

Package ‘rmsb’

September 26, 2023

Title Bayesian Regression Modeling Strategies

Version 1.0-0

Date 2023-09-26

Description

A Bayesian companion to the 'rms' package, 'rmsb' provides Bayesian model fitting, post-fit estimation, and graphics. It implements Bayesian regression models whose fit objects can be processed by 'rms' functions such as 'contrast()', 'summary()', 'Predict()', 'nomogram()', and 'latex()'. The fitting function currently implemented in the package is 'blrm()' for Bayesian logistic binary and ordinal regression with optional clustering, censoring, and departures from the proportional odds assumption using the partial proportional odds model of Peterson and Harrell (1990) <<https://www.jstor.org/stable/2347760>>.

License GPL (>= 3)

Encoding UTF-8

URL <https://hbiostat.org/R/rmsb/>

RoxygenNote 7.2.3

Biarch true

Depends R (>= 3.4.0), rms (>= 6.7-1)

Imports methods, Rcpp (>= 0.12.0), rstan (>= 2.26.23), Hmisc (>= 4.3-0), survival (>= 3.1-12), ggplot2, MASS, cluster, digest, knitr, loo

LinkingTo BH (>= 1.66.0), Rcpp (>= 0.12.0), RcppEigen (>= 0.3.3.3.0), RcppParallel (>= 5.0.1), rstan (>= 2.18.1), StanHeaders (>= 2.18.0)

Suggests cmdstanr, bayesplot, mice

Additional_repositories <https://mc-stan.org/r-packages/>

SystemRequirements GNU make

NeedsCompilation yes

Author Frank Harrell [aut, cre] (<<https://orcid.org/0000-0002-8271-5493>>),
Ben Goodrich [ctb] (contributed Stan code),
Ben Bolker [ctb] (wrote original code that is folded into the

pdensityContour function),
 Doug Bates [ctb] (write original code for highest posterior density
 interval that is folded into the HPDint function)

Maintainer Frank Harrell <fh@fharrell.com>

Repository CRAN

Date/Publication 2023-09-26 13:10:02 UTC

R topics documented:

rmsb-package	3
as.data.frame.Ocens	3
blrm	4
blrmStats	9
cluster	11
coef.rmsb	11
compareBmods	12
distSym	13
ExProb.blrm	13
getParamCoef	14
HPDint	15
Mean.blrm	15
Ocens	16
pdensityContour	17
plot.PostF	18
plot.rmsb	19
PostF	20
predict.blrm	21
print.blrm	23
print.blrmStats	24
print.predict.blrm	25
print.rmsb	25
Quantile.blrm	26
selectedQr	27
stackMI	28
stanDx	29
stanDxplot	30
stanGet	31
tauFetch	32
vcov.rmsb	32
[.Ocens	33

Index

35

rmsb-package

*The 'rmsb' package.***Description**

Regression Modeling Strategies Bayesian

The **rmsb** package is an appendage to the **rms** package that implements Bayesian regression models whose fit objects can be processed by **rms** functions such as `contrast`, `summary`, `Predict`, `nomogram`, and `latex`. The fitting function currently implemented in the package is `blrm` for Bayesian logistic binary and ordinal regression with optional clustering, censoring, and departures from the proportional odds assumption using the partial proportional odds model of Peterson and Harrell (1990).

References

Stan Development Team (2020). RStan: the R interface to Stan. R package version 2.19.3. <https://mc-stan.org>

See Also

- <https://hbiostat.org/R/rmsb/> for the package's main web page
- <https://hbiostat.org/R/examples/blrm/blrm.html> for a vignette with many examples of using the `blrm` function

as.data.frame.Ocens

*Convert Ocens Object to Data Frame to Facilitate Subset***Description**

Converts an `Ocens` object to a data frame so that subsetting will preserve all needed attributes

Usage

```
## S3 method for class 'Ocens'
as.data.frame(x, row.names = NULL, optional = FALSE, ...)
```

Arguments

<code>x</code>	an <code>Ocens</code> object
<code>row.names</code>	optional vector of row names
<code>optional</code>	set to <code>TRUE</code> if needed
<code>...</code>	ignored

Value

data frame containing a 2-column integer matrix with attributes

Author(s)

Frank Harrell

blrm

Bayesian Binary and Ordinal Logistic Regression

Description

Uses `rstan` with pre-compiled Stan code, or `cmdstan` to get posterior draws of parameters from a binary logistic or proportional odds semiparametric ordinal logistic model. The Stan code internally using the qr decomposition on the design matrix so that highly collinear columns of the matrix do not hinder the posterior sampling. The parameters are transformed back to the original scale before returning results to R. Design matrix columns are centered before running Stan, so Stan diagnostic output will have the intercept terms shifted but the results of `blrm()` for intercepts are for the original uncentered data. The only prior distributions for regression betas are normal with mean zero. Priors are specified on contrasts so that they can be specified on a meaningful scale and so that more complex patterns can be imposed. Parameters that are not involved in any contrasts in `pcontrast` have non-informative priors. Contrasts are automatically converted to the QR space used in Stan code.

Usage

```
blrm(
  formula,
  ppo = NULL,
  cppo = NULL,
  keepsep = NULL,
  data = environment(formula),
  subset,
  na.action = na.delete,
  priorsdppo = rep(100, ppo),
  iprior = 0,
  conc = 1/(0.8 + 0.35 * max(k, 3)),
  ascale = 1,
  psigma = 1,
  rsdmean = if (psigma == 1) 0 else 1,
  rsdsd = 1,
  normcppo = TRUE,
  pcontrast = NULL,
  backend = c("rstan", "cmdstan"),
  iter = 2000,
  warmup = iter/2,
  chains = 4,
```

```

refresh = 0,
progress = if (refresh > 0) "stan-progress.txt" else "",
x = TRUE,
y = TRUE,
loo = n <= 1000,
ppairs = NULL,
method = c("both", "sampling", "optimizing"),
inito = if (length(ppo) 0 else "random",
inits = inito,
standata = FALSE,
file = NULL,
debug = FALSE,
sampling.args = NULL,
...
)

```

Arguments

formula	a R formula object that can use rms package enhancements such as the restricted interaction operator
ppo	formula specifying the model predictors for which proportional odds is not assumed
cppo	a function that if present causes a constrained partial PO model to be fit. The function specifies the values in the Gamma vector in Peterson and Harrell (1990) equation (6). To make posterior sampling better behaved, the function should be scaled and centered. This is done by wrapping cppo in a function that scales the cppo result before return the vector value. See the normcco argument for how to prevent this. The default normalization is based on the mean and standard deviation of the function values over the distribution of observed Y. For getting predicted values and estimates post-blrm(), cppo must not reference any functions that are not available at such later times.
keepsep	a single character string containing a regular expression applied to design matrix column names, specifying which columns for non-proportional odds terms are not to be QR-orthonormalized, so that priors for those columns apply to the original parameters. This does not apply to the main part of the model. keepsep is useful for treatment and treatment interaction terms. For example keepsep='treat' will keep separate all design matrix columns containing 'treat' in their names. Some characters such as the caret used in polynomial regression terms will need to be escaped by a double backslash.
data	a data frame; defaults to using objects from the calling environment
subset	a logical vector or integer subscript vector specifying which subset of data should be used
na.action	default is na.delete to remove missings and report on them
priorsdppo	vector of prior standard deviations for non-proportional odds parameters. The last element is the only one for which the SD corresponds to the original data scale.

<code>iprior</code>	specifies whether to use a Dirichlet distribution for the cell probabilities, which induce a more complex prior distribution for the intercepts (<code>iprior=0</code> , the default), non-informative priors (<code>iprior=1</code>) directly on the intercept parameters, or to directly use a t-distribution with 3 d.f. and scale parameter <code>ascale</code> (<code>iprior=2</code>).
<code>conc</code>	the Dirichlet distribution concentration parameter for the prior distribution of cell probabilities at covariate means. The default is the reciprocal of $0.8 + 0.35 \max(k, 3)$ where k is the number of Y categories. The default is chosen to make the posterior mean of the intercepts more closely match the MLE. For optimizing, the concentration parameter is always 1.0 to obtain results very close to the MLE for providing the posterior mode.
<code>ascale</code>	scale parameter for the t-distribution for priors for the intercepts if <code>iprior=2</code> , defaulting to 1.0
<code>psigma</code>	defaults to 1 for a half-t distribution with 4 d.f., location parameter <code>rsdmean</code> and scale parameter <code>rsdsd</code> . Set <code>psigma=2</code> to use the exponential distribution.
<code>rsdmean</code>	the assumed mean of the prior distribution of the standard deviation of random effects. When <code>psigma=2</code> this is the mean of an exponential distribution and defaults to 1. When <code>psigma=1</code> this is the mean of the half-t distribution and defaults to zero.
<code>rsdsd</code>	applies only to <code>psigma=1</code> and is the scale parameter for the half t distribution for the SD of random effects, defaulting to 1.
<code>normcppo</code>	set to <code>FALSE</code> to leave the <code>cppo</code> function as-is without automatically centering and scaling the result
<code>pcontrast</code>	a list specifying contrasts that are to be given Gaussian prior distributions. The predictor combinations specified in <code>pcontrast</code> are run through <code>rms::gendata()</code> so that contrasts are specified in units of original variables, and unspecified variables are set to medians or modes as saved by <code>rms::datadist()</code> . Thanks to Stan, putting priors on combinations and transformations of model parameters has the same effect of putting different priors on the original parameters without figuring out how to do that. The syntax used here allows specification of differences, double differences (e.g., interactions or nonlinearity), triple differences (e.g., to put constraints on nonlinear interactions), etc. The requested predictor combinations must be named so they may be referred to inside <code>contrast</code> . The syntax is <code>pcontrast=list(..., contrast=expression(...), mu=, sd=, weights=, ycut=, expand=)</code> . <code>...</code> denotes one or more <code>list()</code> s with predictor combinations, and each <code>list()</code> must be named, e.g., <code>pcontrast=list(c1=list(sex='female'), c2=list(sex='male'))</code> to set up for a female - male contrast specified as <code>contrast=expression(c1 - c2)</code> . The <code>c1 - c2</code> subtraction will operate on the design matrices generated by the covariate settings in the <code>list()</code> s. For <code>weights</code> , <code>ycut</code> , <code>expand</code> see <code>rms::Xcontrast()</code> and <code>rms::contrast.rms()</code> . <code>mu</code> is a vector of prior means associated with the rows of the stacked contrasts, and <code>sd</code> is a corresponding vector of Gaussian prior SDs. When <code>mu</code> is not given it defaults to 0.0, and <code>sd</code> defaults to 100.0. Values of <code>mu</code> and/or <code>sd</code> are repeated to the number of contrasts if they are of length 1. Full examples are given here .
<code>backend</code>	set to <code>cmdstan</code> to use <code>cmdstan</code> through the R <code>cmdstanr</code> package instead of the default <code>rstan</code> . You can also specify this with a global option <code>rmsb.backend</code> .
<code>iter</code>	number of posterior samples per chain for <code>rstan::sampling()</code> to run, counting warmups

warmup	number of warmup iterations to discard. Default is <code>iter/2</code> .
chains	number of separate chains to run
refresh	see <code>rstan::sampling()</code> and <code>cmdstanr::sample()</code> . The default is 0, indicating that no progress notes are output. If <code>refresh > 0</code> and <code>progress</code> is not <code>''</code> , progress output will be appended to file <code>progress</code> . The default file name is <code>'stan-progress.txt'</code> .
progress	see <code>refresh</code> . Defaults to <code>''</code> if <code>refresh = 0</code> . Note: If running interactively but not under RStudio, <code>rstan</code> will open a browser window for monitoring progress.
x	set to <code>FALSE</code> to not store the design matrix in the fit. <code>x=TRUE</code> is needed if running <code>blrmStats</code> for example.
y	set to <code>FALSE</code> to not store the response variable in the fit
loo	set to <code>FALSE</code> to not run loo and store its result as object <code>loo</code> in the returned object. <code>loo</code> defaults to <code>FALSE</code> if the sample size is greater than 1000, as loo requires the per-observation likelihood components, which creates a matrix <code>N</code> times the number of posterior draws.
ppairs	set to a file name to run <code>rstan pairs</code> or, if <code>backend='cmdstan'</code> <code>bayesplot::mcmc_pairs</code> and store the resulting png plot there. Set to <code>TRUE</code> instead to directly plot these diagnostics. The default is not to run pair plots.
method	set to <code>'optimizing'</code> to run the Stan optimizer and not do posterior sampling, <code>'both'</code> (the default) to run both the optimizer and posterior sampling, or <code>'sampling'</code> to run only the posterior sampling and not compute posterior modes. Running <code>optimizing</code> is a way to obtain maximum likelihood estimates and allows one to quickly study the effect of changing the prior distributions. When <code>method='optimizing'</code> is used the result returned is not a standard <code>blrm()</code> object but is instead the parameter estimates, <code>-2 log likelihood</code> , and optionally the Hessian matrix (if you specify <code>hessian=TRUE</code> in ...). When <code>method='both'</code> is used, <code>rstan::sampling()</code> and <code>rstan::optimizing()</code> are both run, and parameter estimates (posterior modes) from <code>optimizing</code> are stored in a matrix <code>param</code> in the fit object, which also contains the posterior means and medians, and other results from <code>optimizing</code> are stored in object <code>opt</code> in the <code>blrm()</code> fit object. When random effects are present, <code>method</code> is automatically set to <code>'sampling'</code> as maximum likelihood estimates without marginalizing over the random effects do not make sense.
inito	initial value for optimization. The default is the <code>rstan</code> default <code>'random'</code> . Frequently specifying <code>init=0</code> will benefit when the number of distinct <code>Y</code> categories grows or when using <code>ppo</code> hence 0 is the default for that.
inits	initial value for sampling, defaults to <code>inito</code>
standata	set to <code>TRUE</code> to return the Stan data list and not run the model
file	a file name for a <code>saveRDS</code> -created file containing or to contain the saved fit object. If <code>file</code> is specified and the file does not exist, it will be created right before the fit object is returned, less the large <code>rstan</code> object. If the file already exists, its stored <code>md5 hash string datahash</code> fit object component is retrieved and compared to that of the current <code>rstan</code> inputs. If the data to be sent to <code>rstan</code> , the priors, and all sampling and optimization options and stan code are identical, the previously stored fit object is immediately returned and no new calculations are done.

<code>debug</code>	set to TRUE to output timing and progress information to <code>/tmp/debug.txt</code>
<code>sampling.args</code>	a list containing parameters to pass to <code>rstan::sampling()</code> or to the <code>rcmdstan</code> <code>sample</code> function, other than these arguments: <code>iter</code> , <code>warmup</code> , <code>chains</code> , <code>refresh</code> , <code>init</code> which are already arguments to <code>blrm</code>
<code>...</code>	passed to <code>rstan::optimizing()</code> or the <code>rcmdstan</code> <code>optimizing</code> function. The <code>seed</code> parameter is a popular example.

Details

The partial proportional odds model of Peterson and Harrell (1990) is implemented, and is invoked when the user specifies a second model formula as the `ppo` argument. This formula has no left-hand-side variable, and has right-side variables that are a subset of those in formula specifying for which predictors the proportional odds assumption is relaxed.

The Peterson and Harrell (1990) constrained partial proportional odds is also implemented, and is usually preferred to the above unconstrained PPO model as it adds a vector of coefficients instead of a matrix of coefficients. In the constrained PPO model the user provides a function `cppo` that computes a score for all observed values of the dependent variable. For example with a discrete ordinal outcome `cppo` may return a value of 1.0 for a specific value of `Y` and zero otherwise. That will result in a departure from the proportional odds assumption for just that one level of `Y`. The value returned by `cppo` at the lowest `Y` value is never used in any case.

`blrm()` also handles single-level hierarchical random effects models for the case when there are repeated measurements per subject which are reflected as random intercepts, and a different experimental model that allows for AR(1) serial correlation within subject. For both setups, a `cluster` term in the model signals the existence of subject-specific random effects.

When using the `cmdstan` backend, `cmdstanr` will need to compile the Stan code once per computer, only recompiling the code when the Stan source code changes. By default the compiled code is stored in directory `.rmsb` under your home directory. Specify `options(rmsbdir=)` to specify a different location. You should specify `rmsbdir` to be in a project-specific location if you want to archive code for old projects.

If you want to run MCMC sampling even when no inputs or Stan code have changed, i.e., to use a different random number seed for the sampling process, remove the file before running `blrm`.

See [here](#) and [here](#) for multiple examples with results.

Value

an `rms` fit object of class `blrm`, `rmsb`, `rms` that also contains `rstan` or `cmdstanr` results under the name `rstan`. In the `rstan` results, which are also used to produce diagnostics, the intercepts are shifted because of the centering of columns of the design matrix done by `blrm()`. With `method='optimizing'` a class-less list is return with these elements: `coefficients` (MLEs), `beta` (non-intercept parameters on the QR decomposition scale), `deviance` ($-2 \log$ likelihood), `return_code` (see `rstan::optimizing()`), and, if you specified `hessian=TRUE` to `blrm()`, the Hessian matrix. To learn about the scaling of orthogonalized QR design matrix columns, look at the `xqrstd` object in the returned object. This is the vector of SDs for all the columns of the transformed matrix. Those kept out by the `keepsep` argument will have their original SDs. The returned element `sampling_time` is the elapsed time for running posterior samplers, in seconds. This will be just a little more than the time for running one CPU core for one chain.

Author(s)

Frank Harrell and Ben Goodrich

See Also

[print.blrm\(\)](#), [blrmStats\(\)](#), [stanDx\(\)](#), [stanGet\(\)](#), [coef.rmsb\(\)](#), [vcov.rmsb\(\)](#), [print.rmsb\(\)](#), [coef.rmsb\(\)](#)

Examples

```
## Not run:
getHdata(titanic3)
dd <- datadist(titanic3); options(datadist='dd')
f <- blrm(survived ~ (rcs(age, 5) + sex + pclass)^2, data=titanic3)
f # model summary using print.blrm
coef(f) # compute posterior mean parameter values
coef(f, 'median') # compute posterior median values
stanDx(f) # print basic Stan diagnostics
s <- stanGet(f) # extract rstan object from fit
plot(s, pars=f$betas) # Stan posteriors for beta parameters
stanDxplot(s) # Stan diagnostic plots by chain
blrmStats(f) # more details about predictive accuracy measures
ggplot(Predict(...)) # standard rms output
summary(f, ...) # invokes summary.rms
contrast(f, ...) # contrast.rms computes HPD intervals
plot(nomogram(f, ...)) # plot nomogram using posterior mean parameters

# Fit a random effects model to handle multiple observations per
# subject ID using cmdstan
# options(rmsb.backend='cmdstan')
f <- blrm(outcome ~ rcs(age, 5) + sex + cluster(id), data=mydata)

## End(Not run)
```

blrmStats

*Compute Indexes of Predictive Accuracy and Their Uncertainties***Description**

For a binary or ordinal logistic regression fit from [blrm\(\)](#), computes several indexes of predictive accuracy along with highest posterior density intervals for them. Optionally plots their posterior densities. When there are more than two levels of the outcome variable, computes Somers' Dxy and c-index on a random sample of 10,000 observations.

Usage

```
blrmStats(fit, ns = 400, prob = 0.95, pl = FALSE, dist = c("density", "hist"))
```

Arguments

<code>fit</code>	an object produced by <code>blrm()</code>
<code>ns</code>	number of posterior draws to use in the calculations (default is 400)
<code>prob</code>	HPD interval probability (default is 0.95)
<code>pl</code>	set to TRUE to plot the posterior densities using base graphics
<code>dist</code>	if <code>pl</code> is TRUE specifies whether to plot the density estimate (the default) or a histogram

Value

list of class `blrmStats` whose most important element is `Stats`. The indexes computed are defined below, with `gp`, `B`, `EV`, and `vp` computed using the intercept corresponding to the median value of `Y`. See <https://fharrell.com/post/addvalue> for more information.

"Dxy" Somers' Dxy rank correlation between predicted and observed. The concordance probability (c-index; AUROC in the binary `Y` case) may be obtained from the relationship $Dxy=2(c-0.5)$.

"g" Gini's mean difference: the average absolute difference over all pairs of linear predictor values

"gp" Gini's mean difference on the predicted probability scale

"B" Brier score

"EV" explained variation

"v" variance of linear predictor

"vp" variable of estimated probabilities

Author(s)

Frank Harrell

See Also

`Hmisc::rcorr.cens()`

Examples

```
## Not run:
f <- blrm(...)
blrmStats(f, pl=TRUE) # print and plot

## End(Not run)
```

cluster	<i>cluster</i>
---------	----------------

Description

Cluster Function for Random Effects

Usage

```
cluster(x)
```

Arguments

x a vector representing a categorical vector

Details

Used by blrm to signal a categorical variable to generate random effects.

Value

x unchanged

Author(s)

Frank Harrell

coef.rmsb	<i>Extract Bayesian Summary of Coefficients</i>
-----------	---

Description

Computes either the posterior mean (default), posterior median, or posterior mode of the parameters in an rms Bayesian regression model

Usage

```
## S3 method for class 'rmsb'
coef(object, stat = c("mean", "median", "mode"), ...)
```

Arguments

object an object created by an rms package Bayesian fitting function
stat name of measure of posterior distribution central tendency to compute
... ignored

Value

a vector of intercepts and regression coefficients

Author(s)

Frank Harrell

Examples

```
## Not run:  
f <- blrm(...)  
coef(f, stat='mode')  
  
## End(Not run)
```

compareBmods

Compare Bayesian Model Fits

Description

Uses `loo::loo_model_weights()` to compare a series of models such as those created with `blrm()`

Usage

```
compareBmods(..., method = "stacking", r_eff_list = NULL)
```

Arguments

<code>...</code>	a series of model fits
<code>method</code>	see <code>loo::loo_model_weights()</code>
<code>r_eff_list</code>	see <code>loo::loo_model_weights()</code>

Value

a `loo::loo_model_weights()` object

Author(s)

Frank Harrell

distSym	<i>Distribution Symmetry Measure</i>
---------	--------------------------------------

Description

From a sample from a distribution computes a symmetry measure. By default it is the gap between the mean and the 0.95 quantile divided by the gap between the 0.05 quantile and the mean.

Usage

```
distSym(x, prob = 0.9, na.rm = FALSE)
```

Arguments

x	a numeric vector representing a sample from a continuous distribution
prob	quantile interval coverage
na.rm	set to TRUE to remove NAs before proceeding.

Value

a scalar with a value of 1.0 indicating symmetry

Author(s)

Frank Harrell

ExProb.blrm	<i>Function Generator for Exceedance Probabilities for blrm()</i>
-------------	---

Description

For a `blrm()` object generates a function for computing the estimates of the function $\text{Prob}(Y \geq y)$ given one or more values of the linear predictor using the reference (median) intercept. This function can optionally be evaluated at only a set of user-specified y values, otherwise a right-step function is returned. There is a plot method for plotting the step functions, and if more than one linear predictor was evaluated multiple step functions are drawn. `ExProb` is especially useful for `nomogram()`. The linear predictor argument is a posterior summarized linear predictor lp (e.g. using posterior mean of intercepts and slopes) computed at the reference intercept. `lptau` must be provided when call the created function if the model is a partial proportional odds model.

Usage

```
## S3 method for class 'blrm'
ExProb(object, posterior.summary = c("mean", "median"), ...)
```

Arguments

object a `blrm()` fit
 posterior.summary defaults to posterior mean; may also specify "median". Must be consistent with the summary used when creating `lp`.
 ... ignored

Value

an R function

Author(s)

Frank Harrell

getParamCoef *Get a Bayesian Parameter Vector Summary*

Description

Retrieves posterior mean, median, or mode (if available)

Usage

```
getParamCoef(
  fit,
  posterior.summary = c("mean", "median", "mode"),
  what = c("both", "betas", "taus")
)
```

Arguments

fit a Bayesian model fit from `rmsb`
 posterior.summary which summary statistic (Bayesian point estimate) to fetch
 what specifies which coefficients to include. Default is all. Specify `what="betas"` to include only intercepts and betas if the model is a partial proportional odds model (i.e., exclude the tau parameters). Specify `what="taus"` to include only the tau parameters.

Value

vector of regression coefficients

Author(s)

Frank Harrell

HPDint	<i>Highest Posterior Density Interval</i>
--------	---

Description

Adapts code from `coda::HPDinterval()` to compute a highest posterior density interval from posterior samples for a single parameter. Quoting from the coda help file, for each parameter the interval is constructed from the empirical cdf of the sample as the shortest interval for which the difference in the ecdf values of the endpoints is the nominal probability. Assuming that the distribution is not severely multimodal, this is the HPD interval.

Usage

```
HPDint(x, prob = 0.95)
```

Arguments

x	a vector of posterior draws
prob	desired probability coverage

Value

a 2-vector with elements Lower and Upper

Author(s)

Douglas Bates and Frank Harrell

Mean.blrm	<i>Function Generator for Mean Y for blrm()</i>
-----------	---

Description

Creates a function to turn a posterior summarized linear predictor lp (e.g. using posterior mean of intercepts and slopes) computed at the reference intercept into e.g. an estimate of mean Y using the posterior mean of all the intercept. `lptau` must be provided when call the created function if the model is a partial proportional odds model.

Usage

```
## S3 method for class 'blrm'
Mean(object, codes = FALSE, posterior.summary = c("mean", "median"), ...)
```

Arguments

object	a <code>blrm()</code> fit
codes	if TRUE, use the integer codes 1, 2, ..., k for the k -level response in computing the predicted mean response.
posterior.summary	defaults to posterior mean; may also specify "median". Must be consistent with the summary used when creating <code>lp</code> .
...	ignored

Value

an R function

Author(s)

Frank Harrell

Ocens

Censored Ordinal Variable

Description

Creates a 2-column integer matrix that handles left- right- and interval-censored ordinal or continuous values for use in `blrm()`. A pair of values `[a, b]` represents an interval-censored value known to be in the interval `[a, b]` inclusive of `a` and `b`. It is assumed that all distinct values are observed as uncensored for at least one observation. When both input variables are factors it is assumed that the one with the higher number of levels is the one that correctly specifies the order of levels, and that the other variable does not contain any additional levels. If the variables are not factors it is assumed their original values provide the orderings. Since all values that form the left or right endpoints of an interval censored value must be represented in the data, a left-censored point is coded as `a=1` and a right-censored point is coded as `b` equal to the maximum observed value. If the maximum observed value is not really the maximum possible value, everything still works except that predictions involving values above the highest observed value cannot be made. As with most censored-data methods, `blrm()` assumes that censoring is independent of the response variable values that would have been measured had censoring not occurred.

Usage

```
Ocens(a, b = a)
```

Arguments

a	vector representing a factor, numeric, or alphabetically ordered character strings
b	like a. If omitted, it copies a, representing nothing but uncensored values

Value

a 2-column integer matrix of class "Ocens" with an attribute levels (ordered). When the original variables were factors, these are factor levels, otherwise are numerically or alphabetically sorted distinct (over a and b combined) values. When the variables are not factors and are numeric, another attribute median is also returned. This is the median of the uncensored values. When the variables are factor or character, the median of the integer versions of variables for uncensored observations is returned as attribute mid. A final attribute freq is the vector of frequencies of occurrences of all uncensored values. freq aligns with levels.

Author(s)

Frank Harrell

pdensityContour

Bivariate Posterior Contour

Description

Computes coordinates of a highest density contour containing a given probability volume given a sample from a continuous bivariate distribution, and optionally plots. The default method assumes an elliptical shape, but one can optionally use a kernel density estimator. Code adapted from embbook::HPDregionplot. See <https://www.sumsar.net/blog/2014/11/how-to-summarize-a-2d-posterior-using>

Usage

```
pdensityContour(
  x,
  y,
  method = c("ellipse", "kernel"),
  prob = 0.95,
  otherprob = c(0.01, 0.1, 0.25, 0.5, 0.75, 0.9),
  h = c(1.3 * MASS::bandwidth.nrd(x), 1.3 * MASS::bandwidth.nrd(y)),
  n = 70,
  pl = FALSE
)
```

Arguments

x	a numeric vector
y	a numeric vector the same length of x
method	defaults to 'ellipse', can be set to 'kernel'
prob	main probability coverage (the only one for method='ellipse')
otherprob	vector of other probability coverages for method='kernel'
h	vector of bandwidths for x and y. See <code>MASS::kde2d()</code> .
n	number of grid points in each direction, defaulting to normal reference bandwidth (see <code>bandwidth.nrd</code>).
pl	set to TRUE to plot contours

Value

a 2-column matrix with x and y coordinates unless p1=TRUE in which case a ggplot2 graphic is returned

Author(s)

Ben Bolker and Frank Harrell

plot.PostF

Plot Posterior Density of PostF

Description

Computes highest posterior density and posterior mean and median as vertical lines, and plots these on the density function. You can transform the posterior draws while plotting.

Usage

```
## S3 method for class 'PostF'
plot(
  x,
  ...,
  cint = 0.95,
  label = NULL,
  type = c("linetype", "facet"),
  ltitle = ""
)
```

Arguments

x	result of running a function created by PostF
...	other results created by such functions
cint	interval probability
label	x-axis label if not the expression originally evaluated. When more than one result is plotted, label is a vector of character strings, one for each result.
type	when plotting more than one result specifies whether to make one plot distinguishing results by line type, or whether to make separate panels
ltitle	used of type='linetype' to specify name of legend for the line types

Value

ggplot2 object

Author(s)

Frank Harrell

Description

For an rms Bayesian fit object, plots posterior densities for selected parameters along with posterior mode, mean, median, and highest posterior density interval. If the fit was produced by `stackMI` the density represents the distribution after stacking the posterior draws over imputations, and the per-imputation density is also drawn as pale curves. If exactly two parameters are being plotted and `bivar=TRUE`, highest bivariate posterior density contours are plotted instead, for a variety of prob values including the one specified, using

Usage

```
## S3 method for class 'rmsb'
plot(
  x,
  which = NULL,
  nrow = NULL,
  ncol = NULL,
  prob = 0.95,
  bivar = FALSE,
  bivarmethod = c("ellipse", "kernel"),
  ...
)
```

Arguments

<code>x</code>	an rms Bayesian fit object
<code>which</code>	names of parameters to plot, defaulting to all non-intercepts. Can instead be a vector of integers.
<code>nrow</code>	number of rows of plots
<code>ncol</code>	number of columns of plots
<code>prob</code>	probability for HPD interval
<code>bivar</code>	set to TRUE to plot bivariate density contours instead of univariate results (ignored if the number of parameters plotted is not exactly two)
<code>bivarmethod</code>	passed as method argument to <code>pdensityContour</code>
<code>...</code>	passed to <code>pdensityContour</code>

Value

ggplot2 object

Author(s)

Frank Harrell

Description

From a Bayesian fit object such as that from `blrm()` generates an R function for evaluating the probability that an assertion is true. The probability, within simulation error, is the proportion of times the assertion is true over the posterior draws. If the assertion does not evaluate to a logical or 0/1 quantity, it is taken as a continuous derived parameter and the vector of draws for that parameter is returned and can be passed to the PostF plot method. PostF can also be used on objects created by `contrast.rms`

Usage

```
PostF(fit, name = c("short", "orig"), pr = FALSE)
```

Arguments

<code>fit</code>	a Bayesian fit or <code>contrast.rms</code> object
<code>name</code>	specifies whether assertions will refer to shortened parameter names (the default) or original names. Shorted names are of the form <code>a1, ..., ak</code> where <code>k</code> is the number of intercepts in the model, and <code>b1, ..., bp</code> where <code>p</code> is the number of non-intercepts. When using original names that are not legal R variable names, you must enclose them in backticks. For contrast objects, <code>name</code> is ignored and you must use contrast names. The <code>cnames</code> argument to <code>contrast.rms</code> is handy for assigning your own names.
<code>pr</code>	set to TRUE to have a table of short names and original names printed when <code>name='short'</code> . For contrasts the contrast names are printed if <code>pr=TRUE</code> .

Value

an R function

Author(s)

Frank Harrell

Examples

```
## Not run:
f <- blrm(y ~ age + sex)
P <- PostF(f)
P(b2 > 0)      # Model is a1 + b1*age + b2*(sex == 'male')
P(b1 < 0 & b2 > 0) # Post prob of a compound assertion
# To compute probabilities using original parameter names:
P <- PostF(f, name='orig')
P(age < 0)     # Post prob of negative age effect
P(`sex=male` > 0)
```

```

f <- blrm(y ~ sex + pol(age, 2))
P <- PostF(f)
# Compute and plot posterior density of the vertex of the
# quadratic age effect
plot(P(-b2 / (2 * b3)))

# The following would be useful in age and sex interacted
k <- contrast(f, list(age=c(30, 50), sex='male'),
              list(age=c(30, 50), sex='female'),
              cnames=c('age 30 M-F', 'age 50 M-F'))
P <- PostF(k)
P(`age 30 M-F` > 0 & `age 50 M-F` > 0)
##'
## End(Not run)

```

predict.blrm

Make predictions from a `blrm()` fit

Description

Predict method for `blrm()` objects

Usage

```

## S3 method for class 'blrm'
predict(
  object,
  ...,
  kint = NULL,
  ycut = NULL,
  zcppo = TRUE,
  fun = NULL,
  funint = TRUE,
  type = c("lp", "fitted", "fitted.ind", "mean", "x", "data.frame", "terms", "cterms",
           "ccterms", "adjto", "adjto.data.frame", "model.frame"),
  se.fit = FALSE,
  codes = FALSE,
  posterior.summary = c("mean", "median", "all"),
  cint = 0.95
)

```

Arguments

`object`, `...`, `type`, `se.fit`, `codes`
 see [predict.lrm\(\)](#)

`kint` This is only useful in a multiple intercept model such as the ordinal logistic model. There to use to second of three intercepts, for example, specify `kint=2`. The default is the middle intercept corresponding to the median y . You can

	specify <code>ycut</code> instead, and the intercept corresponding to $Y \geq ycut$ will be used for <code>kint</code> .
<code>ycut</code>	for an ordinal model specifies the Y cutoff to use in evaluating departures from proportional odds, when the constrained partial proportional odds model is used. When omitted, <code>ycut</code> is implied by <code>kint</code> . The only time it is absolutely mandatory to specify <code>ycut</code> is when computing an effect (e.g., odds ratio) at a level of the response variable that did not occur in the data. This would only occur when the <code>cppo</code> function given to <code>blrm</code> is a continuous function. If <code>type='x'</code> and neither <code>kint</code> nor <code>ycut</code> are given, the partial PO part of the design matrix is not multiplied by the <code>cppo</code> function. If <code>type='x'</code> , the number of predicted observations is 1, <code>ycut</code> is longer than 1, and <code>zcppo</code> is TRUE, predictions are duplicated to the length of <code>ycut</code> as it is assumed that the user wants to see the effect of varying <code>ycut</code> , e.g., to see cutoff-specific odds ratios.
<code>zcppo</code>	applies only to <code>type='x'</code> for a constrained partial PO model. Set to FALSE to prevent multiplication of Z matrix by <code>cppo(ycut)</code> .
<code>fun</code>	a function to evaluate on the linear predictor, e.g. a function created by <code>Mean()</code> or <code>Quantile()</code>
<code>funint</code>	set to FALSE if <code>fun</code> is not a function such as the result of <code>Mean()</code> , <code>Quantile()</code> , or <code>ExProb()</code> that contains an intercepts argument
<code>posterior.summary</code>	set to 'median' or 'mode' to use posterior median/mode instead of mean. For some types set to 'all' to compute the needed quantity for all posterior draws, and return one more dimension in the array.
<code>cint</code>	probability for highest posterior density interval. Set to FALSE to suppress calculation of the interval.

Value

a data frame, matrix, or vector with posterior summaries for the requested quantity, plus an attribute 'draws' that has all the posterior draws for that quantity. For `type='fitted'` and `type='fitted.ind'` this attribute is a 3-dimensional array representing draws x observations generating predictions x levels of Y.

Author(s)

Frank Harrell

See Also

[predict.lrm\(\)](#)

Examples

```
## Not run:
f <- blrm(...)
predict(f, newdata, type='...', posterior.summary='median')

## End(Not run)
```

print.blrm *Print blrm() Results*

Description

Prints main results from `blrm()` along with indexes and predictive accuracy and their highest posterior density intervals computed from `blrmStats`.

Usage

```
## S3 method for class 'blrm'
print(
  x,
  dec = 4,
  coefs = TRUE,
  intercepts = x$non.slopes < 10,
  prob = 0.95,
  ns = 400,
  title = NULL,
  ...
)
```

Arguments

<code>x</code>	object created by <code>blrm()</code>
<code>dec</code>	number of digits to print to the right of the decimal
<code>coefs</code>	specify <code>FALSE</code> to suppress printing parameter estimates, and in integer <code>k</code> to print only the first <code>k</code>
<code>intercepts</code>	set to <code>FALSE</code> to suppress printing intercepts. Default is to print them unless there are more than 9.
<code>prob</code>	HPD interval probability for summary indexes
<code>ns</code>	number of random samples of the posterior draws for use in computing HPD intervals for accuracy indexes
<code>title</code>	title of output, constructed by default
<code>...</code>	passed to <code>prModFit</code>

Author(s)

Frank Harrell

Examples

```
## Not run:
f <- blrm(...)
options(lang='html') # default is lang='plain'; also can be latex
```

```
f          # print using defaults
print(f, posterior.summary='median') # instead of post. means

## End(Not run)
```

print.blrmStats *Print Details for blrmStats Predictive Accuracy Measures*

Description

Prints results of blrmStats with brief explanations

Usage

```
## S3 method for class 'blrmStats'
print(x, dec = 3, ...)
```

Arguments

x	an object produced by blrmStats
dec	number of digits to round indexes
...	ignored

Author(s)

Frank Harrell

Examples

```
## Not run:
f <- blrm(...)
s <- blrmStats(...)
s # print with defaults
print(s, dec=4)

## End(Not run)
```

```
print.predict.blrm      Print Predictions for blrm()
```

Description

Prints the summary portion of the results of `predict.blrm`

Usage

```
## S3 method for class 'predict.blrm'
print(x, digits = 3, ...)
```

Arguments

<code>x</code>	result from <code>predict.blrm</code>
<code>digits</code>	number of digits to round numeric results
<code>...</code>	ignored

Author(s)

Frank Harrell

```
print.rmsb              Basic Print for Bayesian Parameter Summary
```

Description

For a Bayesian regression fit prints the posterior mean, median, SE, highest posterior density interval, and symmetry coefficient from the posterior draws. For a given parameter, the symmetry measure is computed using the `distSym` function.

Usage

```
## S3 method for class 'rmsb'
print(x, prob = 0.95, dec = 4, intercepts = TRUE, pr = TRUE, ...)
```

Arguments

<code>x</code>	an object created by an rms Bayesian fitting function
<code>prob</code>	HPD interval coverage probability (default is 0.95)
<code>dec</code>	amount of rounding (digits to the right of the decimal)
<code>intercepts</code>	set to FALSE to not print intercepts
<code>pr</code>	set to FALSE to return an unrounded matrix and not print
<code>...</code>	ignored

Value

matrix (rounded if pr=TRUE)

Author(s)

Frank Harrell

Examples

```
## Not run:
f <- blrm(...)
print.rmsb(f)

## End(Not run)
```

Quantile.blrm

Function Generator for Quantiles of Y for `blrm()`

Description

Creates a function to turn a posterior summarized linear predictor lp (e.g. using posterior mean of intercepts and slopes) computed at the reference intercept into e.g. an estimate of a quantile of Y using the posterior mean of all the intercepts. `lptau` must be provided when call the created function if the model is a partial proportional odds model.

Usage

```
## S3 method for class 'blrm'
Quantile(object, codes = FALSE, posterior.summary = c("mean", "median"), ...)
```

Arguments

<code>object</code>	a <code>blrm()</code> fit
<code>codes</code>	if TRUE, use the integer codes $1, 2, \dots, k$ for the k -level response in computing the quantile
<code>posterior.summary</code>	defaults to posterior mean; may also specify "median". Must be consistent with the summary used when creating lp .
<code>...</code>	ignored

Value

an R function

Author(s)

Frank Harrell

Description

Runs a matrix through the QR decomposition and returns the transformed matrix and the forward and inverse transforming matrices R, Rinv. If columns of the input matrix X are centered the QR transformed matrix will be orthogonal. This is helpful in understanding the transformation and in scaling prior distributions on the transformed scale. not can be specified to keep selected columns as-is. cornerQr leaves the last column of X alone (possibly after centering). When not is specified, the square transforming matrices have appropriate identity submatrices inserted so that recreation of original X is automatic.

Usage

```
selectedQr(X, not = NULL, corner = FALSE, center = TRUE)
```

Arguments

X	a numeric matrix
not	an integer vector specifying which columns of X are to be kept with their original values
corner	set to FALSE to not treat the last column specially. You may not specify both not and corner.
center	set to FALSE to not center columns of X first

Value

list with elements X, R, Rinv, xbar where xbar is the vector of means (vector of zeros if center=FALSE)

Author(s)

Ben Goodrich and Frank Harrell

Examples

```
x <- 1 : 10
X <- cbind(x, x^2)
w <- selectedQr(X)
w
with(w, X %**% R) # = scale(X, center=TRUE, scale=FALSE)
Xqr <- w$X
plot(X[, 1], Xqr[, 1])
plot(X[, 1], Xqr[, 2])
cov(X)
cov(Xqr)
```

```
X <- cbind(x, x^3, x^4, x^2)
w <- selectedQr(X, not=2:3)
with(w, X %>% R)
```

stackMI

Bayesian Model Fitting and Stacking for Multiple Imputation

Description

Runs an rmsb package Bayesian fitting function such as `blrm` separately for each completed dataset given a multiple imputation result such as one produced by `Hmisc::aregImpute`. Stacks the posterior draws and diagnostics across all imputations, and computes parameter summaries on the stacked posterior draws.

Usage

```
stackMI(
  formula,
  fitter,
  xtrans,
  data = NULL,
  n.impute = xtrans$n.impute,
  dtrans = NULL,
  derived = NULL,
  subset = NULL,
  refresh = 0,
  progress = if (refresh > 0) "stan-progress.txt" else "",
  file = NULL,
  ...
)
```

Arguments

<code>formula</code>	a model formula
<code>fitter</code>	a Bayesian fitter
<code>xtrans</code>	an object created by <code>transcan</code> , <code>aregImpute</code> , or <code>mice</code>
<code>data</code>	data frame
<code>n.impute</code>	number of imputations to run, default is the number saved in <code>xtrans</code>
<code>dtrans</code>	see <code>Hmisc::fit.mult.impute</code>
<code>derived</code>	see <code>Hmisc::fit.mult.impute</code>
<code>subset</code>	an integer or logical vector specifying the subset of observations to fit
<code>refresh</code>	see <code>rstan::sampling</code> . The default is 0, indicating that no progress notes are output. If <code>refresh > 0</code> and <code>progress</code> is not <code>' '</code> , progress output will be appended to file <code>progress</code> . The default file name is <code>'stan-progress.txt'</code> .

progress	see refresh. Defaults to '' if refresh = 0. Note: If running interactively but not under RStudio, rstan will open a browser window for monitoring progress.
file	optional file name in which to store results in RDS format. If file is given and it already exists, and none of the arguments to stackMI have changed since that fit, the fit object from file is immediately returned. So if the model, data, and imputations have not changed nothing needs to be computed.
...	arguments passed to fitter

Value

an rmsb fit object with expanded posterior draws and diagnostics

Author(s)

Frank Harrell

stanDx *Print Stan Diagnostics*

Description

Retrieves the effect samples sizes and Rhats computed after a fitting function ran rstan, and prepares it for printing. If the fit was created by stackImpute, the diagnostics for all imputations are printed (separately).

Usage

```
stanDx(object)
```

Arguments

object	an object created by an rms package Bayesian fitting function such as blrm() or stackMI()
--------	---

Value

matrix suitable for printing

Author(s)

Frank Harrell

Examples

```
## Not run:
  f <- blrm(...)
  stanDx(f)

## End(Not run)
```

stanDxplot

*Diagnostic Trace Plots***Description**

For an rms Bayesian fit object, uses by default the stored posterior draws to check convergence properties of posterior sampling. If instead `rstan=TRUE`, calls the `rstan` `traceplot` function on the `rstan` object inside the `rmsb` object, to check properties of posterior sampling. If `rstan=TRUE` and the `rstan` object has been removed and `previous=TRUE`, attempts to find an already existing plot created by a previous run of the knitr chunk, assuming it was the `plotno` numbered plot of the chunk.

Usage

```
stanDxplot(
  x,
  which = NULL,
  rstan = FALSE,
  previous = TRUE,
  plotno = 1,
  rev = FALSE,
  stripsize = 8,
  ...
)
```

Arguments

<code>x</code>	an rms Bayesian fit object
<code>which</code>	names of parameters to plot, defaulting to all non-intercepts. When <code>rstan=FALSE</code> these are the friendly rms names, otherwise they are the <code>rstan</code> parameter names. If the model fit was run through <code>stackMI</code> for multiple imputation, the number of traces is multiplied by the number of imputations. Set to 'ALL' to plot all parameters.
<code>rstan</code>	set to <code>TRUE</code> to use <code>rstan::traceplot()</code> on a (presumed) stored <code>rstan</code> object in <code>x</code> , otherwise only real iterations are plotted and parameter values are shown as points instead of lines, with chains separated
<code>previous</code>	see details
<code>plotno</code>	see details
<code>rev</code>	set to <code>TRUE</code> to reverse direction for faceting chains
<code>stripsize</code>	specifies size of chain facet label text, default is 8
<code>...</code>	passed to <code>rstan::traceplot()</code>

Value

ggplot2 object if `rstan` object was in `x`

Author(s)

Frank Harrell

stanGet

Get Stan Output

Description

Extracts the object created by `rstan::sampling()` so that standard Stan diagnostics can be run from it

Usage

```
stanGet(object)
```

Arguments

`object` an object created by an rms package Bayesian fitting function

Value

the object created by `rstan::sampling()`

Author(s)

Frank Harrell

Examples

```
## Not run:  
f <- blrm(...)  
s <- stanGet(f)  
  
## End(Not run)
```

tauFetch

Fetch Partial Proportional Odds Parameters

Description

Fetches matrix of posterior draws for partial proportional odds parameters (taus) for a given intercept. Can also form a matrix containing both regular parameters and taus, or for just non-taus. For the constrained partial proportional odds model the function returns the appropriate cppo function value multiplied by tau (tau being a vector in this case and not a matrix).

Usage

```
tauFetch(fit, intercept, what = c("tau", "nontau", "both"))
```

Arguments

fit	an object created by <code>blrm()</code>
intercept	integer specifying which intercept to fetch
what	specifies the result to return

Value

matrix with number of rows equal to the number of original draws

Author(s)

Frank Harrell

vcov.rmsb

Variance-Covariance Matrix

Description

Computes the variance-covariance matrix from the posterior draws by compute the sample covariance matrix of the draws

Usage

```
## S3 method for class 'rmsb'
vcov(object, regcoef.only = TRUE, intercepts = "all", ...)
```


Arguments

object	an object produced by an rms package Bayesian fitting function
regcoef.only	set to FALSE to also include non-regression coefficients such as shape/scale parameters
intercepts	set to 'all' to include all intercepts (the default), 'none' to exclude them all, or a vector of integers to get selected intercepts
...	ignored

Value

matrix

Author(s)

Frank Harrell

See Also[vcov.rms](#)**Examples**

```
## Not run:
  f <- blrm(...)
  v <- vcov(f)

## End(Not run)
```

[.Ocens

*Ocens***Description**

Subset Method for Ocens Objects

Usage

```
## S3 method for class 'Ocens'
x[rows = 1:d[1], cols = 1:d[2], ...]
```

Arguments

x	an Ocens object
rows	logical or integer vector
cols	logical or integer vector
...	ignored

Details

Subsets an Ocens object, preserving its special attributes. Attributes are not updated. In the future such updating should be implemented.

Value

new Ocens object

Author(s)

Frank Harrell

Index

[.Ocens, 33

aregImpute, 28

as.data.frame.Ocens, 3

blrm, 4

blrm(), 4, 5, 7–10, 12–16, 20, 21, 23, 25, 26, 29, 32

blrmStats, 9

blrmStats(), 9

cluster, 11

cmdstanr::sample(), 7

coda::HPDinterval(), 15

coef.rmsb, 11

coef.rmsb(), 9

compareBmods, 12

distSym, 13

ExProb(), 22

ExProb.blrm, 13

getParamCoef, 14

Hmisc::rcorr.cens(), 10

HPDint, 15

loo::loo_model_weights(), 12

MASS::kde2d(), 17

Mean(), 22

Mean.blrm, 15

mice, 28

Ocens, 16

pdensityContour, 17

plot.PostF, 18

plot.rmsb, 19

PostF, 20

predict.blrm, 21

predict.lrm(), 21, 22

print.blrm, 23

print.blrm(), 9

print.blrmStats, 24

print.predict.blrm, 25

print.rmsb, 25

print.rmsb(), 9

Quantile(), 22

Quantile.blrm, 26

rms::contrast.rms(), 6

rms::datadist(), 6

rms::gendata(), 6

rms::Xcontrast(), 6

rmsb (rmsb-package), 3

rmsb-package, 3

rstan::optimizing(), 7, 8

rstan::sampling, 28

rstan::sampling(), 6–8, 31

rstan::traceplot(), 30

selectedQr, 27

stackMI, 28

stackMI(), 29

stanDx, 29

stanDx(), 9

stanDxplot, 30

stanGet, 31

stanGet(), 9

tauFetch, 32

vcov.rms, 33

vcov.rmsb, 32

vcov.rmsb(), 9