Package ‘bayestestR’

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**Type**  Package

**Title**  Understand and Describe Bayesian Models and Posterior Distributions

**Version**  0.12.1

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**Description**  Provides utilities to describe posterior distributions and Bayesian models. It includes point-estimates such as Maximum A Posteriori (MAP), measures of dispersion (Highest Density Interval - HDI; Kruschke, 2015 <doi:10.1016/C2012-0-00477-2>) and indices used for null-hypothesis testing (such as ROPE percentage, pd and Bayes factors).

**Depends**  R (>= 3.4)

**Imports**  insight (>= 0.17.0), datawizard (>= 0.4.0), graphics, methods, stats, utils

**Suggests**  BayesFactor, bayesQR, blavaan, bridgesampling, brms, dplyr, effectsize, emmeans, gamm4, GGally, ggdist, ggplot2, ggridges, glmTMB, httr, KernSmooth, knitr, lavaan, lme4, logspline, MASS, mclust, mediation, modelbased, parameters, patchwork, performance, poorman, quadprog, posterior, rmarkdown, rstan, rstanarm, see, spelling, stringr, testthat, tidyr, tweedie

**License**  GPL-3

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Description

Based on the DescTools AUC function. It can calculate the area under the curve with a naive algorithm or a more elaborated spline approach. The curve must be given by vectors of xy-coordinates. This function can handle unsorted x values (by sorting x) and ties for the x values (by ignoring duplicates).

Usage

```r
area_under_curve(x, y, method = c("trapezoid", "step", "spline"), ...)
```

```r
auc(x, y, method = c("trapezoid", "step", "spline"), ...)
```

Arguments

- `x` Vector of x values.
- `y` Vector of y values.
as.data.frame.density

Method to compute the Area Under the Curve (AUC). Can be "trapezoid" (default), "step" or "spline". If "trapezoid", the curve is formed by connecting all points by a direct line (composite trapezoid rule). If "step" is chosen then a stepwise connection of two points is used. For calculating the area under a spline interpolation the splinefun function is used in combination with integrate.

Arguments passed to or from other methods.

See Also
DescTools

Examples

library(bayestestR)
posterior <- distribution_normal(1000)

dens <- estimate_density(posterior)
dens <- dens[dens$x > 0,]
x <- dens$x
y <- dens$y

area_under_curve(x, y, method = "trapezoid")
area_under_curve(x, y, method = "step")
area_under_curve(x, y, method = "spline")

as.data.frame.density Coerce to a Data Frame

Description
Coerce to a Data Frame

Usage

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

Arguments

x any R object.
... additional arguments to be passed to or from methods.
Description

Convert to Numeric

Usage

```r
## S3 method for class 'map_estimate' 
as.numeric(x, ...)

## S3 method for class 'p_direction' 
as.numeric(x, ...)

## S3 method for class 'p_map' 
as.numeric(x, ...)

## S3 method for class 'p_significance' 
as.numeric(x, ...)
```

Arguments

- `x` object to be coerced or tested.
- `...` further arguments passed to or from other methods.

bayesfactor  Bayes Factors (BF)

Description

This function compute the Bayes factors (BFs) that are appropriate to the input. For vectors or single models, it will compute BFs for single parameters(), or is hypothesis is specified, BFs for restricted models(). For multiple models, it will return the BF corresponding to comparison between models() and if a model comparison is passed, it will compute the inclusion BF().

For a complete overview of these functions, read the Bayes factor vignette.
Usage

bayesfactor(
  ..., 
  prior = NULL, 
  direction = "two-sided", 
  null = 0, 
  hypothesis = NULL, 
  effects = c("fixed", "random", "all"), 
  verbose = TRUE, 
  denominator = 1, 
  match.models = FALSE, 
  prior_odds = NULL
)

Arguments

... A numeric vector, model object(s), or the output from bayesfactor_models.
prior An object representing a prior distribution (see 'Details').
direction Test type (see 'Details'). One of 0, "two-sided" (default, two tailed), -1, "left" (left tailed) or 1, "right" (right tailed).
null Value of the null, either a scalar (for point-null) or a range (for a interval-null).
hypothesis A character vector specifying the restrictions as logical conditions (see examples below).
effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
verbose Toggle off warnings.
denominator Either an integer indicating which of the models to use as the denominator, or a model to be used as a denominator. Ignored for BFBayesFactor.
match.models See details.
prior_odds Optional vector of prior odds for the models. See BayesFactor::priorOdds<-

Value

Some type of Bayes factor, depending on the input. See bayesfactor_parameters(), bayesfactor_models() or bayesfactor_inclusion()

Note

There is also a plot()-method implemented in the see-package.

Examples

library(bayestestR)

if (require("logspline")) {
  prior <- distribution_normal(1000, mean = 0, sd = 1)


```r
posterior <- distribution_normal(1000, mean = .5, sd = .3)
bayesfactor(posterior, prior = prior)
}
## Not run:
# rstanarm models
# ---------------
if (require("rstanarm")) {
    model <- stan_lmer(extra ~ group + (1 | ID), data = sleep)
    bayesfactor(model)
}
## End(Not run)

if (require("logspline")) {
    # Frequentist models
    # ---------------
    m0 <- lm(extra ~ 1, data = sleep)
    m1 <- lm(extra ~ group, data = sleep)
    m2 <- lm(extra ~ group + ID, data = sleep)

    comparison <- bayesfactor(m0, m1, m2)
    comparison
    bayesfactor(comparison)
}
```

---

**bayesfactor_inclusion**  
*Inclusion Bayes Factors for testing predictors across Bayesian models*

**Description**

The `bf_*` function is an alias of the main function.

For more info, see the [Bayes factors vignette](#).

**Usage**

```r
bayesfactor_inclusion(models, match_models = FALSE, prior_odds = NULL, ...)
bayesfactor_inclusion(models, match_models = FALSE, prior_odds = NULL, ...)
```

**Arguments**

- `models`  
  An object of class `bayesfactor_models()` or `BFBayesFactor`.

- `match_models`  
  See details.

- `prior_odds`  
  Optional vector of prior odds for the models. See `BayesFactor::priorOdds<-`.

- `...`  
  Arguments passed to or from other methods.
Details

Inclusion Bayes factors answer the question: Are the observed data more probable under models with a particular effect, than they are under models without that particular effect? In other words, on average - are models with effect $X$ more likely to have produced the observed data than models without effect $X$?

**Match Models:** If `match_models=FALSE` (default), Inclusion BFs are computed by comparing all models with a term against all models without that term. If `TRUE`, comparison is restricted to models that (1) do not include any interactions with the term of interest; (2) for interaction terms, averaging is done only across models that contain the main effect terms from which the interaction term is comprised.

Value

a data frame containing the prior and posterior probabilities, and log(BF) for each effect (Use `as.numeric()` to extract the non-log Bayes factors; see examples).

Interpreting Bayes Factors

A Bayes factor greater than 1 can be interpreted as evidence against the null, at which one convention is that a Bayes factor greater than 3 can be considered as "substantial" evidence against the null (and vice versa, a Bayes factor smaller than 1/3 indicates substantial evidence in favor of the null-model) (Wetzels et al. 2011).

Note

Random effects in the lmer style are converted to interaction terms: i.e., $(X|G)$ will become the terms $1:G$ and $X:G$.

Author(s)

Mattan S. Ben-Shachar

References


See Also

`weighted_posteriors()` for Bayesian parameter averaging.
Bayes Factors (BF) for model comparison

Description

This function computes or extracts Bayes factors from fitted models.

The `bf_*` function is an alias of the main function.

Usage

`bayesfactor_models(..., denominator = 1, verbose = TRUE)`

`bf_models(..., denominator = 1, verbose = TRUE)`

## Default S3 method:
`bayesfactor_models(..., denominator = 1, verbose = TRUE)`

## S3 method for class 'bayesfactor_models'
`update(object, subset = NULL, reference = NULL, ...)`

Examples

```r
library(bayestestR)

# Using bayesfactor_models:
# ------------------------------
mo0 <- lm(Sepal.Length ~ 1, data = iris)
mo1 <- lm(Sepal.Length ~ Species, data = iris)
mo2 <- lm(Sepal.Length ~ Species + Petal.Length, data = iris)
mo3 <- lm(Sepal.Length ~ Species * Petal.Length, data = iris)

BFmodels <- bayesfactor_models(mo1, mo2, mo3, denominator = mo0)
(bf_inc <- bayesfactor_inclusion(BFmodels))

as.numeric(bf_inc)

## Not run:
# BayesFactor
# -------------------------------
library(BayesFactor)

BF <- generalTestBF(len ~ supp * dose, ToothGrowth, progress = FALSE)

bayesfactor_inclusion(BF)

# compare only matched models:
bayesfactor_inclusion(BF, match_models = TRUE)

## End(Not run)
```
## S3 method for class 'bayesfactor_models'

as.matrix(x, ...)

### Arguments

- **...**
  - Fitted models (see details), all fit on the same data, or a single `BFBayesFactor` object (see 'Details'). Ignored in `as.matrix()`, `update()`. If the following named arguments are present, they are passed to `insight::get_loglikelihood` (see details):
    - `estimator` (defaults to "ML")
    - `check_response` (defaults to FALSE)

- **denominator**
  - Either an integer indicating which of the models to use as the denominator, or a model to be used as a denominator. Ignored for `BFBayesFactor`.

- **verbose**
  - Toggle off warnings.

- **object, x**
  - A `bayesfactor_models()` object.

- **subset**
  - Vector of model indices to keep or remove.

- **reference**
  - Index of model to reference to, or "top" to reference to the best model, or "bottom" to reference to the worst model.

### Details

If the passed models are supported by `insight` the DV of all models will be tested for equality (else this is assumed to be true), and the models’ terms will be extracted (allowing for follow-up analysis with `bayesfactor_inclusion`).

- For `brmsfit` or `stanreg` models, Bayes factors are computed using the `bridgesampling` package.
  - `brmsfit` models must have been fitted with `save_pars = save_pars(all = TRUE)`.
  - `stanreg` models must have been fitted with a defined `diagnostic_file`.

- For `BFBayesFactor`, `bayesfactor_models()` is mostly a wraparound `BayesFactor::extractBF()`.

- For all other model types, Bayes factors are computed using the BIC approximation. Note that BICs are extracted from using `insight::get_loglikelihood`, see documentation there for options for dealing with transformed responses and REML estimation.

In order to correctly and precisely estimate Bayes factors, a rule of thumb are the 4 P’s: **Proper Priors** and **Plentiful Posteriors**. How many? The number of posterior samples needed for testing is substantially larger than for estimation (the default of 4000 samples may not be enough in many cases). A conservative rule of thumb is to obtain 10 times more samples than would be required for estimation (Gronau, Singmann, & Wagenmakers, 2017). If less than 40,000 samples are detected, `bayesfactor_models()` gives a warning.

See also the [Bayes factors vignette](#).

### Value

A data frame containing the models’ formulas (reconstructed fixed and random effects) and their log(BF)s (Use `as.numeric()` to extract the non-log Bayes factors; see examples), that prints nicely.
Interpreting Bayes Factors

A Bayes factor greater than 1 can be interpreted as evidence against the null, at which one convention is that a Bayes factor greater than 3 can be considered as "substantial" evidence against the null (and vice versa, a Bayes factor smaller than 1/3 indicates substantial evidence in favor of the null-model) (Wetzels et al. 2011).

Note

There is also a plot() method implemented in the see-package.

Author(s)

Mattan S. Ben-Shachar

References


Examples

```r
# With lm objects:
# ----------------
lm1 <- lm(mpg ~ 1, data = mtcars)
lm2 <- lm(mpg ~ hp, data = mtcars)
lm3 <- lm(mpg ~ hp + drat, data = mtcars)
lm4 <- lm(mpg ~ hp * drat, data = mtcars)

(BFM <- bayesfactor_models(lm1, lm2, lm3, lm4, denominator = 1))
# bayesfactor_models(lm2, lm3, lm4, denominator = lm1) # same result
# bayesfactor_models(lm1, lm2, lm3, lm4, denominator = lm1) # same result

update(BFM, reference = "bottom")
as.matrix(BFM)
as.numeric(BFM)

lm2b <- lm(sqrt(mpg) ~ hp, data = mtcars)
# Set check_response = TRUE for transformed responses
bayesfactor_models(lm2b, denominator = lm2, check_response = TRUE)
```
## Not run:
# With lmerMod objects:
# ---------------------
if (require("lme4")) {
  lmer1 <- lmer(Sepal.Length ~ Petal.Length + (1 | Species), data = iris)
  lmer2 <- lmer(Sepal.Length ~ Petal.Length + (Petal.Length | Species), data = iris)
  lmer3 <- lmer(Sepal.Length ~ Petal.Length + (Petal.Length | Species) + (1 | Petal.Width),
                data = iris)
  bayesfactor_models(lmer1, lmer2, lmer3,
                     denominator = 1,
                     estimator = "REML")
}

# rstanarm models
# ---------------------
# (note that a unique diagnostic_file MUST be specified in order to work)
if (require("rstanarm")) {
  stan_m0 <- stan_glm(Sepal.Length ~ 1,
                       data = iris,
                       family = gaussian(),
                       diagnostic_file = file.path(tempdir(), "df0.csv")
  )
  stan_m1 <- stan_glm(Sepal.Length ~ Species,
                      data = iris,
                      family = gaussian(),
                      diagnostic_file = file.path(tempdir(), "df1.csv")
  )
  stan_m2 <- stan_glm(Sepal.Length ~ Species + Petal.Length,
                      data = iris,
                      family = gaussian(),
                      diagnostic_file = file.path(tempdir(), "df2.csv")
  )
  bayesfactor_models(stan_m1, stan_m2, denominator = stan_m0)
}

# brms models
# ---------------------
# (note the save_pars MUST be set to save_pars(all = TRUE) in order to work)
if (require("brms")) {
  brm1 <- brm(Sepal.Length ~ 1, data = iris, save_all_pars = TRUE)
  brm2 <- brm(Sepal.Length ~ Species, data = iris, save_all_pars = TRUE)
  brm3 <- brm(
             Sepal.Length ~ Species + Petal.Length,
             data = iris,
             save_pars = save_pars(all = TRUE)
             )
  bayesfactor_models(brm1, brm2, brm3, denominator = 1)
# BayesFactor
# ---------------------------
if (require("BayesFactor")) {
  data(puzzles)
  BF <- anovaBF(RT ~ shape * color + ID,
                data = puzzles,
                whichRandom = "ID", progress = FALSE)
  BF
  bayesfactor_models(BF) # basically the same
}

## End(Not run)

## Bayes Factor for a Single Parameter

### Description

This method computes Bayes factors against the null (either a point or an interval), based on prior and posterior samples of a single parameter. This Bayes factor indicates the degree by which the mass of the posterior distribution has shifted further away from or closer to the null value(s) (relative to the prior distribution), thus indicating if the null value has become less or more likely given the observed data.

When the null is an interval, the Bayes factor is computed by comparing the prior and posterior odds of the parameter falling within or outside the null interval (Morey & Rouder, 2011; Liao et al., 2020); When the null is a point, a Savage-Dickey density ratio is computed, which is also an approximation of a Bayes factor comparing the marginal likelihoods of the model against a model in which the tested parameter has been restricted to the point null (Wagenmakers et al., 2010; Heck, 2019).

Note that the logspline package is used for estimating densities and probabilities, and must be installed for the function to work.

bayesfactor_pointnull() and bayesfactor_rope() are wrappers around bayesfactor_parameters with different defaults for the null to be tested against (a point and a range, respectively). Aliases of the main functions are prefixed with bf_*, like bf_parameters() or bf_pointnull().

**For more info, in particular on specifying correct priors for factors with more than 2 levels, see the Bayes factors vignette.**

### Usage

bayesfactor_parameters(}

## Bayes Factors (BF) for a Single Parameter

Bayes Factors (BF) for a Single Parameter

Description

This method computes Bayes factors against the null (either a point or an interval), based on prior and posterior samples of a single parameter. This Bayes factor indicates the degree by which the mass of the posterior distribution has shifted further away from or closer to the null value(s) (relative to the prior distribution), thus indicating if the null value has become less or more likely given the observed data.

When the null is an interval, the Bayes factor is computed by comparing the prior and posterior odds of the parameter falling within or outside the null interval (Morey & Rouder, 2011; Liao et al., 2020); When the null is a point, a Savage-Dickey density ratio is computed, which is also an approximation of a Bayes factor comparing the marginal likelihoods of the model against a model in which the tested parameter has been restricted to the point null (Wagenmakers et al., 2010; Heck, 2019).

Note that the logspline package is used for estimating densities and probabilities, and must be installed for the function to work.

bayesfactor_pointnull() and bayesfactor_rope() are wrappers around bayesfactor_parameters with different defaults for the null to be tested against (a point and a range, respectively). Aliases of the main functions are prefixed with bf_*, like bf_parameters() or bf_pointnull().

**For more info, in particular on specifying correct priors for factors with more than 2 levels, see the Bayes factors vignette.**

Usage

bayesfactor_parameters(}
bayesfactor_parameters

posterior,
prior = NULL,
direction = "two-sided",
null = 0,
verbose = TRUE,
...
)

bayesfactor_pointnull(
    posterior,
prior = NULL,
direction = "two-sided",
null = 0,
verbose = TRUE,
...
)

bayesfactor_rope(
    posterior,
prior = NULL,
direction = "two-sided",
null = rope_range(posterior),
verbose = TRUE,
...
)

bf_parameters(
    posterior,
prior = NULL,
direction = "two-sided",
null = 0,
verbose = TRUE,
...
)

bf_pointnull(
    posterior,
prior = NULL,
direction = "two-sided",
null = 0,
verbose = TRUE,
...
)

bf_rope(
    posterior,
prior = NULL,
direction = "two-sided",

null = rope_range(posterior),
verbose = TRUE,
...
)

## S3 method for class 'numeric'
bayesfactor_parameters(
  posterior,
  prior = NULL,
  direction = "two-sided",
  null = 0,
  verbose = TRUE,
  ...
)

## S3 method for class 'stanreg'
bayesfactor_parameters(
  posterior,
  prior = NULL,
  direction = "two-sided",
  null = 0,
  verbose = TRUE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "location", "smooth_terms", "sigma", "zi",
                "zero_inflated", "all"),
  parameters = NULL,
  ...
)

## S3 method for class 'brmsfit'
bayesfactor_parameters(
  posterior,
  prior = NULL,
  direction = "two-sided",
  null = 0,
  verbose = TRUE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "location", "smooth_terms", "sigma", "zi",
                "zero_inflated", "all"),
  parameters = NULL,
  ...
)

## S3 method for class 'blavaan'
bayesfactor_parameters(
  posterior,
  prior = NULL,
  direction = "two-sided",
  null = 0,
  verbose = TRUE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "location", "smooth_terms", "sigma", "zi",
                "zero_inflated", "all"),
  parameters = NULL,
  ...
)
null = 0,
verbose = TRUE,
...
)

## S3 method for class 'data.frame'
bayesfactor_parameters(
  posterior,
  prior = NULL,
  direction = "two-sided",
  null = 0,
  verbose = TRUE,
  ...
)

Arguments

- **posterior**: A numerical vector, stanreg / brmsfit object, emmGrid or a data frame - representing a posterior distribution(s) from (see 'Details').
- **prior**: An object representing a prior distribution (see 'Details').
- **direction**: Test type (see 'Details'). One of 0, "two-sided" (default, two tailed), -1, "left" (left tailed) or 1, "right" (right tailed).
- **null**: Value of the null, either a scalar (for point-null) or a range (for a interval-null).
- **verbose**: Toggle off warnings.
- **...**: Arguments passed to and from other methods. (Can be used to pass arguments to internal \texttt{logspline::logspline}().)
- **effects**: Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
- **component**: Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to \texttt{brms}-models.
- **parameters**: Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like \texttt{lp__} or \texttt{prior__}) are filtered by default, so only parameters that typically appear in the \texttt{summary()} are returned. Use \texttt{parameters} to select specific parameters for the output.

Details

This method is used to compute Bayes factors based on prior and posterior distributions.

**One-sided & Dividing Tests (setting an order restriction)**: One sided tests (controlled by direction) are conducted by restricting the prior and posterior of the non-null values (the "alternative") to one side of the null only (Morey & Wagenmakers, 2014). For example, if we have a prior hypothesis that the parameter should be positive, the alternative will be restricted to the region to the right of the null (point or interval). For example, for a Bayes factor comparing the "null" of $0-0.1$ to the alternative $>0.1$, we would set \texttt{bayesfactor_parameters(null = c(0, 0.1), direction = ">")}. 
It is also possible to compute a Bayes factor for dividing hypotheses - that is, for a null and alternative that are complementary, opposing one-sided hypotheses (Morey & Wagenmakers, 2014). For example, for a Bayes factor comparing the "null" of <0 to the alternative >0, we would set `bayesfactor_parameters(null = c(-Inf, 0))`.

**Value**

A data frame containing the (log) Bayes factor representing evidence against the null (Use `as.numeric()` to extract the non-log Bayes factors; see examples).

**Setting the correct prior**

For the computation of Bayes factors, the model priors must be proper priors (at the very least they should be not flat, and it is preferable that they be informative); As the priors for the alternative get wider, the likelihood of the null value(s) increases, to the extreme that for completely flat priors the null is infinitely more favorable than the alternative (this is called the Jeffreys-Lindley-Bartlett paradox). Thus, you should only ever try (or want) to compute a Bayes factor when you have an informed prior.

(Note that by default, `brms::brm()` uses flat priors for fixed-effects; See example below.)

It is important to provide the correct prior for meaningful results.

- When posterior is a numerical vector, prior should also be a numerical vector.
- When posterior is a data.frame, prior should also be a data.frame, with matching column order.
- When posterior is a stanreg, brmsfit or other supported Bayesian model:
  - prior can be set to NULL, in which case prior samples are drawn internally.
  - prior can also be a model equivalent to posterior but with samples from the priors only. See `unupdate()`.
  - **Note:** When posterior is a brmsfit_multiple model, prior must be provided.
- When posterior is an emmGrid / emm_list object:
  - prior should also be an emmGrid / emm_list object equivalent to posterior but created with a model of priors samples only. See `unupdate()`.
  - prior can also be the original (posterior) model. If so, the function will try to update the emmGrid / emm_list to use the `unupdate()`d prior-model. **(This cannot be done for brmsfit models.)**
  - **Note:** When the emmGrid has undergone any transformations ("log", "response", etc.), or regridding, then prior must be an emmGrid object, as stated above.

**Interpreting Bayes Factors**

A Bayes factor greater than 1 can be interpreted as evidence against the null, at which one convention is that a Bayes factor greater than 3 can be considered as "substantial" evidence against the null (and vice versa, a Bayes factor smaller than 1/3 indicates substantial evidence in favor of the null-model) (Wetzels et al. 2011).
Note
There is also a `plot()`-method implemented in the `see-package`.

Author(s)
Mattan S. Ben-Shachar

References


Examples

```r
library(bayestestR)
if (require("logspline")) {
  prior <- distribution_normal(1000, mean = 0, sd = 1)
  posterior <- distribution_normal(1000, mean = .5, sd = .3)
  (BF_pars <- bayesfactor_parameters(posterior, prior))
  as.numeric(BF_pars)
}
## Not run:
# rstanarm models
# -----------------
if (require("rstanarm") & & require("emmeans") & & require("logspline")) {
  contrasts(sleep$group) <- contr.orthonorm # see vignette
  stan_model <- stan_lmer(extra ~ group + (1 | ID), data = sleep)
  bayesfactor_parameters(stan_model)
  bayesfactor_parameters(stan_model, null = rope_range(stan_model))
}
# emmGrid objects
# --------------
  group_diff <- pairs(emmeans(stan_model, ~group))
  bayesfactor_parameters(group_diff, prior = stan_model)
```
# Or
  group_diff_prior <- pairs(emmeans(unupdate(stan_model), ~group))
  bayesfactor_parameters(group_diff, prior = group_diff_prior)
}

# brms models
# ------------
if (require("brms")) {
  contrasts(sleep$group) <- contr.orthonorm # see vignette
  my_custom_priors <-
    set_prior("student_t(3, 0, 1)", class = "b") +
    set_prior("student_t(3, 0, 1)", class = "sd", group = "ID")
  brms_model <- brm(extra ~ group + (1 | ID),
    data = sleep,
    prior = my_custom_priors
  )
  bayesfactor_parameters(brms_model)
}
## End(Not run)

---

**bayesfactor_restricted**

*Bayes Factors (BF) for Order Restricted Models*

---

**Description**

This method computes Bayes factors for comparing a model with an order restrictions on its parameters with the fully unrestricted model. *Note that this method should only be used for confirmatory analyses.*

The `bf_*` function is an alias of the main function.

For more info, in particular on specifying correct priors for factors with more than 2 levels, see the Bayes factors vignette.

**Usage**

```r
bayesfactor_restricted(
  posterior,
  hypothesis,
  prior = NULL,
  verbose = TRUE,
  ...
)
```

`bf_restricted(posterior, hypothesis, prior = NULL, verbose = TRUE, ...)"
## S3 method for class 'stanreg'
bayesfactor_restricted(
    posterior,
    hypothesis,
    prior = NULL,
    verbose = TRUE,
    effects = c("fixed", "random", "all"),
    component = c("conditional", "zi", "zero_inflated", "all"),
    ...
)

## S3 method for class 'brmsfit'
bayesfactor_restricted(
    posterior,
    hypothesis,
    prior = NULL,
    verbose = TRUE,
    effects = c("fixed", "random", "all"),
    component = c("conditional", "zi", "zero_inflated", "all"),
    ...
)

## S3 method for class 'blavaan'
bayesfactor_restricted(
    posterior,
    hypothesis,
    prior = NULL,
    verbose = TRUE,
    ...
)

## S3 method for class 'emmGrid'
bayesfactor_restricted(
    posterior,
    hypothesis,
    prior = NULL,
    verbose = TRUE,
    ...
)

### Arguments

- **posterior**: A `stanreg`/`brmsfit` object, `emmGrid` or a data frame - representing a posterior distribution(s) from (see Details).
- **hypothesis**: A character vector specifying the restrictions as logical conditions (see examples below).
- **prior**: An object representing a prior distribution (see Details).
verbose          Toggle off warnings.
... Currenty not used.
effects          Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
component        Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to **brms**-models.

**Details**

This method is used to compute Bayes factors for order-restricted models vs un-restricted models by setting an order restriction on the prior and posterior distributions (*Morey & Wagenmakers, 2013*).

(Though it is possible to use `bayesfactor_restricted()` to test interval restrictions, it is more suitable for testing order restrictions; see examples).

**Value**

A data frame containing the (log) Bayes factor representing evidence against the un-restricted model (Use `as.numeric()` to extract the non-log Bayes factors; see examples). (A bool_results attribute contains the results for each sample, indicating if they are included or not in the hypothesized restriction.)

**Setting the correct prior**

For the computation of Bayes factors, the model priors must be proper priors (at the very least they should be not flat, and it is preferable that they be informative); As the priors for the alternative get wider, the likelihood of the null value(s) increases, to the extreme that for completely flat priors the null is infinitely more favorable than the alternative (this is called the Jeffreys-Lindley-Bartlett paradox). Thus, you should only ever try (or want) to compute a Bayes factor when you have an informed prior.

(Note that by default, `brms::brm()` uses flat priors for fixed-effects; See example below.)

It is important to provide the correct prior for meaningful results.

- When posterior is a numerical vector, prior should also be a numerical vector.
- When posterior is a `data.frame`, prior should also be a `data.frame`, with matching column order.
- When posterior is a `stanreg`, `brmsfit` or other supported Bayesian model:
  - prior can be set to `NULL`, in which case prior samples are drawn internally.
  - prior can also be a model equivalent to posterior but with samples from the priors only. See `unupdate()`.
  - **Note:** When posterior is a `brmsfit_multiple` model, prior **must** be provided.
- When posterior is an `emmGrid` / `emm_list` object:
  - prior should also be an `emmGrid` / `emm_list` object equivalent to posterior but created with a model of priors samples only. See `unupdate()`.
**Interpreting Bayes Factors**

A Bayes factor greater than 1 can be interpreted as evidence against the null, at which one convention is that a Bayes factor greater than 3 can be considered as "substantial" evidence against the null (and vice versa, a Bayes factor smaller than 1/3 indicates substantial evidence in favor of the null-model) (Wetzels et al. 2011).

**References**


**Examples**

```r
set.seed(444)
library(bayestestR)
prior <- data.frame(
  A = rnorm(1000),
  B = rnorm(1000),
  C = rnorm(1000)
)

posterior <- data.frame(
  A = rnorm(1000, .4, 0.7),
  B = rnorm(1000, -.2, 0.4),
  C = rnorm(1000, 0, 0.5)
)

hyp <- c(
  "A > B & B > C",
  "A > B & A > C",
  "C > A"
)

if (getRversion() > "3.5.0") {
  b <- bayesfactor_restricted(posterior, hypothesis = hyp, prior = prior))
  as.numeric(b)
  if (require("see") && require("patchwork")) {
    i <- attr(b, "bool_results")[["posterior"]]
```
see::plots(
  plot(estimate_density(posterior)),
  # distribution **conditional** on the restrictions
  plot(estimate_density(posterior[[hyps[1]]], )) + ggplot2::ggtitle(hyps[1]),
  plot(estimate_density(posterior[[hyps[2]]], )) + ggplot2::ggtitle(hyps[2]),
  plot(estimate_density(posterior[[hyps[3]]], )) + ggplot2::ggtitle(hyps[3]),
  guides = "collect"
)
)

## Not run:
# rstanarm models
# ---------------
if (require("rstanarm") && require("emmeans")) {
  fit_stan <- stan_glm(mpg ~ wt + cyl + am,
    data = mtcars, refresh = 0
  )
  hyps <- c("am > 0 & cyl < 0", "cyl < 0", "wt - cyl > 0"
  )
  bayesfactor_restricted(fit_stan, hypothesis = hyps)

  # emmGrid objects
  # ---------------
  disgust_data <- read.table(url("http://www.learnbayes.org/disgust_example.txt"), header = TRUE)
  contrasts(disgust_data$condition) <- contr.orthonorm # see vignette
  fit_model <- stan_glm(score ~ condition, data = disgust_data, family = gaussian())
  em_condition <- emmeans(fit_model, ~condition)
  hyps <- c("lemon < control & control < sulfur")
  bayesfactor_restricted(em_condition, prior = fit_model, hypothesis = hyps)
  # > # Bayes Factor (Order-Restriction)
  # > # Hypothesis P(Prior) P(Posterior) BF
  # > lemon < control & control < sulfur 0.17 0.75 4.49
  # > ---
  # > Bayes factors for the restricted model vs. the un-restricted model.
}

## End(Not run)

---

Bias Corrected and Accelerated Interval (BCa)
Description

Compute the Bias Corrected and Accelerated Interval (BCa) of posterior distributions.

Usage

bci(x, ...)

bcai(x, ...)

## S3 method for class 'numeric'
bci(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'data.frame'
bci(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'MCMCglmm'
bci(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'sim.merMod'
bci(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  parameters = NULL,
  verbose = TRUE,
  ...
)

## S3 method for class 'sim'
bci(x, ci = 0.95, parameters = NULL, verbose = TRUE, ...)

## S3 method for class 'emmGrid'
bci(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'stanreg'
bci(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                "distributional", "auxiliary"),
  parameters = NULL,
  verbose = TRUE,
  ...
)

## S3 method for class 'brmsfit'
bci(

\begin{verbatim}
x,  
ci = 0.95,  
effects = c("fixed", "random", "all"),  
component = c("conditional", "zi", "zero_inflated", "all"),  
parameters = NULL,  
verbose = TRUE,  
...
)

## S3 method for class 'BFBayesFactor'

bci(x, ci = 0.95, verbose = TRUE, ...)
\end{verbatim}

**Arguments**

- **x**: Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. bayestestR supports a wide range of models (see, for example, methods("hdi")) and not all of those are documented in the 'Usage' section, because methods for other classes mostly resemble the arguments of the .numeric or .data.frame methods.

- **...**: Currently not used.

- **ci**: Value or vector of probability of the (credible) interval - CI (between 0 and 1) to be estimated. Default to .95 (95%).

- **verbose**: Toggle off warnings.

- **effects**: Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

- **parameters**: Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior__) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

- **component**: Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

**Details**

Unlike equal-tailed intervals (see eti()) that typically exclude 2.5% from each tail of the distribution and always include the median, the HDI is not equal-tailed and therefore always includes the mode(s) of posterior distributions. While this can be useful to better represent the credibility mass of a distribution, the HDI also has some limitations. See spi() for details.

The 95% or 89% Credible Intervals (CI) are two reasonable ranges to characterize the uncertainty related to the estimation (see here for a discussion about the differences between these two values). The 89% intervals (ci = 0.89) are deemed to be more stable than, for instance, 95% intervals (Kruschke, 2014). An effective sample size of at least 10.000 is recommended if one wants to estimate 95% intervals with high precision (Kruschke, 2014, p. 183ff). Unfortunately, the default number of posterior samples for most Bayes packages (e.g., rstanarm or brms) is only 4.000 (thus, you might want to increase it when fitting your model). Moreover, 89 indicates the arbitrariness of interval
limits - its only remarkable property is being the highest prime number that does not exceed the already unstable 95% threshold (McElreath, 2015). However, 95% has some advantages too. For instance, it shares (in the case of a normal posterior distribution) an intuitive relationship with the standard deviation and it conveys a more accurate image of the (artificial) bounds of the distribution. Also, because it is wider, it makes analyses more conservative (i.e., the probability of covering 0 is larger for the 95% CI than for lower ranges such as 89%), which is a good thing in the context of the reproducibility crisis.

A 95% equal-tailed interval (ETI) has 2.5% of the distribution on either side of its limits. It indicates the 2.5th percentile and the 97.5th percentile. In symmetric distributions, the two methods of computing credible intervals, the ETI and the HDI, return similar results.

This is not the case for skewed distributions. Indeed, it is possible that parameter values in the ETI have lower credibility (are less probable) than parameter values outside the ETI. This property seems undesirable as a summary of the credible values in a distribution.

On the other hand, the ETI range does change when transformations are applied to the distribution (for instance, for a log odds scale to probabilities): the lower and higher bounds of the transformed distribution will correspond to the transformed lower and higher bounds of the original distribution. On the contrary, applying transformations to the distribution will change the resulting HDI.

Value

A data frame with following columns:

- **Parameter** The model parameter(s), if x is a model-object. If x is a vector, this column is missing.
- **CI** The probability of the credible interval.
- **CI_low, CI_high** The lower and upper credible interval limits for the parameters.

References


See Also

Other ci: `ci()`, `cwi()`, `eti()`, `hdi()`, `si()`, `spi()`

Examples

```r
posterior <- rnorm(1000)
bci(posterior)
bci(posterior, ci = c(.80, .89, .95))
```
bic_to_bf

Convert BIC indices to Bayes Factors via the BIC-approximation method.

Description

The difference between two Bayesian information criterion (BIC) indices of two models can be used to approximate Bayes factors via:

\[ BF_{10} = e^{(BIC_0 - BIC_1)/2} \]

Usage

bic_to_bf(bic, denominator, log = FALSE)

Arguments

bic A vector of BIC values.
denominator The BIC value to use as a denominator (to test against).
log If TRUE, return the log(BF).

Value

The Bayes Factors corresponding to the BIC values against the denominator.

References


Examples

bic1 <- BIC(lm(Sepal.Length ~ 1, data = iris))
bic2 <- BIC(lm(Sepal.Length ~ Species, data = iris))
bic3 <- BIC(lm(Sepal.Length ~ Species + Petal.Length, data = iris))
bic4 <- BIC(lm(Sepal.Length ~ Species * Petal.Length, data = iris))

bic_to_bf(c(bic1, bic2, bic3, bic4), denominator = bic1)
check_prior  

Check if Prior is Informative

Description

Performs a simple test to check whether the prior is informative to the posterior. This idea, and the accompanying heuristics, were discussed in this blogpost.

Usage

```r
check_prior(model, method = "gelman", simulate_priors = TRUE, ...)
```

Arguments

- **model**: A `stanreg`, `stanfit`, `brmsfit`, `blavaan`, or `MCMCglmm` object.
- **method**: Can be "gelman" or "lakeland". For the "gelman" method, if the SD of the posterior is more than 0.1 times the SD of the prior, then the prior is considered as informative. For the "lakeland" method, the prior is considered as informative if the posterior falls within the 95% HDI of the prior.
- **simulate_priors**: Should prior distributions be simulated using `simulate_prior()` (default; faster) or sampled via `unupdate()` (slower, more accurate).
- **...**: Currently not used.

Value

A data frame with two columns: The parameter names and the quality of the prior (which might be "informative", "uninformative") or "not determinable" if the prior distribution could not be determined.

References

https://statmodeling.stat.columbia.edu/2019/08/10/

Examples

```r
## Not run:
library(bayestestR)
if (require("rstanarm")) {
  model <- stan_glm(mpg ~ wt + am, data = mtcars, chains = 1, refresh = 0)
  check_prior(model, method = "gelman")
  check_prior(model, method = "lakeland")

  # An extreme example where both methods diverge:
  model <- stan_glm(mpg ~ wt,
                   data = mtcars[1:3, ],
                   prior = normal(-3.3, 1, FALSE),
                   prior_intercept = normal(0, 1000, FALSE),
```
refresh = 0
)
check_prior(model, method = "gelman")
check_prior(model, method = "lakeland")
plot(si(model)) # can provide visual confirmation to the Lakeland method
}

## End(Not run)

---

ci

Confidence/Credible/Compatibility Interval (CI)

Description

Compute Confidence/Credible/Compatibility Intervals (CI) or Support Intervals (SI) for Bayesian and frequentist models. The Documentation is accessible for:

Usage

`ci(x, ...)`

## S3 method for class 'numeric'
`ci(x, ci = 0.95, method = "ETI", verbose = TRUE, BF = 1, ...)`

## S3 method for class 'data.frame'
`ci(x, ci = 0.95, method = "ETI", verbose = TRUE, BF = 1, ...)`

## S3 method for class 'sim.merMod'
`ci(`
  `x,`
  `ci = 0.95,`
  `method = "ETI",`
  `effects = c("fixed", "random", "all"),`
  `parameters = NULL,`
  `verbose = TRUE,`
  `...`
`)

## S3 method for class 'sim'
`ci(x, ci = 0.95, method = "ETI", parameters = NULL, verbose = TRUE, ...)`

## S3 method for class 'stanreg'
`ci(`
  `x,`
  `ci = 0.95,`
  `method = "ETI",`
  `effects = c("fixed", "random", "all"),`
  `component = c("location", "all", "conditional", "smooth_terms", "sigma",`
Arguments

x A stanreg or brmsfit model, or a vector representing a posterior distribution.

... Currently not used.

ci Value or vector of probability of the CI (between 0 and 1) to be estimated. Default to 0.95 (95%).

method Can be 'ETI' (default), 'HDI', 'BCI', 'SPI' or 'SI'.

verbose Toggle off warnings.

BF The amount of support required to be included in the support interval.

effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior_) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.
Details

- Bayesian models
- Frequentist models

Value

A data frame with following columns:

- Parameter The model parameter(s), if x is a model-object. If x is a vector, this column is missing.
- CI The probability of the credible interval.
- CI_low, CI_high The lower and upper credible interval limits for the parameters.

Note

When it comes to interpretation, we recommend thinking of the CI in terms of an "uncertainty" or "compatibility" interval, the latter being defined as “Given any value in the interval and the background assumptions, the data should not seem very surprising” (Gelman & Greenland 2019).

There is also a plot()-method implemented in the see-package.

References

Gelman A, Greenland S. Are confidence intervals better termed "uncertainty intervals"? BMJ 2019;1:5381. 10.1136/bmj.l5381

See Also

Other ci: bci(), ci(), eti(), hdi(), si(), spi()

Examples

library(bayestestR)

posterior <- rnorm(1000)
ci(posterior, method = "ETI")
ci(posterior, method = "HDI")

df <- data.frame(replicate(4, rnorm(100)))
ci(df, method = "ETI", ci = c(.80, .89, .95))
ci(df, method = "HDI", ci = c(.80, .89, .95))
## Not run:
if (require("rstanarm")) {
  model <- stan_glm(mpg ~ wt, data = mtcars, chains = 2, iter = 200, refresh = 0)
  ci(model, method = "ETI", ci = c(.80, .89))
  ci(model, method = "HDI", ci = c(.80, .89))
  ci(model, method = "SI")
}
if (require("brms")) {
model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
ci(model, method = "ETI")
ci(model, method = "HDI")
ci(model, method = "SI")
}

if (require("BayesFactor")) {
  bf <- ttestBF(x = rnorm(100, 1, 1))
ci(bf, method = "ETI")
ci(bf, method = "HDI")
}

if (require("emmeans")) {
  model <- emtrends(model, ~1, "wt")
ci(model, method = "ETI")
ci(model, method = "HDI")
ci(model, method = "SI")
}

## End(Not run)

---

### contr.orthonorm

**Orthonormal Contrast Matrices for Bayesian Estimation**

**Description**

Returns a design or model matrix of orthonormal contrasts such that the marginal prior on all effects is identical (see 'Details'). Implementation from Singmann & Gronau’s `bfrms`, following the description in Rouder, Morey, Speckman, & Province (2012, p. 363).

Though using this factor coding scheme might obscure the interpretation of parameters, it is essential for correct estimation of Bayes factors for contrasts and order restrictions of multi-level factors (where k>2). See info on specifying correct priors for factors with more than 2 levels in the Bayes factors vignette.

**Usage**

```r
contr.orthonorm(n, contrasts = TRUE, sparse = FALSE)
```

**Arguments**

- `n`: a vector of levels for a factor, or the number of levels.
- `contrasts`: a logical indicating whether contrasts should be computed.
- `sparse`: logical indicating if the result should be sparse (of class `dgCMatrix`), using package `Matrix`. 
Details

When contrasts = FALSE, the returned contrasts are equivalent to contr.treatment(, contrasts = FALSE), as suggested by McElreath (also known as one-hot encoding).

**Setting Priors:**
It is recommended to set 0-centered, identically-scaled priors on the dummy coded variables produced by this method. These priors then represent the distance the mean of one of the levels might have from the overall mean.

**Contrasts:**
This method guarantees that any set of contrasts between the k groups will have the same multivariate prior regardless of level order; However, different contrasts within a set contrasts can have different univariate prior shapes/scales.

For example the contrasts A - B will have the same prior as B - C, as will (A + C) - B and (B + A) - C, but A - B and (A + C) - B will differ.

Value

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

References


Examples

```r
contr.orthonorm(2) # Q_2 in Rouder et al. (2012, p. 363)
contr.orthonorm(5) # equivalent to Q_5 in Rouder et al. (2012, p. 363)
```

```r
## check decomposition
Q3 <- contr.orthonorm(3)
Q3 %*% t(Q3) ## 2/3 on diagonal and -1/3 on off-diagonal elements
```

**Description**

Refit Bayesian model as frequentist. Can be useful for comparisons.
Usage

convert_bayesian_as_frequentist(model, data = NULL, REML = TRUE)

bayesian_as_frequentist(model, data = NULL, REML = TRUE)

Arguments

model A Bayesian model.

data Data used by the model. If NULL, will try to extract it from the model.

REML For mixed effects, should models be estimated using restricted maximum likelihood (REML) (TRUE, default) or maximum likelihood (FALSE)?)

Examples

# Rstanarm ----------------------
if (require("rstanarm")) {
    # Simple regressions
    model <- stan_glm(Sepal.Length ~ Species,
                      data = iris, chains = 2, refresh = 0
        )
    bayesian_as_frequentist(model)
}

## Not run:
if (require("rstanarm")) {
    model <- stan_glm(vs ~ mpg,
                      family = "binomial",
                      data = mtcars, chains = 2, refresh = 0
        )
    bayesian_as_frequentist(model)
    # Mixed models
    model <- stan_glmer(Sepal.Length ~ Petal.Length + (1 | Species),
                        data = iris, chains = 2, refresh = 0
        )
    bayesian_as_frequentist(model)
    model <- stan_glmer(vs ~ mpg + (1 | cyl),
                        family = "binomial",
                        data = mtcars, chains = 2, refresh = 0
        )
    bayesian_as_frequentist(model)
}

## End(Not run)
**Description**

Compute the **Curvewise interval (CWI)** (also called the "simultaneous interval" or "joint interval") of posterior distributions using `ggdist::curve_interval()`. Whereas the more typical "pointwise intervals" contain xx% of the posterior for a single parameter, joint/curvewise intervals contain xx% of the posterior distribution for all parameters.

**Usage**

cwi(x, ...)

```r
## S3 method for class 'data.frame'
cwi(x, ci = 0.95, ...)
```

**Arguments**

- `x` Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. `bayestestR` supports a wide range of models (see, for example, `methods("hdi")`) and not all of those are documented in the 'Usage' section, because methods for other classes mostly resemble the arguments of the `.numeric` or `.data.frame` methods.
- `...` Currently not used.
- `ci` Value or vector of probability of the (credible) interval - CI (between 0 and 1) to be estimated. Default to .95 (95%).

**Details**

Applied model predictions, pointwise intervals contain xx% of the predicted response values conditional on specific predictor values. In contrast, curvewise intervals contain xx% of the predicted response values across all predictor values. Put another way, curvewise intervals contain xx% of the full prediction lines from the model.

For more details, see the `ggdist` documentation on curvewise intervals.

**Value**

A data frame with following columns:

- **Parameter** The model parameter(s), if x is a model-object. If x is a vector, this column is missing.
- **CI** The probability of the credible interval.
- **CI_low, CI_high** The lower and upper credible interval limits for the parameters.
density_at

Density Probability at a Given Value

See Also

Other ci: bci(), ci(), eti(), hdi(), si(), spi()

Examples

library(bayestestR)

if (require("ggplot2") && require("rstanarm") && require("ggdist")) {

  # Generate data =============================================
  k <- 11 # number of curves (iterations)
  n <- 201 # number of rows
  data <- data.frame(x = seq(-15, 15, length.out = n))

  # Simulate iterations as new columns
  for (i in 1:k) {
    data[paste0("iter_", i)] <- dnorm(data$x, seq(-5, 5, length.out = k)[i], 3)
  }

  # Note: first, we need to transpose the data to have iters as rows
  iters <- datawizard::data_transpose(data[paste0("iter_", 1:k)])

  # Compute Median
  data$Median <- point_estimate(iters)["Median"]

  # Compute Credible Intervals ================================
  # Compute ETI (default type of CI)
  data[c("ETI_low", "ETI_high")]<- eti(iters, ci = 0.5)[c("CI_low", "CI_high")]

  # Compute CWI
  ggdist::curve_interval(reshape_iterations(data), iter_value .width = c(.5))

  # Visualization =============================================
  ggplot(data, aes(x = x, y = Median)) +
    geom_ribbon(aes(ymin = ETI_low, ymax = ETI_high), fill = "red", alpha = 0.3) +
    geom_line(size = 1) +
    geom_line(
      data = reshape_iterations(data),
      aes(y = iter_value, group = iter_group),
      alpha = 0.3
    )
}
describe_posterior

Description

Compute the density value at a given point of a distribution (i.e., the value of the y axis of a value x of a distribution).

Usage

density_at(posterior, x, precision = 2^10, method = "kernel", ...)

Arguments

- posterior: Vector representing a posterior distribution.
- x: The value of which to get the approximate probability.
- precision: Number of points of density data. See the n parameter in density.
- method: Density estimation method. Can be "kernel" (default), "logspline" or "KernSmooth".
- ...: Currently not used.

Examples

library(bayestestR)
posterior <- distribution_normal(n = 10)
density_at(posterior, 0)
density_at(posterior, c(0, 1))

describe_posterior

Describe Posterior Distributions

Description

Compute indices relevant to describe and characterize the posterior distributions.

Usage

describe_posterior(posteriors, ...)

## S3 method for class 'numeric'
describe_posterior(
  posteriors,
  centrality = "median",
  dispersion = FALSE,
  ci = 0.95,
  ci_method = "eti",
  test = c("p_direction", "rope"),
  rope_range = "default",
  rope_ci = 0.95,
  keep_iterations = FALSE,
  bf_prior = NULL,
)
BF = 1,
...
)

## S3 method for class 'stanreg'
describe_posterior(
  posteriors,
  centrality = "median",
  dispersion = FALSE,
  ci = 0.95,
  ci_method = "eti",
  test = c("p_direction", "rope"),
  rope_range = "default",
  rope_ci = 0.95,
  keep_iterations = FALSE,
  bf_prior = NULL,
  diagnostic = c("ESS", "Rhat"),
  priors = FALSE,
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                    "distributional", "auxiliary"),
  parameters = NULL,
  BF = 1,
  ...
)

## S3 method for class 'brmsfit'
describe_posterior(
  posteriors,
  centrality = "median",
  dispersion = FALSE,
  ci = 0.95,
  ci_method = "eti",
  test = c("p_direction", "rope"),
  rope_range = "default",
  rope_ci = 0.95,
  keep_iterations = FALSE,
  bf_prior = NULL,
  diagnostic = c("ESS", "Rhat"),
  priors = FALSE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all", "location",
                    "distributional", "auxiliary"),
  parameters = NULL,
  BF = 1,
  priors = FALSE,
  ...
)
Arguments

posteriors  A vector, data frame or model of posterior draws. bayestestR supports a wide range of models (see methods("describe_posterior")) and not all of those are documented in the 'Usage' section, because methods for other classes mostly resemble the arguments of the .numeric method.

...  Additional arguments to be passed to or from methods.

centrality  The point-estimates (centrality indices) to compute. Character (vector) or list with one or more of these options: "median", "mean", "MAP" or "all".

dispersion  Logical, if TRUE, computes indices of dispersion related to the estimate(s) (SD and MAD for mean and median, respectively).

ci  Value or vector of probability of the CI (between 0 and 1) to be estimated. Default to .95 (95%).

ci_method  The type of index used for Credible Interval. Can be "ETI" (default, see eti()), "HDI" (see hdi()), "BCI" (see bci()), "SPI" (see spi()), or "SI" (see si()).

test  The indices of effect existence to compute. Character (vector) or list with one or more of these options: "p_direction" (or "pd"), "rope", "p_map", "equivalence_test" (or "equitest"), "bayesfactor" (or "bf") or "all" to compute all tests. For each "test", the corresponding bayestestR function is called (e.g. rope() or p_direction()) and its results included in the summary output.

rope_range  ROPE’s lower and higher bounds. Should be a list of two values (e.g., c(-0.1, 0.1)) or "default". If "default", the bounds are set to x ± 0.1*SD(response).

rope_ci  The Credible Interval (CI) probability, corresponding to the proportion of HDI, to use for the percentage in ROPE.

keep_iterations  If TRUE, will keep all iterations (draws) of bootstrapped or Bayesian models. They will be added as additional columns named iter_1, iter_2, .... You can reshape them to a long format by running reshape_iterations().

bf_prior  Distribution representing a prior for the computation of Bayes factors / SI. Used if the input is a posterior, otherwise (in the case of models) ignored.

BF  The amount of support required to be included in the support interval.

diagnostic  Diagnostic metrics to compute. Character (vector) or list with one or more of these options: "ESS", "Rhat", "MCSE" or "all".

priors  Add the prior used for each parameter.

effects  Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component  Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

parameters  Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior_) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.
Details

One or more components of point estimates (like posterior mean or median), intervals and tests can be omitted from the summary output by setting the related argument to NULL. For example, test = NULL and centrality = NULL would only return the HDI (or CI).

References

- Region of Practical Equivalence (ROPE)
- Bayes factors

Examples

```r
library(bayestestR)
if (require("logspline")) {
  x <- rnorm(1000)
  describe_posterior(x)
  describe_posterior(x, centrality = "all", dispersion = TRUE, test = "all")
  describe_posterior(x, ci = c(0.80, 0.90))

  df <- data.frame(replace(4, rnorm(100)))
  describe_posterior(df)
  describe_posterior(df, centrality = "all", dispersion = TRUE, test = "all")
  describe_posterior(df, ci = c(0.80, 0.90))

  df <- data.frame(replace(4, rnorm(20)))
  head(reshape_iterations(describe_posterior(df, keep_iterations = TRUE)))
}
## Not run:
# rstanarm models
# -----------------------------------------------
if (require("rstanarm") && require("emmeans")) {
  model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
  describe_posterior(model)
  describe_posterior(model, centrality = "all", dispersion = TRUE, test = "all")
  describe_posterior(model, ci = c(0.80, 0.90))

  # emmeans estimates
  # -----------------------------------------------
  describe_posterior(emtrends(model, ~1, "wt"))
}

# brms models
# -----------------------------------------------
if (require("brms")) {
  model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
  describe_posterior(model)
  describe_posterior(model, centrality = "all", dispersion = TRUE, test = "all")
}
describe_posterior(model, ci = c(0.80, 0.90))
}

# BayesFactor objects
# -----------------------------------------------
if (require("BayesFactor")) {
  bf <- ttestBF(x = rnorm(100, 1, 1))
  describe_posterior(bf)
  describe_posterior(bf, centrality = "all", dispersion = TRUE, test = "all")
  describe_posterior(bf, ci = c(0.80, 0.90))
}

## End(Not run)

---

### describe_prior

**Describe Priors**

Returns a summary of the priors used in the model.

#### Description

Returns a summary of the priors used in the model.

#### Usage

```r
describe_prior(model, 
```

```r
## S3 method for class 'brmsfit'
```

```r
describe_prior( 
  model, 
  effects = c("fixed", "random", "all"), 
  component = c("conditional", "zi", "zero_inflated", "all", "location", 
                 "distributional", "auxiliary"), 
  parameters = NULL,  
  ...
)
```

#### Arguments

- **model**: A Bayesian model.
- **effects**: Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
- **component**: Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to `brms`-models.
- **parameters**: Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like `lp__` or `prior__`) are filtered by default, so only parameters that typically appear in the `summary()` are returned. Use `parameters` to select specific parameters for the output.
Examples

```r
## Not run:
library(bayestestR)

# rstanarm models
# -----------------------------------------------
if (require("rstanarm")) {
  model <- rstanarm::stan_glm(mpg ~ wt + cyl, data = mtcars)
  describe_prior(model)
}

# brms models
# -----------------------------------------------
if (require("brms")) {
  model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
  describe_prior(model)
}

# BayesFactor objects
# -----------------------------------------------
if (require("BayesFactor")) {
  bf <- ttestBF(x = rnorm(100, 1, 1))
  describe_prior(bf)
}

## End(Not run)
```

diagnostic_draws

**Diagnostic values for each iteration**

### Description

Returns the accumulated log-posterior, the average Metropolis acceptance rate, divergent transitions, treedepth rather than terminated its evolution normally.

### Usage

```r
diagnostic_draws(posteriors, ...)
```

### Arguments

- `posteriors` A `stanreg`, `stanfit`, `brmsfit`, or `blavaan` object.
- `...` Currently not used.
Examples

```r
## Not run:
set.seed(333)

if (require("brms", quietly = TRUE)) {
  model <- brm(mpg ~ wt * cyl * vs,
               data = mtcars,
               iter = 100, control = list(adapt_delta = 0.80),
               refresh = 0
  )
  diagnostic_draws(model)
}

## End(Not run)
```

---

diagnostic_posterior  

Posterior Sampling Diagnostic

Description

Extract diagnostic metrics (Effective Sample Size (ESS), Rhat and Monte Carlo Standard Error MCSE).

Usage

```r
diagnostic_posterior(posteriors, diagnostic = c("ESS", "Rhat"), ...)
```

```r
## S3 method for class 'stanreg'
diagnostic_posterior(
  posteriors,
  diagnostic = "all",
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                "distributional", "auxiliary"),
  parameters = NULL,
  ...
)

## S3 method for class 'brmsfit'
diagnostic_posterior(
  posteriors,
  diagnostic = "all",
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL,
  ...
)
```
Arguments

- **posterior**: A `stanreg`, `stanfit`, `brmsfit`, or `blavaan` object.
- **diagnostic**: Diagnostic metrics to compute. Character (vector) or list with one or more of these options: "ESS", "Rhat", "MCSE" or "all".
- **...**: Currently not used.
- **effects**: Should parameters for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
- **component**: Which type of parameters to return, such as parameters for the conditional model, the zero-inflated part of the model, the dispersion term, the instrumental variables or marginal effects be returned? Applies to models with zero-inflated and/or dispersion formula, or to models with instrumental variables (so called fixed-effects regressions), or models with marginal effects from `mfx`. May be abbreviated. Note that the **conditional** component is also called **count or mean** component, depending on the model. There are three convenient shortcuts: component = "all" returns all possible parameters. If component = "location", location parameters such as conditional, zero_inflated, smooth_terms, or instruments are returned (everything that are fixed or random effects - depending on the effects argument - but no auxiliary parameters). For component = "distributional" (or "auxiliary"), components like sigma, dispersion, beta or precision (and other auxiliary parameters) are returned.
- **parameters**: Regular expression pattern that describes the parameters that should be returned.

Details

**Effective Sample (ESS)** should be as large as possible, although for most applications, an effective sample size greater than 1000 is sufficient for stable estimates (Bürkner, 2017). The ESS corresponds to the number of independent samples with the same estimation power as the N autocorrelated samples. It is is a measure of “how much independent information there is in autocorrelated chains” (Kruschke 2015, p182-3).

**Rhat** should be the closest to 1. It should not be larger than 1.1 (Gelman and Rubin, 1992) or 1.01 (Vehtari et al., 2019). The split Rhat statistic quantifies the consistency of an ensemble of Markov chains.

**Monte Carlo Standard Error (MCSE)** is another measure of accuracy of the chains. It is defined as standard deviation of the chains divided by their effective sample size (the formula for `mcse()` is from Kruschke 2015, p. 187). The MCSE “provides a quantitative suggestion of how big the estimation noise is”.

**References**


Examples

```r
## Not run:
# rstanarm models
# -----------------------------------------------
if (require("rstanarm", quietly = TRUE)) {
  model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
  diagnostic_posterior(model)
}

# brms models
# -----------------------------------------------
if (require("brms", quietly = TRUE)) {
  model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
  diagnostic_posterior(model)
}

## End(Not run)
```

distribution

**Empirical Distributions**

**Description**

Generate a sequence of n-quantiles, i.e., a sample of size n with a near-perfect distribution.

**Usage**

distribution(type = "normal", ...)
distribution_custom(n, type = "norm", ..., random = FALSE)
distribution_beta(n, shape1, shape2, ncp = 0, random = FALSE, ...)
distribution_binomial(n, size = 1, prob = 0.5, random = FALSE, ...)
distribution_binom(n, size = 1, prob = 0.5, random = FALSE, ...)
distribution_cauchy(n, location = 0, scale = 1, random = FALSE, ...)
distribution_chisquared(n, df, ncp = 0, random = FALSE, ...)
distribution_chisq(n, df, ncp = 0, random = FALSE, ...)
distribution_gamma(n, shape, scale = 1, random = FALSE, ...)
distribution_mixture_normal(n, mean = c(-3, 3), sd = 1, random = FALSE, ...)
distribution_normal(n, mean = 0, sd = 1, random = FALSE, ...)
distribution_gaussian(n, mean = 0, sd = 1, random = FALSE, ...)
distribution_nbinom(n, size, prob, mu, phi, random = FALSE, ...)
distribution_poisson(n, lambda = 1, random = FALSE, ...)
distribution_student(n, df, ncp, random = FALSE, ...)
distribution_t(n, df, ncp, random = FALSE, ...)
distribution_student_t(n, df, ncp, random = FALSE, ...)
distribution_tweedie(n, xi = NULL, mu, phi, power = NULL, random = FALSE, ...)
distribution_uniform(n, min = 0, max = 1, random = FALSE, ...)
rnorm_perfect(n, mean = 0, sd = 1)

Arguments

- type: Can be any of the names from base R’s Distributions, like “cauchy”, “pois” or “beta”.
- ...: Arguments passed to or from other methods.
- n: the number of observations.
- random: Generate near-perfect or random (simple wrappers for the base R r* functions) distributions.
- shape1: non-negative parameters of the Beta distribution.
- shape2: non-negative parameters of the Beta distribution.
- ncp: non-centrality parameter.
- size: number of trials (zero or more).
- prob: probability of success on each trial.
- location: location and scale parameters.
- scale: location and scale parameters.
- df: degrees of freedom (non-negative, but can be non-integer).
- shape: shape and scale parameters. Must be positive, scale strictly.
- mean: vector of means.
- sd: vector of standard deviations.
- mu: the mean
- phi: Corresponding to glmmTMB’s implementation of nbinom distribution, where size=mu/phi.
- lambda: vector of (non-negative) means.
For tweedie distributions, the value of \( x_i \) such that the variance is \( \text{var}(Y) = \phi \cdot \mu^{x_i} \).

**power**

Alias for \( x_i \).

**min**

lower and upper limits of the distribution. Must be finite.

**max**

lower and upper limits of the distribution. Must be finite.

Details

When \( \text{random} = \text{FALSE} \), these function return \( q*(\text{ppoints}(n), ...) \).

Examples

```r
library(bayestestR)
x <- distribution(n = 10)
plot(density(x))
x <- distribution(type = "gamma", n = 100, shape = 2)
plot(density(x))
```

---

**effective_sample**  
Effective Sample Size (ESS)

Description

This function returns the effective sample size (ESS).

Usage

```r
effective_sample(model, ...)
```

## S3 method for class 'brmsfit'
effective_sample(  
  model,  
  effects = c("fixed", "random", "all"),  
  component = c("conditional", "zi", "zero_inflated", "all"),  
  parameters = NULL,  
  ...
)

## S3 method for class 'stanreg'
effective_sample(  
  model,  
  effects = c("fixed", "random", "all"),  
  component = c("location", "all", "conditional", "smooth_terms", "sigma",  
    "distributional", "auxiliary"),  
  parameters = NULL,  
  ...
)
Arguments

model A stanreg, stanfit, brmsfit, blavaan, or MCMCglmm object.

... Currently not used.

effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior__) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

Details

**Effective Sample (ESS)** should be as large as possible, although for most applications, an effective sample size greater than 1,000 is sufficient for stable estimates (Bürkner, 2017). The ESS corresponds to the number of independent samples with the same estimation power as the N autocorrelated samples. It is a measure of “how much independent information there is in autocorrelated chains” (*Kruschke 2015, p182-3*).

Value

A data frame with two columns: Parameter name and effective sample size (ESS).

References


Examples

```r
## Not run:
library(rstanarm)
model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
effective_sample(model)
## End(Not run)
```
equivalence_test

Test for Practical Equivalence

Description

Perform a Test for Practical Equivalence for Bayesian and frequentist models.

Usage

equivalence_test(x, ...)

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

## S3 method for class 'data.frame'
equivalence_test(x, range = "default", ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'stanreg'
equivalence_test(
x, range = "default",
    ci = 0.95,
    effects = c("fixed", "random", "all"),
    component = c("location", "all", "conditional", "smooth_terms", "sigma",
        "distributional", "auxiliary"),
    parameters = NULL,
    verbose = TRUE,
    ...
)

## S3 method for class 'brmsfit'
equivalence_test(
x, range = "default",
    ci = 0.95,
    effects = c("fixed", "random", "all"),
    component = c("conditional", "zi", "zero_inflated", "all"),
    parameters = NULL,
    verbose = TRUE,
    ...
)

Arguments

x Vector representing a posterior distribution. Can also be a stanreg or brmsfit model.

... Currently not used.
range ROPE’s lower and higher bounds. Should be "default" or depending on the number of outcome variables a vector or a list. In models with one response, range should be a vector of length two (e.g., c(-0.1, 0.1)). In multivariate models, range should be a list with a numeric vectors for each response variable. Vector names should correspond to the name of the response variables. If "default" and input is a vector, the range is set to c(-0.1, 0.1). If "default" and input is a Bayesian model, rope_range() is used.

ci The Credible Interval (CI) probability, corresponding to the proportion of HDI, to use for the percentage in ROPE.

verbose Toggle off warnings.

effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior_) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

Details

Documentation is accessible for:

- Bayesian models
- Frequentist models

For Bayesian models, the Test for Practical Equivalence is based on the "HDI+ROPE decision rule" (Kruschke, 2014, 2018) to check whether parameter values should be accepted or rejected against an explicitly formulated "null hypothesis" (i.e., a ROPE). In other words, it checks the percentage of the 89\% HDI that is the null region (the ROPE). If this percentage is sufficiently low, the null hypothesis is rejected. If this percentage is sufficiently high, the null hypothesis is accepted.

Using the ROPE and the HDI, Kruschke (2018) suggests using the percentage of the 95\% (or 89\%, considered more stable) HDI that falls within the ROPE as a decision rule. If the HDI is completely outside the ROPE, the "null hypothesis" for this parameter is "rejected". If the ROPE completely covers the HDI, i.e., all most credible values of a parameter are inside the region of practical equivalence, the null hypothesis is accepted. Else, it’s undecided whether to accept or reject the null hypothesis. If the full ROPE is used (i.e., 100\% of the HDI), then the null hypothesis is rejected or accepted if the percentage of the posterior within the ROPE is smaller than to 2.5\% or greater than 97.5\%. Desirable results are low proportions inside the ROPE (the closer to zero the better).

Some attention is required for finding suitable values for the ROPE limits (argument range). See 'Details' in rope_range() for further information.

Multicollinearity: Non-independent covariates

When parameters show strong correlations, i.e. when covariates are not independent, the joint
parameter distributions may shift towards or away from the ROPE. In such cases, the test for practical equivalence may have inappropriate results. Collinearity invalidates ROPE and hypothesis testing based on univariate marginals, as the probabilities are conditional on independence. Most problematic are the results of the "undecided" parameters, which may either move further towards "rejection" or away from it (Kruschke 2014, 340f).

equivalence_test() performs a simple check for pairwise correlations between parameters, but as there can be collinearity between more than two variables, a first step to check the assumptions of this hypothesis testing is to look at different pair plots. An even more sophisticated check is the projection predictive variable selection (Piironen and Vehtari 2017).

**Value**

A data frame with following columns:

- Parameter The model parameter(s), if x is a model-object. If x is a vector, this column is missing.
- CI The probability of the HDI.
- ROPE_low, ROPE_high The limits of the ROPE. These values are identical for all parameters.
- ROPE_Percentage The proportion of the HDI that lies inside the ROPE.
- ROPE_Equivalence The "test result", as character. Either "rejected", "accepted" or "undecided".
- HDI_low, HDI_high The lower and upper HDI limits for the parameters.

**Note**

There is a print()-method with a digits-argument to control the amount of digits in the output, and there is a plot()-method to visualize the results from the equivalence-test (for models only).

**References**


**Examples**

library(bayestestR)
equivalence_test(x = rnorm(1000, 0, 0.01), range = c(-0.1, 0.1))
equivalence_test(x = rnorm(1000, 0, 1), range = c(-0.1, 0.1))
equivalence_test(x = rnorm(1000, 1, 0.01), range = c(-0.1, 0.1))
equivalence_test(x = rnorm(1000, 1, 1), ci = c(.50, .99))
```r
# print more digits
test <- equivalence_test(x = rnorm(1000, 1, 1), ci = c(.50, .99))
print(test, digits = 4)
## Not run:
library(rstanarm)
model <- rstanarm::stan_glm(mpg ~ wt + cyl, data = mtcars)
equivalence_test(model)
equivalence_test(model, ci = c(.50, 1))

# plot result
test <- equivalence_test(model)
plot(test)

library(emmeans)
equivalence_test(emtrends(model, ~1, "wt"))

library(brms)
model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
equivalence_test(model)
equivalence_test(model, ci = c(.50, .99))

library(BayesFactor)
bf <- ttestBF(x = rnorm(100, 1, 1))
equivalence_test(bf)
equivalence_test(bf, ci = c(.50, .99))
## End(Not run)
```

---

**estimate_density**

**Density Estimation**

**Description**

This function is a wrapper over different methods of density estimation. By default, it uses the base R `density` with by default uses a different smoothing bandwidth ("SJ") from the legacy default implemented the base R density function ("nrd0"). However, Deng & Wickham suggest that `method = "KernSmooth"` is the fastest and the most accurate.

**Usage**

```r
estimate_density(x, ...)
```

## S3 method for class 'data.frame'
estimate_density(
  x,
  method = "kernel",
  precision = 2^10,
  extend = FALSE,
  extend_scale = 0.1,
)
estimate_density

bw = "SJ",
ci = NULL,
select = NULL,
at = NULL,
group_by = NULL,
...
)

Arguments

x Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. bayestestR supports a wide range of models (see, for example, methods("hdi")) and not all of those are documented in the 'Usage' section, because methods for other classes mostly resemble the arguments of the .numeric or .data.frame methods.

... Currently not used.

method Density estimation method. Can be "kernel" (default), "logspline" or "KernSmooth".

precision Number of points of density data. See the n parameter in density.

extend Extend the range of the x axis by a factor of extend_scale.

extend_scale Ratio of range by which to extend the x axis. A value of 0.1 means that the x axis will be extended by 1/10 of the range of the data.

bw See the eponymous argument in density. Here, the default has been changed for "SJ", which is recommended.

ci The confidence interval threshold. Only used when method = "kernel". This feature is experimental, use with caution.

select Character vector of column names. If NULL (the default), all numeric variables will be selected. Other arguments from datawizard::find_columns() (such as exclude) can also be used.

at Optional character vector. If not NULL and input is a data frame, density estimation is performed for each group (subsets) indicated by at. See examples.

group_by Deprecated in favour of at.

Note

There is also a plot()-method implemented in the see-package.

References


Examples

library(bayestestR)

set.seed(1)
x <- rnorm(250, mean = 1)
# Basic usage
density_kernel <- estimate_density(x)  # default method is "kernel"

hist(x, prob = TRUE)
lines(density_kernel$x, density_kernel$y, col = "black", lwd = 2)
lines(density_kernel$x, density_kernel$CI_low, col = "gray", lty = 2)
lines(density_kernel$x, density_kernel$CI_high, col = "gray", lty = 2)
legend("topright",
    legend = c("Estimate", "95% CI"),
    col = c("black", "gray"), lwd = 2, lty = c(1, 2))

# Other Methods
density_logspline <- estimate_density(x, method = "logspline")
density_KernSmooth <- estimate_density(x, method = "KernSmooth")
density_mixture <- estimate_density(x, method = "mixture")

hist(x, prob = TRUE)
lines(density_kernel$x, density_kernel$y, col = "black", lwd = 2)
lines(density_logspline$x, density_logspline$y, col = "red", lwd = 2)
lines(density_KernSmooth$x, density_KernSmooth$y, col = "blue", lwd = 2)
lines(density_mixture$x, density_mixture$y, col = "green", lwd = 2)

# Extension
density_extended <- estimate_density(x, extend = TRUE)
density_default <- estimate_density(x, extend = FALSE)

hist(x, prob = TRUE)
lines(density_extended$x, density_extended$y, col = "red", lwd = 3)
lines(density_default$x, density_default$y, col = "black", lwd = 3)

# Multiple columns
head(estimate_density(iris))
head(estimate_density(iris, select = "Sepal.Width"))

# Grouped data
head(estimate_density(iris, at = "Species"))
head(estimate_density(iris$Petal.Width, at = iris$Species))
## Not run:
# rstanarm models
# -----------------------------------------------
library(rstanarm)
model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
head(estimate_density(model))

library(emmeans)
head(estimate_density(emtrends(model, ~1, "wt")))

# brms models
# -----------------------------------------------
library(brms)
model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
estimate_density(model)
Description

Compute the Equal-Tailed Interval (ETI) of posterior distributions using the quantiles method. The probability of being below this interval is equal to the probability of being above it. The ETI can be used in the context of uncertainty characterisation of posterior distributions as Credible Interval (CI).

Usage

eti(x, ...)

## S3 method for class 'numeric'
eti(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'stanreg'
eti(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                "distributional", "auxiliary"),
  parameters = NULL,
  verbose = TRUE,
  ...
)

## S3 method for class 'brmsfit'
eti(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL,
  verbose = TRUE,
  ...
)

Arguments

x  Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. bayestestR supports a wide range of models (see, for
example, methods("hdi") and not all of those are documented in the 'Usage' section, because methods for other classes mostly resemble the arguments of the .numeric or .data.frame methods.

... Currently not used.

ci Value or vector of probability of the (credible) interval - CI (between 0 and 1) to be estimated. Default to .95 (95%).

verbose Toggle off warnings.

effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior_) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

Details

Unlike equal-tailed intervals (see eti()) that typically exclude 2.5% from each tail of the distribution and always include the median, the HDI is not equal-tailed and therefore always includes the mode(s) of posterior distributions. While this can be useful to better represent the credibility mass of a distribution, the HDI also has some limitations. See spi() for details.

The 95% or 89% Credible Intervals (CI) are two reasonable ranges to characterize the uncertainty related to the estimation (see here for a discussion about the differences between these two values). The 89% intervals (ci = 0.89) are deemed to be more stable than, for instance, 95% intervals (Kruschke, 2014). An effective sample size of at least 10,000 is recommended if one wants to estimate 95% intervals with high precision (Kruschke, 2014, p. 183ff). Unfortunately, the default number of posterior samples for most Bayes packages (e.g., rstanarm or brms) is only 4,000 (thus, you might want to increase it when fitting your model). Moreover, 89 indicates the arbitrariness of interval limits - its only remarkable property is being the highest prime number that does not exceed the already unstable 95% threshold (McElreath, 2015). However, 95% has some advantages too. For instance, it shares (in the case of a normal posterior distribution) an intuitive relationship with the standard deviation and it conveys a more accurate image of the (artificial) bounds of the distribution. Also, because it is wider, it makes analyses more conservative (i.e., the probability of covering 0 is larger for the 95% CI than for lower ranges such as 89%), which is a good thing in the context of the reproducibility crisis.

A 95% equal-tailed interval (ETI) has 2.5% of the distribution on either side of its limits. It indicates the 2.5th percentile and the 97.5th percentile. In symmetric distributions, the two methods of computing credible intervals, the ETI and the HDI, return similar results.

This is not the case for skewed distributions. Indeed, it is possible that parameter values in the ETI have lower credibility (are less probable) than parameter values outside the ETI. This property seems undesirable as a summary of the credible values in a distribution.
On the other hand, the ETI range does change when transformations are applied to the distribution (for instance, for a log odds scale to probabilities): the lower and higher bounds of the transformed distribution will correspond to the transformed lower and higher bounds of the original distribution. On the contrary, applying transformations to the distribution will change the resulting HDI.

Value

A data frame with following columns:

- **Parameter** The model parameter(s), if x is a model-object. If x is a vector, this column is missing.
- **CI** The probability of the credible interval.
- **CI_low, CI_high** The lower and upper credible interval limits for the parameters.

See Also

Other ci: `bci()`, `ci()`, `cwi()`, `hdi()`, `si()`, `spi()`

Examples

```r
library(bayestestR)

posterior <- rnorm(1000)
eti(posterior)
eti(posterior, ci = c(.80, .89, .95))

df <- data.frame(replicate(4, rnorm(100)))
eti(df)
eti(df, ci = c(.80, .89, .95))
## Not run:
library(rstanarm)
model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
eti(model)
eti(model, ci = c(.80, .89, .95))
library(emmeans)
eti(emtrends(model, ~1, "wt"))

library(brms)
model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
eti(model)
eti(model, ci = c(.80, .89, .95))

library(BayesFactor)
bf <- ttestBF(x = rnorm(100, 1, 1))
eti(bf)
eti(bf, ci = c(.80, .89, .95))
## End(Not run)
```
**hdi**  
*Highest Density Interval (HDI)*

**Description**

Compute the **Highest Density Interval (HDI)** of posterior distributions. All points within this interval have a higher probability density than points outside the interval. The HDI can be used in the context of uncertainty characterisation of posterior distributions as **Credible Interval (CI)**.

**Usage**

```r
hdi(x, ...)  

## S3 method for class 'numeric'
hdi(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'data.frame'
hdi(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'stanreg'
hdi(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma", "distributional", "auxiliary"),
  parameters = NULL,
  verbose = TRUE,
  ...
)

## S3 method for class 'brmsfit'
hdi(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL,
  verbose = TRUE,
  ...
)
```

**Arguments**

- **x**: Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. **bayestestR** supports a wide range of models (see, for example, `methods("hdi")`) and not all of those are documented in the 'Usage'
section, because methods for other classes mostly resemble the arguments of the `numeric` or `data.frame` methods.

... Currently not used.

ci Value or vector of probability of the (credible) interval - CI (between 0 and 1) to be estimated. Default to .95 (95%).

verbose Toggle off warnings.

effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to `brms`-models.

parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like `lp__` or `prior_`) are filtered by default, so only parameters that typically appear in the `summary()` are returned. Use `parameters` to select specific parameters for the output.

Details

Unlike equal-tailed intervals (see `eti()`) that typically exclude 2.5% from each tail of the distribution and always include the median, the HDI is not equal-tailed and therefore always includes the mode(s) of posterior distributions. While this can be useful to better represent the credibility mass of a distribution, the HDI also has some limitations. See `spi()` for details.

The 95% or 89% Credible Intervals (CI) are two reasonable ranges to characterize the uncertainty related to the estimation (see here for a discussion about the differences between these two values). The 89% intervals (ci = 0.89) are deemed to be more stable than, for instance, 95% intervals (Kruschke, 2014). An effective sample size of at least 10,000 is recommended if one wants to estimate 95% intervals with high precision (Kruschke, 2014, p. 183ff). Unfortunately, the default number of posterior samples for most Bayes packages (e.g., `rstanarm` or `brms`) is only 4,000 (thus, you might want to increase it when fitting your model). Moreover, 89 indicates the arbitrariness of interval limits - its only remarkable property is being the highest prime number that does not exceed the already unstable 95% threshold (McElreath, 2015).

However, 95% has some advantages too. For instance, it shares (in the case of a normal posterior distribution) an intuitive relationship with the standard deviation and it conveys a more accurate image of the (artificial) bounds of the distribution. Also, because it is wider, it makes analyses more conservative (i.e., the probability of covering 0 is larger for the 95% CI than for lower ranges such as 89%), which is a good thing in the context of the reproducibility crisis.

A 95% equal-tailed interval (ETI) has 2.5% of the distribution on either side of its limits. It indicates the 2.5th percentile and the 97.5th percentile. In symmetric distributions, the two methods of computing credible intervals, the ETI and the HDI, return similar results.

This is not the case for skewed distributions. Indeed, it is possible that parameter values in the ETI have lower credibility (are less probable) than parameter values outside the ETI. This property seems undesirable as a summary of the credible values in a distribution.

On the other hand, the ETI range does change when transformations are applied to the distribution
(for instance, for a log odds scale to probabilities): the lower and higher bounds of the transformed distribution will correspond to the transformed lower and higher bounds of the original distribution. On the contrary, applying transformations to the distribution will change the resulting HDI.

Value

A data frame with following columns:

- **Parameter**: The model parameter(s), if x is a model-object. If x is a vector, this column is missing.
- **CI**: The probability of the credible interval.
- **CI_low, CI_high**: The lower and upper credible interval limits for the parameters.

Note

There is also a `plot()`-method implemented in the `see-package`.

Author(s)

Credits go to `ggdistribute` and `HDInterval`.

References


See Also

Other interval functions, such as `hdi()`, `eti()`, `bci()`, `spi()`, `si()`, `cwi()`.

Other ci: `bci()`, `ci()`, `cwi()`, `eti()`, `si()`, `spi()`

Examples

```r
library(bayestestR)

posterior <- rnorm(1000)
hdi(posterior, ci = .89)
hdi(posterior, ci = c(.80, .90, .95))

df <- data.frame(replicate(4, rnorm(100)))
hdi(df)
hdi(df, ci = c(.80, .90, .95))
## Not run:
library(rstanarm)
model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
hdi(model)
hdi(model, ci = c(.80, .90, .95))
```
library(emmeans)
hdni(emtrends(model, ~1, "wt"))

library(brms)
model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
hdni(model)
hdni(model, ci = c(.80, .90, .95))

library(BayesFactor)
bf <- ttestBF(x = rnorm(100, 1, 1))
hdni(bf)
hdni(bf, ci = c(.80, .90, .95))

## End(Not run)

map_estimate

Maximum A Posteriori probability estimate (MAP)

Description

Find the **Highest Maximum A Posteriori probability estimate (MAP)** of a posterior, i.e., the value associated with the highest probability density (the "peak" of the posterior distribution). In other words, it is an estimation of the **mode** for continuous parameters. Note that this function relies on estimate_density, which by default uses a different smoothing bandwidth ("SJ") compared to the legacy default implemented the base R density function ("nrd0").

Usage

map_estimate(x, precision = 2^10, method = "kernel", ...)

## S3 method for class 'numeric'
map_estimate(x, precision = 2^10, method = "kernel", ...)

## S3 method for class 'stanreg'
map_estimate(
x, 
precision = 2^10, 
method = "kernel", 
effects = c("fixed", "random", "all"), 
component = c("location", "all", "conditional", "smooth_terms", "sigma", 
  "distributional", "auxiliary"), 
parameters = NULL, 
... 
)

## S3 method for class 'brmsfit'
map_estimate(
x, 
... 
)
map_estimate

```r
precision = 2^10,
method = "kernel",
effects = c("fixed", "random", "all"),
component = c("conditional", "zi", "zero_inflated", "all"),
parameters = NULL,
...
)
```

## S3 method for class 'data.frame'
map_estimate(x, precision = 2^10, method = "kernel", ...)

### Arguments

- **x**: Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. **bayestestR** supports a wide range of models (see, for example, `methods("hdi")`) and not all of those are documented in the 'Usage' section, because methods for other classes mostly resemble the arguments of the `.numeric` or `.data.frame` methods.
- **precision**: Number of points of density data. See the `n` parameter in `density`.
- **method**: Density estimation method. Can be "kernel" (default), "logspline" or "KernSmooth".
- **...**: Currently not used.
- **effects**: Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
- **component**: Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to **brms**-models.
- **parameters**: Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like `lp__` or `prior_`) are filtered by default, so only parameters that typically appear in the `summary()` are returned. Use parameters to select specific parameters for the output.

### Value

A numeric value if `x` is a vector. If `x` is a model-object, returns a data frame with following columns:

- **Parameter** The model parameter(s), if `x` is a model-object. If `x` is a vector, this column is missing.
- **MAP_Estimate** The MAP estimate for the posterior or each model parameter.

### Examples

```r
## Not run:
library(bayestestR)
posterior <- rnorm(10000)
map_estimate(posterior)
plot(density(posterior))
```
mcse

Monte-Carlo Standard Error (MCSE)

Description

This function returns the Monte Carlo Standard Error (MCSE).

Usage

mcse(model, ...)

## S3 method for class 'stanreg'
mcse(model, 
    effects = c("fixed", "random", "all"),
    component = c("location", "all", "conditional", "smooth_terms", "sigma", "distributional", "auxiliary"),
    parameters = NULL,
    ...
)

Arguments

model A stanreg, stanfit, brmsfit, blavaan, or MCMCglmm object.

... Currently not used.
effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior__) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.
Details

**Monte Carlo Standard Error (MCSE)** is another measure of accuracy of the chains. It is defined as standard deviation of the chains divided by their effective sample size (the formula for `mcse()` is from Kruschke 2015, p. 187). The MCSE “provides a quantitative suggestion of how big the estimation noise is”.

References


Examples

```r
## Not run:
library(bayestestR)
library(rstanarm)

model <- stan_glm(mpg ~ wt + am, data = mtcars, chains = 1, refresh = 0)
mcse(model)
## End(Not run)
```

mediation

Summary of Bayesian multivariate-response mediation-models

Description

`mediation()` is a short summary for multivariate-response mediation-models, i.e. this function computes average direct and average causal mediation effects of multivariate response models.

Usage

```r
mediation(model, ...)
```

## S3 method for class 'brmsfit'
```r
mediation(
  model,
  treatment,
  mediator,
  response = NULL,
  centrality = "median",
  ci = 0.95,
  method = "ETI",
  ...
)
```

## S3 method for class 'stanmvreg'

mediation(
  model,
  treatment,
  mediator,
  response = NULL,
  centrality = "median",
  ci = 0.95,
  method = "ETI",
  ...
)

Arguments

model A brmsfit or stanmvreg object.
... Not used.
treatment Character, name of the treatment variable (or direct effect) in a (multivariate response) mediator-model. If missing, mediation() tries to find the treatment variable automatically, however, this may fail.
mediator Character, name of the mediator variable in a (multivariate response) mediator-model. If missing, mediation() tries to find the treatment variable automatically, however, this may fail.
response A named character vector, indicating the names of the response variables to be used for the mediation analysis. Usually can be NULL, in which case these variables are retrieved automatically. If not NULL, names should match the names of the model formulas, names(insight::find_response(model, combine = TRUE)). This can be useful if, for instance, the mediator variable used as predictor has a different name from the mediator variable used as response. This might occur when the mediator is transformed in one model, but used "as is" as response variable in the other model. Example: The mediator m is used as response variable, but the centered version m_center is used as mediator variable. The second response variable (for the treatment model, with the mediator as additional predictor), y, is not transformed. Then we could use response like this: mediation(model, response = c(m = "m_center", y = "y"))
centrality The point-estimates (centrality indices) to compute. Character (vector) or list with one or more of these options: "median", "mean", "MAP" or "all".
ci Value or vector of probability of the CI (between 0 and 1) to be estimated. Default to .95 (95%).
method Can be 'ETI' (default), 'HDI', 'BCI', 'SPI' or 'SI'.

Details

mediation() returns a data frame with information on the direct effect (mean value of posterior samples from treatment of the outcome model), mediator effect (mean value of posterior samples from mediator of the outcome model), indirect effect (mean value of the multiplication of the posterior samples from mediator of the outcome model and the posterior samples from treatment of the mediation model) and the total effect (mean value of sums of posterior samples used for the direct and indirect effect). The proportion mediated is the indirect effect divided by the total effect.
For all values, the 89% credible intervals are calculated by default. Use ci to calculate a different interval.

The arguments treatment and mediator do not necessarily need to be specified. If missing, mediation() tries to find the treatment and mediator variable automatically. If this does not work, specify these variables.

The direct effect is also called *average direct effect* (ADE), the indirect effect is also called *average causal mediation effects* (ACME). See also Tingley et al. 2014 and Imai et al. 2010.

**Value**

A data frame with direct, indirect, mediator and total effect of a multivariate-response mediation-model, as well as the proportion mediated. The effect sizes are median values of the posterior samples (use centrality for other centrality indices).

**Note**

There is an as.data.frame() method that returns the posterior samples of the effects, which can be used for further processing in the different bayestestR package.

**References**


**See Also**

The mediation package for a causal mediation analysis in the frequentist framework.

**Examples**

```r
## Not run:
library(mediation)
library(brms)
library(rstanarm)
# load sample data
data(jobs)
set.seed(123)

# linear models, for mediation analysis
b1 <- lm(job_seek ~ treat + econ_hard + sex + age, data = jobs)
b2 <- lm(depress2 ~ treat + job_seek + econ_hard + sex + age, data = jobs)
# mediation analysis, for comparison with Stan models
m1 <- mediate(b1, b2, sims = 1000, treat = "treat", mediator = "job_seek")

# Fit Bayesian mediation model in brms
```
f1 <- bf(job_seek ~ treat + econ_hard + sex + age)
f2 <- bf(depress2 ~ treat + job_seek + econ_hard + sex + age)
m2 <- brm(f1 + f2 + set_rescor(FALSE), data = jobs, cores = 4, refresh = 0)

# Fit Bayesian mediation model in rstanarm
m3 <- stan_mvmer(
  list(
    job_seek ~ treat + econ_hard + sex + age + (1 | occp),
    depress2 ~ treat + job_seek + econ_hard + sex + age + (1 | occp)
  ),
  data = jobs,
  cores = 4,
  refresh = 0
)

summary(m1)
mediation(m2, centrality = "mean", ci = .95)
mediation(m3, centrality = "mean", ci = .95)

## End(Not run)

---

### model_to_priors

Convert model’s posteriors to priors *(EXPERIMENTAL)*

#### Description

Convert model’s posteriors to (normal) priors.

#### Usage

```r
model_to_priors(model, scale_multiply = 3, ...)
```

#### Arguments

- `model` A Bayesian model.
- `scale_multiply` The SD of the posterior will be multiplied by this amount before being set as a prior to avoid overly narrow priors.
- `...` Other arguments for `insight::get_prior()` or `describe_posterior`.

#### Examples

```r
## Not run:
# brms models
# -----------------------------------------------
if (require("brms")) {
  formula <- brms::brmsformula(mpg ~ wt + cyl, center = FALSE)
  model <- brms::brm(formula, data = mtcars, refresh = 0)
  priors <- model_to_priors(model)
}
```r
priors <- brms::validate_prior(priors, formula, data = mtcars)
priors

model2 <- brms::brm(formula, data = mtcars, prior = priors, refresh = 0)
}

## End(Not run)
```

---

**overlap**

*Overlap Coefficient*

**Description**

A method to calculate the overlap coefficient between two empirical distributions (that can be used as a measure of similarity between two samples).

**Usage**

```r
overlap(
  x, y,
  method_density = "kernel",
  method_auc = "trapezoid",
  precision = 2^10,
  extend = TRUE,
  extend_scale = 0.1,
  ...
)
```

**Arguments**

- `x` Vector of x values.
- `y` Vector of x values.
- `method_density` Density estimation method. See `estimate_density()`.
- `method_auc` Area Under the Curve (AUC) estimation method. See `area_under_curve()`.
- `precision` Number of points of density data. See the n parameter in `density`.
- `extend` Extend the range of the x axis by a factor of `extend_scale`.
- `extend_scale` Ratio of range by which to extend the x axis. A value of 0.1 means that the x axis will be extended by 1/10 of the range of the data.
- `...` Currently not used.
Examples

library(bayestestR)

x <- distribution_normal(1000, 2, 0.5)
y <- distribution_normal(1000, 0, 1)

overlap(x, y)
plot(overlap(x, y))

---

**pd_to_p**

*Convert between Probability of Direction (pd) and p-value.*

**Description**

Enables a conversion between Probability of Direction (pd) and p-value.

**Usage**

- pd_to_p(pd, direction = "two-sided", ...)
- p_to_pd(p, direction = "two-sided", ...)
- convert_p_to_pd(p, direction = "two-sided", ...)
- convert_pd_to_p(pd, direction = "two-sided", ...)

**Arguments**

- **pd**  
  A Probability of Direction (pd) value (between 0 and 1).
- **direction**  
  What type of p-value is requested or provided. Can be "two-sided" (default, two tailed) or "one-sided" (one tailed).
- **...**  
  Arguments passed to or from other methods.
- **p**  
  A p-value.

**Examples**

- pd_to_p(pd = 0.95)
- pd_to_p(pd = 0.95, direction = "one-sided")
Description

Compute various point-estimates, such as the mean, the median or the MAP, to describe posterior distributions.

Usage

point_estimate(x, ...)

## S3 method for class 'numeric'
point_estimate(x, centrality = "all", dispersion = FALSE, threshold = 0.1, ...)

## S3 method for class 'stanreg'
point_estimate(
x,
  centrality = "all",
  dispersion = FALSE,
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                "distributional", "auxiliary"),
  parameters = NULL,
  ...)

## S3 method for class 'brmsfit'
point_estimate(
x,
  centrality = "all",
  dispersion = FALSE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL,
  ...)

## S3 method for class 'BFBayesFactor'
point_estimate(x, centrality = "all", dispersion = FALSE, ...)

Arguments

x Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. bayestestR supports a wide range of models (see, for example, methods("hdi")) and not all of those are documented in the 'Usage'
section, because methods for other classes mostly resemble the arguments of the 
\texttt{.numeric} or \texttt{.data.frame} methods.

... Additional arguments to be passed to or from methods.

\textbf{centrality} The point-estimates (centrality indices) to compute. Character (vector) or list with one or more of these options: "median", "mean", "MAP" or "all".

\textbf{dispersion} Logical, if \texttt{TRUE}, computes indices of dispersion related to the estimate(s) (SD and MAD for mean and median, respectively).

\textbf{threshold} For \texttt{centrality = "trimmed"} (i.e. trimmed mean), indicates the fraction (0 to 0.5) of observations to be trimmed from each end of the vector before the mean is computed.

\textbf{effects} Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

\textbf{component} Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to \texttt{brms}-models.

\textbf{parameters} Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like \texttt{lp__} or \texttt{prior__}) are filtered by default, so only parameters that typically appear in the \texttt{summary()} are returned. Use \texttt{parameters} to select specific parameters for the output.

\textbf{Note}

There is also a \texttt{plot()}-method implemented in the \texttt{see}-package.

\textbf{References}


\textbf{Examples}

library(bayestestR)

point_estimate(rnorm(1000))
point_estimate(rnorm(1000), centrality = "all", dispersion = \texttt{TRUE})
point_estimate(rnorm(1000), centrality = c("median", "MAP"))

df <- data.frame(replicate(4, rnorm(100)))
point_estimate(df, centrality = "all", dispersion = \texttt{TRUE})
point_estimate(df, centrality = c("median", "MAP"))

## Not run:
# rstanarm models
# -----------------------------------------------
library(rstanarm)
model <- rstanarm::stan_glm(mpg ~ wt + cyl, data = mtcars)
point_estimate(model, centrality = "all", dispersion = \texttt{TRUE})
point_estimate(model, centrality = c("median", "MAP"))
p_direction

Probability of Direction (pd)

Description

Compute the **Probability of Direction** (pd, also known as the Maximum Probability of Effect - MPE). It varies between 50% and 100% (i.e., 0.5 and 1) and can be interpreted as the probability (expressed in percentage) that a parameter (described by its posterior distribution) is strictly positive or negative (whichever is the most probable). It is mathematically defined as the proportion of the posterior distribution that is of the median’s sign. Although differently expressed, this index is fairly similar (i.e., is strongly correlated) to the frequentist p-value.

Note that in some (rare) cases, especially when used with model averaged posteriors (see `weighted_posteriors()`) or `brms::posterior_average`, `pd` can be smaller than 0.5, reflecting high credibility of 0.

Usage

```r
p_direction(x, ...)
pd(x, ...)
```

## S3 method for class 'numeric'
`p_direction(x, method = "direct", null = 0, ...)`

## S3 method for class 'data.frame'
p_direction(x, method = "direct", null = 0, ...)

## S3 method for class 'MCMCglmm'
p_direction(x, method = "direct", null = 0, ...)

## S3 method for class 'emmGrid'
p_direction(x, method = "direct", null = 0, ...)

## S3 method for class 'stanreg'
p_direction(x, effects = c("fixed", "random", "all"),
            component = c("location", "all", "conditional", "smooth_terms", "sigma",
                          "distributional", "auxiliary"),
            parameters = NULL,
            method = "direct",
            null = 0,
            ...
        )

## S3 method for class 'brmsfit'
p_direction(x, effects = c("fixed", "random", "all"),
            component = c("conditional", "zi", "zero_inflated", "all"),
            parameters = NULL,
            method = "direct",
            null = 0,
            ...
        )

## S3 method for class 'BFBayesFactor'
p_direction(x, method = "direct", null = 0, ...)

Arguments

x Vector representing a posterior distribution. Can also be a Bayesian model (stanreg, brmsfit or BayesFactor).

... Currently not used.

method Can be "direct" or one of methods of density estimation, such as "kernel", "logspline" or "KernSmooth". If "direct" (default), the computation is based on the raw ratio of samples superior and inferior to 0. Else, the result is based on the Area under the Curve (AUC) of the estimated density function.

null The value considered as a "null" effect. Traditionally 0, but could also be 1 in the case of ratios.

effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
component

Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to `brms`-models.

parameters

Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like `lp__` or `prior_`) are filtered by default, so only parameters that typically appear in the `summary()` are returned. Use `parameters` to select specific parameters for the output.

Details

What is the pd?: The Probability of Direction (pd) is an index of effect existence, ranging from 50% to 100%, representing the certainty with which an effect goes in a particular direction (i.e., is positive or negative). Beyond its simplicity of interpretation, understanding and computation, this index also presents other interesting properties:

- It is independent from the model: It is solely based on the posterior distributions and does not require any additional information from the data or the model.
- It is robust to the scale of both the response variable and the predictors.
- It is strongly correlated with the frequentist p-value, and can thus be used to draw parallels and give some reference to readers non-familiar with Bayesian statistics.

Relationship with the p-value: In most cases, it seems that the `pd` has a direct correspondence with the frequentist one-sided p-value through the formula $p_{onesided} = 1 - \frac{pd}{100}$ and to the two-sided p-value (the most commonly reported one) through the formula $p_{twosided} = 2 \times (1 - \frac{pd}{100})$. Thus, a two-sided p-value of respectively .1, .05, .01 and .001 would correspond approximately to a `pd` of 95%, 97.5%, 99.5% and 99.95%. See also `pd_to_p()`.

Methods of computation: The most simple and direct way to compute the `pd` is to 1) look at the median’s sign, 2) select the portion of the posterior of the same sign and 3) compute the percentage that this portion represents. This "simple" method is the most straightforward, but its precision is directly tied to the number of posterior draws. The second approach relies on density estimation. It starts by estimating the density function (for which many methods are available), and then computing the area under the curve (AUC) of the density curve on the other side of 0.

Strengths and Limitations: Strengths: Straightforward computation and interpretation. Objective property of the posterior distribution. 1:1 correspondence with the frequentist p-value.

Limitations: Limited information favoring the null hypothesis.

Value

Values between 0.5 and 1 corresponding to the probability of direction (pd).

Note that in some (rare) cases, especially when used with model averaged posteriors (see `weighted_posteriors()` or `brms::posterior_average`), pd can be smaller than 0.5, reflecting high credibility of 0. To detect such cases, the method = "direct" must be used.

Note

There is also a `plot()`-method implemented in the `see-package`. 

p_direction
References


See Also

pd_to_p() to convert between Probability of Direction (pd) and p-value.

Examples

library(bayestestR)

# Simulate a posterior distribution of mean 1 and SD 1
# ----------------------------------------------------
posterior <- rnorm(1000, mean = 1, sd = 1)
p_direction(posterior)
p_direction(posterior, method = "kernel")

# Simulate a dataframe of posterior distributions
# -----------------------------------------------
df <- data.frame(replicate(4, rnorm(100)))
p_direction(df)
p_direction(df, method = "kernel")
## Not run:
# rstanarm models
# -----------------------------------------------
if (require("rstanarm")) {
  model <- rstanarm::stan_glm(mpg ~ wt + cyl,
    data = mtcars,
    chains = 2, refresh = 0
  )
p_direction(model)
p_direction(model, method = "kernel")
}

# emmeans
# -----------------------------------------------
if (require("emmeans")) {
  p_direction(emtrends(model, ~1, "wt"))
}

# brms models
# -----------------------------------------------
if (require("brms")) {
  model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
p_direction(model)
p_direction(model, method = "kernel")
}

# BayesFactor objects
# -----------------------------------------------
### p_map

Bayesian p-value based on the density at the Maximum A Posteriori (MAP)

#### Description

Compute a Bayesian equivalent of the p-value, related to the odds that a parameter (described by its posterior distribution) has against the null hypothesis (h0) using Mills' (2014, 2017) *Objective Bayesian Hypothesis Testing* framework. It corresponds to the density value at 0 divided by the density at the Maximum A Posteriori (MAP).

#### Usage

```r
p_map(x, precision = 2^10, method = "kernel", ...)
p_pointnull(x, precision = 2^10, method = "kernel", ...)
```

#### S3 method for class 'stanreg'

```r
p_map(
    x,
    precision = 2^10,
    method = "kernel",
    effects = c("fixed", "random", "all"),
    component = c("location", "all", "conditional", "smooth_terms", "sigma",
                   "distributional", "auxiliary"),
    parameters = NULL,
    ...
)
```

#### S3 method for class 'brmsfit'

```r
p_map(
    x,
    precision = 2^10,
    method = "kernel",
    effects = c("fixed", "random", "all"),
    component = c("conditional", "zi", "zero_inflated", "all"),
    parameters = NULL,
    ...
)
```
Arguments

- **x**: Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. `bayestestR` supports a wide range of models (see, for example, `methods("hdi")`) and not all of those are documented in the 'Usage' section, because methods for other classes mostly resemble the arguments of the `numeric` or `data.frame` methods.

- **precision**: Number of points of density data. See the `n` parameter in `density`.

- **method**: Density estimation method. Can be "kernel" (default), "logspline" or "KernSmooth".

- **...**: Currently not used.

- **effects**: Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

- **component**: Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to `brms`-models.

- **parameters**: Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like `lp__` or `prior_`) are filtered by default, so only parameters that typically appear in the `summary()` are returned. Use `parameters` to select specific parameters for the output.

Details

Note that this method is sensitive to the density estimation method (see the section in the examples below).

**Strengths and Limitations**: **Strengths**: Straightforward computation. Objective property of the posterior distribution.

**Limitations**: Limited information favoring the null hypothesis. Relates on density approximation. Indirect relationship between mathematical definition and interpretation. Only suitable for weak / very diffused priors.

References


See Also

Jeff Mill’s talk

Examples

```r
library(bayestestR)
p_map(rnorm(1000, 0, 1))
```
p_rope

p_map(rnorm(1000, 10, 1))
## Not run:
library(rstanarm)
model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
p_map(model)

library(emmeans)
p_map(emtrends(model, ~1, "wt"))

library(brms)
model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
p_map(model)

library(BayesFactor)
bf <- ttestBF(x = rnorm(100, 1, 1))
p_map(bf)

# ---------------------------------------
# Robustness to density estimation method
set.seed(333)
data <- data.frame()
for (iteration in 1:250) {
  x <- rnorm(1000, 1, 1)
  result <- data.frame(
    "Kernel" = p_map(x, method = "kernel"),
    "KernSmooth" = p_map(x, method = "KernSmooth"),
    "logspline" = p_map(x, method = "logspline")
  )
  data <- rbind(data, result)
}
data$KernSmooth <- data$Kernel - data$KernSmooth
data$logspline <- data$Kernel - data$logspline

summary(data$KernSmooth)
summary(data$logspline)
boxplot(data[,c("KernSmooth", "logspline")])
## End(Not run)

p_rope

Probability of being in the ROPE

Description
Compute the proportion of the whole posterior distribution that doesn’t lie within a region of practical equivalence (ROPE). It is equivalent to running rope(..., ci = 1).

Usage
p_rope(x, ...)
## S3 method for class 'numeric'
p_rope(x, range = "default", ...)

## S3 method for class 'stanreg'
p_rope(
  x,
  range = "default",
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                "distributional", "auxiliary"),
  parameters = NULL,
  ...
)

## S3 method for class 'brmsfit'
p_rope(
  x,
  range = "default",
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL,
  ...
)

### Arguments

- **x**
  Vector representing a posterior distribution. Can also be a stanreg or brmsfit model.

- **...**
  Currently not used.

- **range**
  ROPE's lower and higher bounds. Should be "default" or depending on the number of outcome variables a vector or a list. In models with one response, range should be a vector of length two (e.g., `c(-0.1, 0.1)`). In multivariate models, range should be a list with a numeric vectors for each response variable. Vector names should correspond to the name of the response variables. If "default" and input is a vector, the range is set to `c(-0.1, 0.1)`. If "default" and input is a Bayesian model, `rope_range()` is used.

- **effects**
  Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

- **component**
  Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

- **parameters**
  Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior_) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.
Examples

```r
library(bayestestR)

p_rope(x = rnorm(1000, 0, 0.01), range = c(-0.1, 0.1))
p_rope(x = mtcars, range = c(-0.1, 0.1))
```

---

**p_significance**  
Practical Significance (ps)

Description

Compute the probability of Practical Significance (ps), which can be conceptualized as a unidirectional equivalence test. It returns the probability that effect is above a given threshold corresponding to a negligible effect in the median’s direction. Mathematically, it is defined as the proportion of the posterior distribution of the median sign above the threshold.

Usage

```r
p_significance(x, ...)
```

## S3 method for class 'numeric'

```r
p_significance(x, threshold = "default", ...)
```

## S3 method for class 'stanreg'

```r
p_significance(
  x,
  threshold = "default",
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                  "distributional", "auxiliary"),
  parameters = NULL,
  verbose = TRUE,
  ...
)
```

## S3 method for class 'brmsfit'

```r
p_significance(
  x,
  threshold = "default",
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL,
  verbose = TRUE,
  ...
)
```
Arguments

x  Vector representing a posterior distribution. Can also be a stanreg or brmsfit model.

...  Currently not used.

threshold  The threshold value that separates significant from negligible effect. If "default", the range is set to 0.1 if input is a vector, and based on rope_range() if a Bayesian model is provided.

effects  Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component  Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.

parameters  Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior_) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

verbose  Toggle off warnings.

Details

p_significance() returns the proportion of a probability distribution (x) that is outside a certain range (the negligible effect, or ROPE, see argument threshold). If there are values of the distribution both below and above the ROPE, p_significance() returns the higher probability of a value being outside the ROPE. Typically, this value should be larger than 0.5 to indicate practical significance. However, if the range of the negligible effect is rather large compared to the range of the probability distribution x, p_significance() will be less than 0.5, which indicates no clear practical significance.

Value

Values between 0 and 1 corresponding to the probability of practical significance (ps).

Note

There is also a plot()-method implemented in the see-package.

Examples

library(bayestestR)

# Simulate a posterior distribution of mean 1 and SD 1
# -----------------------------------------------
posterior <- rnorm(1000, mean = 1, sd = 1)
p_significance(posterior)

# Simulate a dataframe of posterior distributions
# -----------------------------------------------
df <- data.frame(replicate(4, rnorm(100)))
## Not run:
# rstanarm models
# -----------------------------------------------
if (require("rstanarm")) {
  model <- rstanarm::stan_glm(mpg ~ wt + cyl,
    data = mtcars,
    chains = 2, refresh = 0
  )
  p_significance(model)
}
## End(Not run)

### reshape_iterations

Reshape estimations with multiple iterations (draws) to long format

**Description**

Reshape a wide data.frame of iterations (such as posterior draws or bootstrapped samples) as columns to long format. Instead of having all iterations as columns (e.g., iter_1, iter_2, ...), will return 3 columns with the \*_index (the previous index of the row), the \*_group (the iteration number) and the \*_value (the value of said iteration).

**Usage**

```r
reshape_iterations(x, prefix = c("draw", "iter", "iteration", "sim"))
```

```r
reshape_draws(x, prefix = c("draw", "iter", "iteration", "sim"))
```

**Arguments**

- **x**: A data.frame containing posterior draws obtained from `estimate_response` or `estimate_link`.
- **prefix**: The prefix of the draws (for instance, "iter_" for columns named as iter_1, iter_2, iter_3). If more than one are provided, will search for the first one that matches.

**Value**

Data frame of reshaped draws in long format.

**Examples**

```r
if (require("rstanarm")) {
  model <- stan_glm(mpg ~ am, data = mtcars, refresh = 0)
  draws <- insight::get_predicted(model)
  long_format <- reshape_iterations(draws)
  head(long_format)
```
Region of Practical Equivalence (ROPE)

Description

Compute the proportion of the HDI (default to the 89% HDI) of a posterior distribution that lies within a region of practical equivalence.

Usage

rope(x, ...)

## S3 method for class 'numeric'
rope(x, range = "default", ci = 0.95, ci_method = "ETI", verbose = TRUE, ...)

## S3 method for class 'stanreg'
rope(
  x,
  range = "default",
  ci = 0.95,
  ci_method = "ETI",
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                "distributional", "auxiliary"),
  parameters = NULL,
  verbose = TRUE,
  ...
)

## S3 method for class 'brmsfit'
rope(
  x,
  range = "default",
  ci = 0.95,
  ci_method = "ETI",
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL,
  verbose = TRUE,
  ...
)
Arguments

- **Arguments**

- **x**: Vector representing a posterior distribution. Can also be a `stanreg` or `brmsfit` model.

- **...**: Currently not used.

- **range**: ROPE’s lower and higher bounds. Should be "default" or depending on the number of outcome variables a vector or a list. In models with one response, range should be a vector of length two (e.g., `c(-0.1, 0.1)`). In multivariate models, range should be a list with a numeric vector for each response variable. Vector names should correspond to the name of the response variables. If "default" and input is a vector, the range is set to `c(-0.1, 0.1)`. If "default" and input is a Bayesian model, `rope_range()` is used.

- **ci**: The Credible Interval (CI) probability, corresponding to the proportion of HDI, to use for the percentage in ROPE.

- **ci_method**: The type of interval to use to quantify the percentage in ROPE. Can be 'HDI' (default) or 'ETI'. See `ci()`.

- **verbose**: Toggle off warnings.

- **effects**: Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

- **component**: Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to `brms`-models.

- **parameters**: Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like `lp__` or `prior_`) are filtered by default, so only parameters that typically appear in the `summary()` are returned. Use parameters to select specific parameters for the output.

Details

**ROPE**: Statistically, the probability of a posterior distribution of being different from 0 does not make much sense (the probability of a single value null hypothesis in a continuous distribution is 0). Therefore, the idea underlining ROPE is to let the user define an area around the null value enclosing values that are equivalent to the null value for practical purposes (Kruschke 2010, 2011, 2014).

Kruschke (2018) suggests that such null value could be set, by default, to the -0.1 to 0.1 range of a standardized parameter (negligible effect size according to Cohen, 1988). This could be generalized: For instance, for linear models, the ROPE could be set as `0 +/- .1 * sd(y)`. This ROPE range can be automatically computed for models using the `rope_range` function.

Kruschke (2010, 2011, 2014) suggests using the proportion of the 95% (or 89%, considered more stable) HDI that falls within the ROPE as an index for "null-hypothesis" testing (as understood under the Bayesian framework, see `equivalence_test()`).

**Sensitivity to parameter’s scale**: It is important to consider the unit (i.e., the scale) of the predictors when using an index based on the ROPE, as the correct interpretation of the ROPE as representing a region of practical equivalence to zero is dependent on the scale of the predictors.
Indeed, the percentage in ROPE depend on the unit of its parameter. In other words, as the ROPE represents a fixed portion of the response’s scale, its proximity with a coefficient depends on the scale of the coefficient itself.

**Multicollinearity: Non-independent covariates:** When parameters show strong correlations, i.e. when covariates are not independent, the joint parameter distributions may shift towards or away from the ROPE. Collinearity invalidates ROPE and hypothesis testing based on univariate marginals, as the probabilities are conditional on independence. Most problematic are parameters that only have partial overlap with the ROPE region. In case of collinearity, the (joint) distributions of these parameters may either get an increased or decreased ROPE, which means that inferences based on `rope()` are inappropriate (Kruschke 2014, 340f).

`rope()` performs a simple check for pairwise correlations between parameters, but as there can be collinearity between more than two variables, a first step to check the assumptions of this hypothesis testing is to look at different pair plots. An even more sophisticated check is the projection predictive variable selection (Piironen and Vehtari 2017).

**Strengths and Limitations:**

**Strengths:** Provides information related to the practical relevance of the effects.

**Limitations:** A ROPE range needs to be arbitrarily defined. Sensitive to the scale (the unit) of the predictors. Not sensitive to highly significant effects.

**Note**

There is also a `plot()`-method implemented in the `see-package`.

**References**

Examples

```r
library(bayestestR)

rope(x = rnorm(1000, 0, 0.01), range = c(-0.1, 0.1))
rope(x = rnorm(1000, 0, 1), range = c(-0.1, 0.1))
rope(x = rnorm(1000, 1, 0.01), range = c(-0.1, 0.1))
rope(x = rnorm(1000, 1, 1), ci = c(.90, .95))
## Not run:
library(rstanarm)
model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
rope(model)
rope(model, ci = c(.90, .95))

library(emmeans)
rope(emtrends(model, ~1, "wt"), ci = c(.90, .95))

library(brms)
model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
rope(model)
rope(model, ci = c(.90, .95))

library(brms)
model <- brms::brm(brms::mvbind(mpg, disp) ~ wt + cyl, data = mtcars)
rope(model)
rope(model, ci = c(.90, .95))

library(BayesFactor)
bf <- ttestBF(x = rnorm(100, 1, 1))
rope(bf)
rope(bf, ci = c(.90, .95))
## End(Not run)
```

---

**rope_range**

| Find Default Equivalence (ROPE) Region Bounds |

**Description**

This function attempts at automatically finding suitable "default" values for the Region Of Practical Equivalence (ROPE).

**Usage**

```r
rope_range(x, 
...)
```

## Default S3 method:
rope_range(x, verbose = TRUE, 
...)

Arguments

x A stanreg, brmsfit or BFBayesFactor object.

... Currently not used.

verbose Toggle warnings.

Details

Kruschke (2018) suggests that the region of practical equivalence could be set, by default, to a range from -0.1 to 0.1 of a standardized parameter (negligible effect size according to Cohen, 1988).

- For **linear models (lm)**, this can be generalised to \([-0.1 \times SD_y, 0.1 \times SD_y]\).

  \item For **logistic models**, the parameters expressed in log odds ratio can be converted to standardized difference through the formula \(\pi/\sqrt{3}\), resulting in a range of -0.18 to 0.18.

  \item For **logistic models**, it is strongly recommended to manually specify the rope argument. Currently, the same default is applied that for logistic models.

  \item For **count data**, the residual variance is used. This is a rather experimental threshold and is probably often similar to `-0.1, 0.1`, but should be used with care!

  \item For **t-tests**, the standard deviation of the response is used, similarly to linear models (see above).

  \item For **correlations**, `-0.05, 0.05` is used, i.e., half the value of a negligible correlation as suggested by Cohen's (1988) rules of thumb.

  \item For all other models, `-0.1, 0.1` is used to determine the ROPE limits, but it is strongly advised to specify it manually.

References


Examples

```r
## Not run:
if (require("rstanarm")) {
  model <- stan_glm(
    mpg ~ wt + gear,
    data = mtcars,
    chains = 2,
    iter = 200,
```

```r
```
refresh = 0
}
rope_range(model)

model <- stan_glm(vs ~ mpg, data = mtcars, family = "binomial", refresh = 0)
rope_range(model)
}

if (require("brms")) {
  model <- brm(mpg ~ wt + cyl, data = mtcars)
  rope_range(model)
}

if (require("BayesFactor")) {
  model <- ttestBF(mtcars[mtcars$vs == 1, "mpg"], mtcars[mtcars$vs == 0, "mpg"])
  rope_range(model)
  model <- lmBF(mpg ~ vs, data = mtcars)
  rope_range(model)
}

## End(Not run)

---

**sensitivity_to_prior**  
**Sensitivity to Prior**

**Description**

Computes the sensitivity to priors specification. This represents the proportion of change in some indices when the model is fitted with an antagonistic prior (a prior of same shape located on the opposite of the effect).

**Usage**

```r
sensitivity_to_prior(model, index = "Median", magnitude = 10, ...)
```

**Arguments**

- **model**: A Bayesian model (`stanreg` or `brmsfit`).
- **index**: The indices from which to compute the sensitivity. Can be one or multiple names of the columns returned by `describe_posterior`. The case is important here (e.g., write 'Median' instead of 'median').
- **magnitude**: This represent the magnitude by which to shift the antagonistic prior (to test the sensitivity). For instance, a magnitude of 10 (default) means that the mode will be updated with a prior located at 10 standard deviations from its original location.
- **...**: Arguments passed to or from other methods.
See Also

DescTools

Examples

```r
## Not run:
library(bayestestR)

# rstanarm models
# -----------------------------------------------
if (require("rstanarm")) {
  model <- rstanarm::stan_glm(mpg ~ wt, data = mtcars)
  sensitivity_to_prior(model)
  model <- rstanarm::stan_glm(mpg ~ wt + cyl, data = mtcars)
  sensitivity_to_prior(model, index = c("Median", "MAP"))
}

# brms models
# -----------------------------------------------
if (require("brms")) {
  model <- brms::brm(mpg ~ wt + cyl, data = mtcars)
  sensitivity_to_prior(model)
}
## End(Not run)
```

sexit

Sequential Effect eXistence and sIgnificance Testing (SEXXIT)

Description

The SEXIT is a new framework to describe Bayesian effects, guiding which indices to use. Accordingly, the sexit() function returns the minimal (and optimal) required information to describe models’ parameters under a Bayesian framework. It includes the following indices:

- **Centrality**: the median of the posterior distribution. In probabilistic terms, there is 50% of probability that the effect is higher and lower. See `point_estimate()`.

- **Uncertainty**: the 95% Highest Density Interval (HDI). In probabilistic terms, there is 95% of probability that the effect is within this confidence interval. See `ci()`.

- **Existence**: The probability of direction allows to quantify the certainty by which an effect is positive or negative. It is a critical index to show that an effect of some manipulation is not harmful (for instance in clinical studies) or to assess the direction of a link. See `p_direction()`.

- **Significance**: Once existence is demonstrated with high certainty, we can assess whether the effect is of sufficient size to be considered as significant (i.e., not negligible). This is a useful index to determine which effects are actually important and worthy of discussion in a given process. See `p_significance()`.
- **Size:** Finally, this index gives an idea about the strength of an effect. However, beware, as studies have shown that a big effect size can be also suggestive of low statistical power (see details section).

Usage

```r
sexit(x, significant = "default", large = "default", ci = 0.95, ...)
```

Arguments

- `x` Vector representing a posterior distribution. Can also be a Bayesian model (stanreg, brmsfit or BayesFactor).
- `significant`, `large` The threshold values to use for significant and large probabilities. If left to 'default', will be selected through `sexit_thresholds()`. See the details section below.
- `ci` Value or vector of probability of the (credible) interval - CI (between 0 and 1) to be estimated. Default to .95 (95%).
- `...` Currently not used.

Details

**Rationale:** The assessment of "significance" (in its broadest meaning) is a pervasive issue in science, and its historical index, the p-value, has been strongly criticized and deemed to have played an important role in the replicability crisis. In reaction, more and more scientists have tuned to Bayesian methods, offering an alternative set of tools to answer their questions. However, the Bayesian framework offers a wide variety of possible indices related to "significance", and the debate has been raging about which index is the best, and which one to report.

This situation can lead to the mindless reporting of all possible indices (with the hopes that with that the reader will be satisfied), but often without having the writer understanding and interpreting them. It is indeed complicated to juggle between many indices with complicated definitions and subtle differences.

SEXIT aims at offering a practical framework for Bayesian effects reporting, in which the focus is put on intuitiveness, explicitness and usefulness of the indices’ interpretation. To that end, we suggest a system of description of parameters that would be intuitive, easy to learn and apply, mathematically accurate and useful for taking decision.

Once the thresholds for significance (i.e., the ROPE) and the one for a "large" effect are explicitly defined, the SEXIT framework does not make any interpretation, i.e., it does not label the effects, but just sequentially gives 3 probabilities (of direction, of significance and of being large, respectively) as-is on top of the characteristics of the posterior (using the median and HDI for centrality and uncertainty description). Thus, it provides a lot of information about the posterior distribution (through the mass of different 'sections' of the posterior) in a clear and meaningful way.

**Threshold selection:** One of the most important thing about the SEXIT framework is that it relies on two "arbitrary" thresholds (i.e., that have no absolute meaning). They are the ones related to effect size (an inherently subjective notion), namely the thresholds for significant and large effects. They are set, by default, to .05 and .3 of the standard deviation of the outcome variable (tiny and large effect sizes for correlations according to Funder & Ozer, 2019). However,
these defaults were chosen by lack of a better option, and might not be adapted to your case. Thus, they are to be handled with care, and the chosen thresholds should always be explicitly reported and justified.

- For **linear models (lm)**, this can be generalised to \([0.05 \times SD_y]\) and \([0.3 \times SD_y]\) for significant and large effects, respectively.
- For **logistic models**, the parameters expressed in log odds ratio can be converted to standardised difference through the formula \(\pi/\sqrt{3}\), resulting a threshold of 0.09 and 0.54.
- For other models with **binary outcome**, it is strongly recommended to manually specify the rope argument. Currently, the same default is applied that for logistic models.
- For models from **count data**, the residual variance is used. This is a rather experimental threshold and is probably often similar to 0.05 and 0.3, but should be used with care!
- For **t-tests**, the standard deviation of the response is used, similarly to linear models (see above).
- For **correlations**, 0.05 and 0.3 are used.
- For all other models, 0.05 and 0.3 are used, but it is strongly advised to specify it manually.

**Examples:** The three values for existence, significance and size provide a useful description of the posterior distribution of the effects. Some possible scenarios include:

- The probability of existence is low, but the probability of being large is high: it suggests that the posterior is very wide (covering large territories on both side of 0). The statistical power might be too low, which should warrant any confident conclusion.
- The probability of existence and significance is high, but the probability of being large is very small: it suggests that the effect is, with high confidence, not large (the posterior is mostly contained between the significance and the large thresholds).
- The 3 indices are very low: this suggests that the effect is null with high confidence (the posterior is closely centred around 0).

**Value**

A dataframe and text as attribute.

**References**


**Examples**

```r
## Not run:
library(bayestestR)

s <- sexit(rnorm(1000, -1, 1))
s
print(s, summary = TRUE)
```
sexit_thresholds

Find Effect Size Thresholds

Description
This function attempts at automatically finding suitable default values for a "significant" (i.e., non-negligible) and "large" effect. This is to be used with care, and the chosen threshold should always be explicitly reported and justified. See the detail section in sexit() for more information.

Usage
sexit_thresholds(x, ...)

Arguments
x Vector representing a posterior distribution. Can also be a stanreg or brmsfit model.

... Currently not used.

References

Examples
sexit_thresholds(rnorm(1000))
## Not run:
if (require("rstanarm")) {
  model <- rstanarm::stan_glm(mpg ~ wt * cyl,
    data = mtcars,
    iter = 400, refresh = 0
  )
  s <- sexit(model)
  s
  print(s, summary = TRUE)
}
## End(Not run)
chains = 2,
iter = 200,
refresh = 0
)
sexit_thresholds(model)

model <- stan_glm(vs ~ mpg, data = mtcars, family = "binomial", refresh = 0)
sexit_thresholds(model)
}

if (require("brms")) {
  model <- brm(mpg ~ wt + cyl, data = mtcars)
  sexit_thresholds(model)
}

if (require("BayesFactor")) {
  bf <- ttestBF(x = rnorm(100, 1, 1))
  sexit_thresholds(bf)
}
## End(Not run)

---

### `si`  
Compute Support Intervals

**Description**

A support interval contains only the values of the parameter that predict the observed data better than average, by some degree $k$; these are values of the parameter that are associated with an updating factor greater or equal than $k$. From the perspective of the Savage-Dickey Bayes factor, testing against a point null hypothesis for any value within the support interval will yield a Bayes factor smaller than $1/k$.

**For more info, in particular on specifying correct priors for factors with more than 2 levels, see the Bayes factors vignette.**

**Usage**

```r
si(posterior, prior = NULL, BF = 1, verbose = TRUE, ...)
```

```
## S3 method for class 'numeric'
si(posterior, prior = NULL, BF = 1, verbose = TRUE, ...)
```

```
## S3 method for class 'stanreg'
si(
  posterior,
  prior = NULL,
  BF = 1,
  verbose = TRUE,
)```
effects = c("fixed", "random", "all"),
component = c("conditional", "location", "zi", "zero_inflated", "all",
  "smooth_terms", "sigma", "distributional", "auxiliary"),
parameters = NULL,
...)

## S3 method for class 'brmsfit'
si(
  posterior,
  prior = NULL,
  BF = 1,
  verbose = TRUE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "location", "zi", "zero_inflated", "all",
    "smooth_terms", "sigma", "distributional", "auxiliary"),
  parameters = NULL,
  ...
)

## S3 method for class 'blavaan'
si(
  posterior,
  prior = NULL,
  BF = 1,
  verbose = TRUE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "location", "zi", "zero_inflated", "all",
    "smooth_terms", "sigma", "distributional", "auxiliary"),
  parameters = NULL,
  ...
)

## S3 method for class 'emmGrid'
si(posterior, prior = NULL, BF = 1, verbose = TRUE, ...)

## S3 method for class 'data.frame'
si(posterior, prior = NULL, BF = 1, verbose = TRUE, ...)

Arguments

- `posterior`: A numerical vector, `stanreg / brmsfit` object, `emmGrid` or a data frame - representing a posterior distribution(s) from (see 'Details').
- `prior`: An object representing a prior distribution (see 'Details').
- `BF`: The amount of support required to be included in the support interval.
- `verbose`: Toggle off warnings.
- `...`: Arguments passed to and from other methods. (Can be used to pass arguments to internal `logspline:::logspline()`.)
effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.

component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms models.

parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior_) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

Details
This method is used to compute support intervals based on prior and posterior distributions. For the computation of support intervals, the model priors must be proper priors (at the very least they should be not flat, and it is preferable that they be informative - note that by default, brms::brm() uses flat priors for fixed-effects; see example below).

Choosing a value of BF: The choice of BF (the level of support) depends on what we want our interval to represent:

• A BF = 1 contains values whose credibility is not decreased by observing the data.
• A BF > 1 contains values who received more impressive support from the data.
• A BF < 1 contains values whose credibility has not been impressively decreased by observing the data. Testing against values outside this interval will produce a Bayes factor larger than 1/BF in support of the alternative. E.g., if an SI (BF = 1/3) excludes 0, the Bayes factor against the point-null will be larger than 3.

Value
A data frame containing the lower and upper bounds of the SI.
Note that if the level of requested support is higher than observed in the data, the interval will be [NA, NA].

Setting the correct prior
For the computation of Bayes factors, the model priors must be proper priors (at the very least they should be not flat, and it is preferable that they be informative); As the priors for the alternative get wider, the likelihood of the null value(s) increases, to the extreme that for completely flat priors the null is infinitely more favorable than the alternative (this is called the Jeffreys-Lindley-Bartlett paradox). Thus, you should only ever try (or want) to compute a Bayes factor when you have an informed prior.

(Note that by default, brms::brm() uses flat priors for fixed-effects; See example below.)

It is important to provide the correct prior for meaningful results.

• When posterior is a numerical vector, prior should also be a numerical vector.
• When posterior is a data.frame, prior should also be a data.frame, with matching column order.
• When posterior is a stanreg, brmsfit or other supported Bayesian model:
  – prior can be set to NULL, in which case prior samples are drawn internally.
  – prior can also be a model equivalent to posterior but with samples from the priors only. See `unupdate()`.
  – **Note:** When posterior is a brmsfit_multiple model, prior **must** be provided.

• When posterior is an emmGrid / emm_list object:
  – prior should also be an emmGrid / emm_list object equivalent to posterior but created with a model of priors samples only. See `unupdate()`.
  – prior can also be the original (posterior) model. If so, the function will try to update the emmGrid / emm_list to use the `unupdate()`d prior-model. (*This cannot be done for brmsfit models.*)
  – **Note:** When the emmGrid has undergone any transformations ("log", "response", etc.), or regridding, then prior must be an emmGrid object, as stated above.

**Note**

There is also a `plot()`-method implemented in the see-package.

**References**


**See Also**

Other ci: `bci()`, `ci()`, `cwi()`, `eti()`, `hdi()`, `spi()`

**Examples**

```r
library(bayestestR)

prior <- distribution_normal(1000, mean = 0, sd = 1)
posterior <- distribution_normal(1000, mean = .5, sd = .3)

si(posterior, prior)
## Not run:
# rstanarm models
# ---------------
library(rstanarm)
contrasts(sleep$group) <- contr.orthonorm # see vinette
stan_model <- stan_lmer(extra ~ group + (1 | ID), data = sleep)
si(stan_model)
si(stan_model, BF = 3)

# emmGrid objects
# ---------------
library(emmeans)
group_diff <- pairs(emmeans(stan_model, ~group))
si(group_diff, prior = stan_model)
```
## simulate_correlation

### Data Simulation

**Description**

Simulate data with specific characteristics.

**Usage**

```r
simulate_correlation(n = 100, r = 0.5, mean = 0, sd = 1, names = NULL, ...)
```

### Arguments

- `n` The number of observations to be generated.
- `r` A value or vector corresponding to the desired correlation coefficients.
- `mean` A value or vector corresponding to the mean of the variables.
- `sd` A value or vector corresponding to the SD of the variables.
- `names` A character vector of desired variable names.
- `...` Arguments passed to or from other methods.
- `d` A value or vector corresponding to the desired difference between the groups.

### Examples

```r
# Correlation --------------------------------
data <- simulate_correlation(r = 0.5)
plot(data$V1, data$V2)
```
cor.test(data$V1, data$V2)
summary(lm(V2 ~ V1, data = data))

# Specify mean and SD
data <- simulate_correlation(r = 0.5, n = 50, mean = c(0, 1), sd = c(0.7, 1.7))
cor.test(data$V1, data$V2)
round(c(mean(data$V1), sd(data$V1)), 1)
round(c(mean(data$V2), sd(data$V2)), 1)
summary(lm(V2 ~ V1, data = data))

# Generate multiple variables
cor_matrix <- matrix(c(1.0, 0.2, 0.4,
                        0.2, 1.0, 0.3,
                        0.4, 0.3, 1.0
                      ),
                      nrow = 3
                      )
data <- simulate_correlation(r = cor_matrix, names = c("y", "x1", "x2"))
cor(data)
summary(lm(y ~ x1, data = data))

# t-test --------------------------------
data <- simulate_ttest(n = 30, d = 0.3)
plot(data$V1, data$V0)
round(c(mean(data$V1), sd(data$V1)), 1)
diff(t.test(data$V1 ~ data$V0)$estimate)
summary(lm(V1 ~ V0, data = data))
summary(glm(V0 ~ V1, data = data, family = "binomial"))

# Difference --------------------------------
data <- simulate_difference(n = 30, d = 0.3)
plot(data$V1, data$V0)
round(c(mean(data$V1), sd(data$V1)), 1)
diff(t.test(data$V1 ~ data$V0)$estimate)
summary(lm(V1 ~ V0, data = data))
summary(glm(V0 ~ V1, data = data, family = "binomial"))

---

**simulate_prior**

*Returns Priors of a Model as Empirical Distributions*

**Description**

Transforms priors information to actual distributions.

**Usage**

simulate_prior(model, n = 1000, ...)

simulate_simpson

Simpson's paradox dataset simulation

Description

Simpson's paradox, or the Yule-Simpson effect, is a phenomenon in probability and statistics, in which a trend appears in several different groups of data but disappears or reverses when these groups are combined.

Usage

```r
simulate_simpson(
  n = 100,
  r = 0.5,
  groups = 3,
  difference = 1,
  group_prefix = "G_"
)
```

Arguments

- **n**: The number of observations for each group to be generated (minimum 4).
- **r**: A value or vector corresponding to the desired correlation coefficients.
- **groups**: Number of groups (groups can be participants, clusters, anything).
- **difference**: Difference between groups.
- **group_prefix**: The prefix of the group name (e.g., "G_1", "G_2", "G_3", ...).

See Also

unupdate() for directly sampling from the prior distribution (useful for complex priors and designs).

Examples

```r
## Not run:
library(bayestestR)
if (require("rstanarm")) {
  model <- stan_glm(mpg ~ wt + am, data = mtcars, chains = 1, refresh = 0)
  simulate_prior(model)
}
## End(Not run)
```
Value

A dataset.

Examples

data <- simulate_simpson(n = 10, groups = 5, r = 0.5)

if (require("ggplot2")) {
  ggplot(data, aes(x = V1, y = V2)) +
  geom_point(aes(color = Group)) +
  geom_smooth(aes(color = Group), method = "lm") +
  geom_smooth(method = "lm")
}

spi

Shortest Probability Interval (SPI)

Description

Compute the Shortest Probability Interval (SPI) of posterior distributions. The SPI is a more computationally stable HDI. The implementation is based on the algorithm from the SPIn package.

Usage

spi(x, ...)

## S3 method for class 'numeric'
spi(x, ci = 0.95, verbose = TRUE, ...)

## S3 method for class 'stanreg'
spi(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  component = c("location", "all", "conditional", "smooth_terms", "sigma",
                "distributional", "auxiliary"),
  parameters = NULL,
  verbose = TRUE,
  ...
)

## S3 method for class 'brmsfit'
spi(
  x,
  ci = 0.95,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
spi

parameters = NULL,
verbose = TRUE,
...
)}

Arguments

x Vector representing a posterior distribution, or a data frame of such vectors. Can also be a Bayesian model. bayestestR supports a wide range of models (see, for example, methods("hdi")) and not all of those are documented in the ‘Usage’ section, because methods for other classes mostly resemble the arguments of the .numeric or .data.frame methods.
...
Currently not used.
ci Value or vector of probability of the (credible) interval - CI (between 0 and 1) to be estimated. Default to .95 (95%).
verbose Toggle off warnings.
effects Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
component Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to brms-models.
parameters Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like lp__ or prior__) are filtered by default, so only parameters that typically appear in the summary() are returned. Use parameters to select specific parameters for the output.

Details

The SPI is an alternative method to the HDI (hdi()) to quantify uncertainty of (posterior) distributions. The SPI is said to be more stable than the HDI, because, the “HDI can be noisy (that is, have a high Monte Carlo error)” (Liu et al. 2015). Furthermore, the HDI is sensitive to additional assumptions, in particular assumptions related to the different estimation methods, which can make the HDI less accurate or reliable (see also discussion here).

Value

A data frame with following columns:

- Parameter The model parameter(s), if x is a model-object. If x is a vector, this column is missing.
- CI The probability of the credible interval.
- CI_low, CI_high The lower and upper credible interval limits for the parameters.

Note

The code to compute the SPI was adapted from the SPIin package, and slightly modified to be more robust for Stan models. Thus, credits go to Ying Liu for the original SPI algorithm and R implementation.
References


See Also

Other ci: bci(), ci(), cwi(), eti(), hdi(), si()

Examples

library(bayestestR)

posterior <- rnorm(1000)
spi(posterior)
spi(posterior, ci = c(.80, .89, .95))

df <- data.frame(replicate(4, rnorm(100)))
spi(df)
spi(df, ci = c(.80, .89, .95))
## Not run:
library(rstanarm)
model <- stan_glm(mpg ~ wt + gear, data = mtcars, chains = 2, iter = 200, refresh = 0)
spi(model)
## End(Not run)

---

**weighted_posteriors**

Generate posterior distributions weighted across models

**Description**

Extract posterior samples of parameters, weighted across models. Weighting is done by comparing posterior model probabilities, via bayesfactor_models().

**Usage**

weighted_posteriors(..., prior_odds = NULL, missing = 0, verbose = TRUE)

## S3 method for class 'data.frame'
weighted_posteriors(..., prior_odds = NULL, missing = 0, verbose = TRUE)

## S3 method for class 'stanreg'
weighted_posteriors(
  ...,
  prior_odds = NULL,
  missing = 0,
  verbose = TRUE,
)
weighted_posteriors

```r
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL
)

## S3 method for class 'brmsfit'
weighted_posteriors(
  ..., 
  prior_odds = NULL,
  missing = 0,
  verbose = TRUE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL
)

## S3 method for class 'blavaan'
weighted_posteriors(
  ..., 
  prior_odds = NULL,
  missing = 0,
  verbose = TRUE,
  effects = c("fixed", "random", "all"),
  component = c("conditional", "zi", "zero_inflated", "all"),
  parameters = NULL
)

## S3 method for class 'BFBayesFactor'
weighted_posteriors(
  ..., 
  prior_odds = NULL,
  missing = 0,
  verbose = TRUE,
  iterations = 4000
)
```

Arguments

- `...`: Fitted models (see details), all fit on the same data, or a single BFBayesFactor object.
- `prior_odds`: Optional vector of prior odds for the models compared to the first model (or the denominator, for BFBayesFactor objects). For data.frames, this will be used as the basis of weighting.
- `missing`: An optional numeric value to use if a model does not contain a parameter that appears in other models. Defaults to 0.
- `verbose`: Toggle off warnings.
- `effects`: Should results for fixed effects, random effects or both be returned? Only applies to mixed models. May be abbreviated.
component

Should results for all parameters, parameters for the conditional model or the zero-inflated part of the model be returned? May be abbreviated. Only applies to \texttt{brms}-models.

parameters

Regular expression pattern that describes the parameters that should be returned. Meta-parameters (like \texttt{lp__} or \texttt{prior__}) are filtered by default, so only parameters that typically appear in the \texttt{summary()} are returned. Use parameters to select specific parameters for the output.

iterations

For \texttt{BayesFactor} models, how many posterior samples to draw.

Details

Note that across models some parameters might play different roles. For example, the parameter A plays a different role in the model $Y \sim A + B$ (where it is a main effect) than it does in the model $Y \sim A + B + A:B$ (where it is a simple effect). In many cases centering of predictors (mean subtracting for continuous variables, and effects coding via \texttt{contr.sum} or orthonormal coding via \texttt{contr.orthonorm()} for factors) can reduce this issue. In any case you should be mindful of this issue.

See \texttt{bayesfactor_models()} details for more info on passed models.

Note that for \texttt{BayesFactor} models, posterior samples cannot be generated from intercept only models.

This function is similar in function to \texttt{brms::posterior_average}.

Value

A data frame with posterior distributions (weighted across models).

Note

For \texttt{BayesFactor} < 0.9.12-4.3, in some instances there might be some problems of duplicate columns of random effects in the resulting data frame.

References


See Also

\texttt{bayesfactor_inclusion()} for Bayesian model averaging.
Examples

```r
if (require("rstanarm") && require("see")) {
  stan_m0 <- stan_glm(extra ~ 1,
                     data = sleep,
                     family = gaussian(),
                     refresh = 0,
                     diagnostic_file = file.path(tempdir(), "df0.csv")
  )

  stan_m1 <- stan_glm(extra ~ group,
                     data = sleep,
                     family = gaussian(),
                     refresh = 0,
                     diagnostic_file = file.path(tempdir(), "df1.csv")
  )

  res <- weighted_posteriors(stan_m0, stan_m1)

  plot(eti(res))
}

## With BayesFactor
if (require("BayesFactor")) {
  extra_sleep <- ttestBF(formula = extra ~ group, data = sleep)

  wp <- weighted_posteriors(extra_sleep)

  describe_posterior(extra_sleep, test = NULL)
  describe_posterior(wp$delta, test = NULL) # also considers the null
}

## weighted prediction distributions via data.frames
if (require("rstanarm")) {
  m0 <- stan_glm(
                mpg ~ 1,
                data = mtcars,
                family = gaussian(),
                diagnostic_file = file.path(tempdir(), "df0.csv"),
                refresh = 0
  )

  m1 <- stan_glm(
                mpg ~ carb,
                data = mtcars,
                family = gaussian(),
                diagnostic_file = file.path(tempdir(), "df1.csv"),
                refresh = 0
  )

```
# Predictions:
pred_m0 <- data.frame(posterior_predict(m0))
pred_m1 <- data.frame(posterior_predict(m1))

BFmods <- bayesfactor_models(m0, m1)

wp <- weighted_posteriors(pred_m0, pred_m1,
    prior_odds = as.numeric(BFmods)[2]
)

# look at first 5 prediction intervals
hdi(pred_m0[1:5])
hdi(pred_m1[1:5])
hdi(wp[1:5])  # between, but closer to pred_m1
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