list of indices mapping the publication weights  $\omega$  from the individual weightfunctions into a joint weightfunction.

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