

# Package ‘countSTAR’

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**Type** Package

**Title** Flexible Modeling of Count Data

**Version** 1.2.0

**Description** For Bayesian and classical inference and prediction with count-valued data, Simultaneous Transformation and Rounding (STAR) Models provide a flexible, interpretable, and easy-to-use approach. STAR models the observed count data using a rounded continuous data model and incorporates a transformation for greater flexibility. Implicitly, STAR formalizes the commonly-applied yet incoherent procedure of (i) transforming count-valued data and subsequently (ii) modeling the transformed data using Gaussian models. STAR is well-defined for count-valued data, which is reflected in predictive accuracy, and is designed to account for zero-inflation, bounded or censored data, and over- or underdispersion. Importantly, STAR is easy to combine with existing MCMC or point estimation methods for continuous data, which allows seamless adaptation of continuous data models (such as linear regressions, additive models, BART, random forests, and gradient boosting machines) for count-valued data. The package also includes several methods for modeling count time series data, namely via warped Dynamic Linear Models. For more details and background on these methodologies, see the works of Kowal and Canale (2020) <[doi:10.1214/20-EJS1707](https://doi.org/10.1214/20-EJS1707)>, Kowal and Wu (2022) <[doi:10.1111/biom.13617](https://doi.org/10.1111/biom.13617)>, King and Kowal (2023) <[doi:10.1214/23-BA1394](https://doi.org/10.1214/23-BA1394)>, and Kowal and Wu (2023) <[doi:10.48550/arXiv.2110.12316](https://doi.org/10.48550/arXiv.2110.12316)>.

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<https://github.com/bking124/countSTAR>

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## Contents

a_j	3
bam_star	4
bart_star	6
blm_star	10
confint.lmstar	13
credBands	14
ergMean	14
gbm_star	15
genEM_star	18
genMCMC_star	20
getEffSize	23
g_bc	24
g_cdf	25
g_inv	26
g_inv_approx	26
g_inv_bc	27
HPDregion	28
init_lm_gprior	28
lm_star	29
plot_coef	31
plot_fitted	32
plot_pmf	33
predict.lmstar	33
pvals	35
randomForest_star	36
rdir	38
roaches	39
round_floor	40
sample_lm_gprior	40
simBaS	41

<code>a_j</code>	3
<code>simulate_nb_friedman</code> . . . . .	42
<code>simulate_nb_lm</code> . . . . .	43
<code>spline_star</code> . . . . .	45
<code>warpDLM</code> . . . . .	47
<b>Index</b>	<b>49</b>

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<code>a_j</code>	<i>Inverse rounding function</i>
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### Description

Define the intervals associated with  $y = j$  based on the flooring function. The function returns  $-\text{Inf}$  for  $j = 0$  (or smaller) and  $\text{Inf}$  for any  $j \geq y_{\text{max}} + 1$ , where  $y_{\text{max}}$  is a known upper bound on the data  $y$  (if specified).

### Usage

```
a_j(j, y_max = Inf)
```

### Arguments

<code>j</code>	the integer-valued input(s)
<code>y_max</code>	a fixed and known upper bound for all observations; default is $\text{Inf}$

### Value

The (lower) interval endpoint(s) associated with  $j$ .

### Examples

```
# Standard cases:
a_j(1)
a_j(20)

# Boundary cases:
a_j(0)
a_j(20, y_max = 15)
```

bam\_star

*Fit Bayesian Additive STAR Model with MCMC***Description**

Run the MCMC algorithm for a STAR Bayesian additive model. The transformation can be known (e.g., log or sqrt) or unknown (Box-Cox or estimated nonparametrically) for greater flexibility.

**Usage**

```
bam_star(
  y,
  X_lin,
  X_nonlin,
  splinetype = "orthogonal",
  transformation = "np",
  y_max = Inf,
  nsave = 1000,
  nburn = 1000,
  nskip = 0,
  save_y_hat = FALSE,
  verbose = TRUE
)
```

**Arguments**

y	n x 1 vector of observed counts
X_lin	n x pL matrix of predictors to be modelled as linear
X_nonlin	n x pNL matrix of predictors to be modelled as nonlinear
splinetype	Type of spline to use for modelling the nonlinear predictors; must be either "orthogonal" (orthogonalized splines—the default) or "thinplate" (low-rank thin plate splines)
transformation	transformation to use for the latent data; must be one of <ul style="list-style-type: none"> <li>• "identity" (identity transformation)</li> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>• "box-cox" (box-cox transformation with learned parameter)</li> <li>• "ispline" (transformation is modeled as unknown, monotone function using I-splines)</li> </ul>
y_max	a fixed and known upper bound for all observations; default is Inf

nsave	number of MCMC iterations to save
nburn	number of MCMC iterations to discard
nskip	number of MCMC iterations to skip between saving iterations, i.e., save every (nskip + 1)th draw
save_y_hat	logical; if TRUE, compute and save the posterior draws of the expected counts, $E(y)$ , which may be slow to compute
verbose	logical; if TRUE, print time remaining

## Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous *latent* data and (2) connecting the latent data to the observed data via a *transformation and rounding* operation.

Posterior and predictive inference is obtained via a Gibbs sampler that combines (i) a latent data augmentation step (like in probit regression) and (ii) an existing sampler for a continuous data model.

There are several options for the transformation. First, the transformation can belong to the *Box-Cox* family, which includes the known transformations 'identity', 'log', and 'sqrt', as well as a version in which the Box-Cox parameter is inferred within the MCMC sampler ('box-cox'). Second, the transformation can be estimated (before model fitting) using the empirical distribution of the data  $y$ . Options in this case include the empirical cumulative distribution function (CDF), which is fully nonparametric ('np'), or the parametric alternatives based on Poisson ('pois') or Negative-Binomial ('neg-bin') distributions. For the parametric distributions, the parameters of the distribution are estimated using moments (means and variances) of  $y$ . Third, the transformation can be modeled as an unknown, monotone function using I-splines ('ispline'). The Robust Adaptive Metropolis (RAM) sampler is used for drawing the parameter of the transformation function.

## Value

a list with at least the following elements:

- `coefficients`: the posterior mean of the coefficients
- `fitted.values`: the posterior mean of the conditional expectation of the counts  $y$
- `post.coefficients`: posterior draws of the coefficients
- `post.fitted.values`: posterior draws of the conditional mean of the counts  $y$
- `post.pred`: draws from the posterior predictive distribution of  $y$
- `post.lambda`: draws from the posterior distribution of  $\lambda$
- `post.sigma`: draws from the posterior distribution of  $\sigma$
- `post.log.like.point`: draws of the log-likelihood for each of the  $n$  observations
- WAIC: Widely-Applicable/Watanabe-Akaike Information Criterion
- `p_waic`: Effective number of parameters based on WAIC

In the case of `transformation="ispline"`, the list also contains

- `post.g`: draws from the posterior distribution of the transformation  $g$
- `post.sigma.gamma`: draws from the posterior distribution of `sigma.gamma`, the prior standard deviation of the transformation  $g()$  coefficients

**Examples**

```

# Simulate data with count-valued response y:
sim_dat = simulate_nb_friedman(n = 100, p = 5, seed=32)
y = sim_dat$y; X = sim_dat$X

# Linear and nonlinear components:
X_lin = as.matrix(X[,-(1:3)])
X_nonlin = as.matrix(X[, (1:3)])

# STAR: nonparametric transformation
library(spikeSlabGAM)
fit = bam_star(y = y, X_lin = X_lin, X_nonlin = X_nonlin)

# What is included:
names(fit)

# Posterior mean of each coefficient:
coef(fit)

# WAIC:
fit$WAIC

# MCMC diagnostics:
plot(as.ts(fit$post.coefficients[,1:3]))

# Posterior predictive check:
hist(apply(fit$post.pred, 1,
          function(x) mean(x==0)), main = 'Proportion of Zeros', xlab='');
abline(v = mean(y==0), lwd=4, col = 'blue')

```

---

bart\_star

*MCMC Algorithm for BART-STAR*


---

**Description**

Run the MCMC algorithm for a BART model for count-valued responses using STAR. The transformation can be known (e.g., log or sqrt) or unknown (Box-Cox or estimated nonparametrically) for greater flexibility.

**Usage**

```

bart_star(
  y,
  X,
  X_test = NULL,
  y_test = NULL,
  transformation = "np",

```

```

y_max = Inf,
n.trees = 200,
sigest = NULL,
sigdf = 3,
sigquant = 0.9,
k = 2,
power = 2,
base = 0.95,
nsave = 1000,
nburn = 1000,
nskip = 0,
save_y_hat = FALSE,
verbose = TRUE
)

```

### Arguments

y	n x 1 vector of observed counts
X	n x p matrix of predictors
X_test	n_test x p matrix of predictors for test data
y_test	n_test x 1 vector of the test data responses (used for computing log-predictive scores)
transformation	transformation to use for the latent process; must be one of <ul style="list-style-type: none"> <li>• "identity" (identity transformation)</li> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>• "box-cox" (box-cox transformation with learned parameter)</li> <li>• "ispline" (transformation is modeled as unknown, monotone function using I-splines)</li> </ul>
y_max	a fixed and known upper bound for all observations; default is Inf
n.trees	number of trees to use in BART; default is 200
sigest	positive numeric estimate of the residual standard deviation (see ?bart)
sigdf	degrees of freedom for error variance prior (see ?bart)
sigquant	quantile of the error variance prior that the rough estimate (sigest) is placed at. The closer the quantile is to 1, the more aggressive the fit will be (see ?bart)
k	the number of prior standard deviations $E(Y x) = f(x)$ is away from +/- 0.5. The response is internally scaled to range from -0.5 to 0.5. The bigger k is, the more conservative the fitting will be (see ?bart)
power	power parameter for tree prior (see ?bart)
base	base parameter for tree prior (see ?bart)

nsave	number of MCMC iterations to save
nburn	number of MCMC iterations to discard
nskip	number of MCMC iterations to skip between saving iterations, i.e., save every (nskip + 1)th draw
save_y_hat	logical; if TRUE, compute and save the posterior draws of the expected counts, $E(y)$ , which may be slow to compute
verbose	logical; if TRUE, print time remaining

### Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous *latent* data and (2) connecting the latent data to the observed data via a *transformation and rounding* operation. Here, the model in (1) is a Bayesian additive regression tree (BART) model.

Posterior and predictive inference is obtained via a Gibbs sampler that combines (i) a latent data augmentation step (like in probit regression) and (ii) an existing sampler for a continuous data model.

There are several options for the transformation. First, the transformation can belong to the *Box-Cox* family, which includes the known transformations 'identity', 'log', and 'sqrt', as well as a version in which the Box-Cox parameter is inferred within the MCMC sampler ('box-cox'). Second, the transformation can be estimated (before model fitting) using the empirical distribution of the data  $y$ . Options in this case include the empirical cumulative distribution function (CDF), which is fully nonparametric ('np'), or the parametric alternatives based on Poisson ('pois') or Negative-Binomial ('neg-bin') distributions. For the parametric distributions, the parameters of the distribution are estimated using moments (means and variances) of  $y$ . Third, the transformation can be modeled as an unknown, monotone function using I-splines ('ispline'). The Robust Adaptive Metropolis (RAM) sampler is used for drawing the parameter of the transformation function.

### Value

a list with the following elements:

- `post.pred`: draws from the posterior predictive distribution of  $y$
- `post.sigma`: draws from the posterior distribution of  $\sigma$
- `post.log.like.point`: draws of the log-likelihood for each of the  $n$  observations
- WAIC: Widely-Applicable/Watanabe-Akaike Information Criterion
- `p_waic`: Effective number of parameters based on WAIC
- `post.pred.test`: draws from the posterior predictive distribution at the test points  $X_{\text{test}}$  (NULL if  $X_{\text{test}}$  is not given)
- `post.fitted.values.test`: posterior draws of the conditional mean at the test points  $X_{\text{test}}$  (NULL if  $X_{\text{test}}$  is not given)
- `post.mu.test`: draws of the conditional mean of  $z_{\text{star}}$  at the test points  $X_{\text{test}}$  (NULL if  $X_{\text{test}}$  is not given)
- `post.log.pred.test`: draws of the log-predictive distribution for each of the  $n_{\text{test}}$  test cases (NULL if  $X_{\text{test}}$  is not given)

- `fitted.values`: the posterior mean of the conditional expectation of the counts  $y$  (NULL if `save_y_hat=FALSE`)
- `post.fitted.values`: posterior draws of the conditional mean of the counts  $y$  (NULL if `save_y_hat=FALSE`)

In the case of `transformation="ispline"`, the list also contains

- `post.g`: draws from the posterior distribution of the transformation  $g$
- `post.sigma.gamma`: draws from the posterior distribution of `sigma.gamma`, the prior standard deviation of the transformation  $g()$  coefficients

If `transformation="box-cox"`, then the list also contains

- `post.lambda`: draws from the posterior distribution of  $\lambda$

## Examples

```
# Simulate data with count-valued response y:
sim_dat = simulate_nb_friedman(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X

# BART-STAR with log-transformation:
fit_log = bart_star(y = y, X = X, transformation = 'log',
                   save_y_hat = TRUE, nburn=1000, nskip=0)

# Fitted values
plot_fitted(y = sim_dat$Ey,
            post_y = fit_log$post.fitted.values,
            main = 'Fitted Values: BART-STAR-log')

# WAIC for BART-STAR-log:
fit_log$WAIC

# MCMC diagnostics:
plot(as.ts(fit_log$post.fitted.values[,1:10]))

# Posterior predictive check:
hist(apply(fit_log$post.pred, 1,
          function(x) mean(x==0)), main = 'Proportion of Zeros', xlab='');
abline(v = mean(y==0), lwd=4, col = 'blue')

# BART-STAR with nonparametric transformation:
fit = bart_star(y = y, X = X,
               transformation = 'np', save_y_hat = TRUE)

# Fitted values
plot_fitted(y = sim_dat$Ey,
            post_y = fit$post.fitted.values,
            main = 'Fitted Values: BART-STAR-np')

# WAIC for BART-STAR-np:
fit$WAIC
```

```

# MCMC diagnostics:
plot(as.ts(fit$post.fitted.values[,1:10]))

# Posterior predictive check:
hist(apply(fit$post.pred, 1,
          function(x) mean(x==0)), main = 'Proportion of Zeros', xlab='');
abline(v = mean(y==0), lwd=4, col = 'blue')

```

---

blm\_star

*STAR Bayesian Linear Regression*


---

### Description

Posterior and predictive inference for STAR linear model

### Usage

```

blm_star(
  y,
  X,
  X_test = X,
  transformation = "bnp",
  y_max = Inf,
  prior = "gprior",
  use_MCMC = TRUE,
  nsave = 1000,
  nburn = 1000,
  nskip = 0,
  psi = length(y),
  alpha = 1,
  F0 = NULL,
  compute_marg = FALSE,
  verbose = FALSE
)

```

### Arguments

y	n x 1 vector of observed counts
X	n x p matrix of predictors
X_test	n_test x p matrix of predictors for test data; default is the observed covariates X
transformation	transformation to use for the latent process; must be one of <ul style="list-style-type: none"> <li>"identity" (identity transformation)</li> <li>"log" (log transformation)</li> </ul>

- "sqrt" (square root transformation)
- "np" (nonparametric transformation estimated from empirical CDF)
- "pois" (transformation for moment-matched marginal Poisson CDF)
- "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)
- "box-cox" (box-cox transformation with learned parameter)
- "ispline" (transformation is modeled as unknown, monotone function using I-splines)
- "bnp" (Bayesian nonparametric transformation)

y_max	a fixed and known upper bound for all observations; default is Inf
prior	prior to use for the latent linear regression; currently implemented options are "gprior", "horseshoe", and "ridge"
use_MCMC	logical; whether to run Gibbs sampler or Monte Carlo (default is TRUE)
nsave	number of MC(MC) iterations to save
nburn	number of MCMC iterations to discard
nskip	number of MCMC iterations to skip between saving iterations, i.e., save every (nskip + 1)th draw
psi	prior variance (g-prior)
alpha	prior precision for the Dirichlet Process prior ('bnp' transformation only); default is one
F0	function to evaluate the base measure CDF supported on $\{0, \dots, y_{\max}\}$ ('bnp' transformation only)
compute_marg	logical; if TRUE, compute and return the marginal likelihood (only available when using exact sampler, i.e. use_MCMC=FALSE)
verbose	logical; if TRUE, print time remaining

## Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous *latent* data and (2) connecting the latent data to the observed data via a *transformation and rounding* operation. Here, the continuous latent data model is a linear regression.

There are several options for the transformation. First, the transformation can belong to the *Box-Cox* family, which includes the known transformations 'identity', 'log', and 'sqrt', as well as a version in which the Box-Cox parameter is inferred within the MCMC sampler ('box-cox').

Second, the transformation can be estimated (before model fitting) using the the data  $y$ . Options in this case include the empirical cumulative distribution function (ECDF), which is fully nonparametric ('np'), or the parametric alternatives based on Poisson ('pois') or Negative-Binomial ('neg-bin') distributions. For the parametric distributions, the parameters of the distribution are estimated using moments (means and variances) of  $y$ .

Lastly, the transformation can be modeled nonparametrically using (monotone) splines ('ispline') or Bayesian nonparametrics via Dirichlet processes ('bnp'). The 'bnp' option is the default because it is highly flexible, accounts for uncertainty when the transformation is unknown, and is computationally efficient.

The Monte Carlo sampler (`use_MCMC=FALSE`) produces direct, joint draws from the posterior predictive distribution under a g-prior. When `n` is moderate to large, or to use other priors, MCMC sampling (`use_MCMC=TRUE`) is much faster and more convenient.

## Value

a list with at least the following elements:

- `coefficients`: the posterior mean of the regression coefficients
- `post.beta`: posterior draws of the regression coefficients
- `post.pred`: draws from the posterior predictive distribution of `y`
- `post.log.like.point`: draws of the log-likelihood for each of the `n` observations
- `WAIC`: Widely-Applicable/Watanabe-Akaike Information Criterion
- `p_waic`: Effective number of parameters based on WAIC

Other elements may be present depending on the choice of prior, transformation, and sampling approach.

## Examples

```
# Simulate data with count-valued response y:
sim_dat = simulate_nb_lm(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X

# Fit the Bayesian STAR linear model:
fit = blm_star(y = y, X = X)

# What is included:
names(fit)

# Posterior mean of each coefficient:
coef(fit)

# WAIC:
fit$WAIC

# MCMC diagnostics:
plot(as.ts(fit$post.beta))

# Posterior predictive check:
hist(apply(fit$post.pred, 1,
          function(x) mean(x==0)), main = 'Proportion of Zeros', xlab='');
abline(v = mean(y==0), lwd=4, col = 'blue')
```

---

confint.lmstar	<i>Compute asymptotic confidence intervals for STAR linear regression</i>
----------------	---

---

## Description

For a linear regression model within the STAR framework, compute (asymptotic) confidence intervals for a regression coefficient of interest. Confidence intervals are computed by inverting the likelihood ratio test and profiling the log-likelihood.

## Usage

```
## S3 method for class 'lmstar'
confint(object, parm, level = 0.95, ...)
```

## Arguments

object	Object of class "lmstar" as output by <code>lm_star</code>
parm	a specification of which parameters are to be given confidence intervals, either a vector of numbers or a vector of names. If missing, all parameters are considered.
level	confidence level; default is 0.95
...	Ignored

## Value

A matrix (or vector) with columns giving lower and upper confidence limits for each parameter. These will be labelled as  $(1-\text{level})/2$  and  $1 - (1-\text{level})/2$  in

## Examples

```
#Simulate data with count-valued response y:
sim_dat = simulate_nb_lm(n = 100, p = 2)
y = sim_dat$y; X = sim_dat$X[,-1] # remove intercept

# Select a transformation:
transformation = 'np'

#Estimate model
fit = lm_star(y~X, transformation = transformation)

#Confidence interval for all parameters
confint(fit)
```

---

credBands	<i>Compute Simultaneous Credible Bands</i>
-----------	--

---

**Description**

Compute (1-alpha)% credible BANDS for a function based on MCMC samples using Crainiceanu et al. (2007)

**Usage**

```
credBands(sampFuns, alpha = 0.05)
```

**Arguments**

sampFuns	Nsims x m matrix of Nsims MCMC samples and m points along the curve
alpha	confidence level

**Value**

m x 2 matrix of credible bands; the first column is the lower band, the second is the upper band

**Note**

The input needs not be curves: the simultaneous credible "bands" may be computed for vectors. The resulting credible intervals will provide joint coverage at the (1-alpha) level across all components of the vector.

---

ergMean	<i>Compute the ergodic (running) mean.</i>
---------	--

---

**Description**

Compute the ergodic (running) mean.

**Usage**

```
ergMean(x)
```

**Arguments**

x	vector for which to compute the running mean
---	--

**Value**

A vector y with each element defined by  $y[i] = \text{mean}(x[1:i])$

**Examples**

```
# Compare:
ergMean(1:10)
mean(1:10)

# Running mean for iid N(5, 1) samples:
x = rnorm(n = 10^4, mean = 5, sd = 1)
plot(ergMean(x))
abline(h=5)
```

gbm\_star

*Fitting STAR Gradient Boosting Machines via EM algorithm***Description**

Compute the MLEs and log-likelihood for the Gradient Boosting Machines (GBM) STAR model. The STAR model requires a *transformation* and an *estimation function* for the conditional mean given observed data. The transformation can be known (e.g., log or sqrt) or unknown (Box-Cox or estimated nonparametrically) for greater flexibility. The estimator in this case is a GBM. Standard function calls including `fitted` and `residuals` apply.

**Usage**

```
gbm_star(
  y,
  X,
  X.test = NULL,
  transformation = "np",
  y_max = Inf,
  sd_init = 10,
  tol = 10^-3,
  max_iters = 1000,
  n.trees = 100,
  interaction.depth = 1,
  shrinkage = 0.1,
  bag.fraction = 1
)
```

**Arguments**

<code>y</code>	<code>n x 1</code> vector of observed counts
<code>X</code>	<code>n x p</code> matrix of predictors
<code>X.test</code>	<code>m x p</code> matrix of out-of-sample predictors
<code>transformation</code>	transformation to use for the latent data; must be one of <ul style="list-style-type: none"> <li>"identity" (identity transformation)</li> </ul>

	<ul style="list-style-type: none"> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>• "box-cox" (box-cox transformation with learned parameter)</li> </ul>
y_max	a fixed and known upper bound for all observations; default is Inf
sd_init	add random noise for EM algorithm initialization scaled by sd_init times the Gaussian MLE standard deviation; default is 10
tol	tolerance for stopping the EM algorithm
max_iters	maximum number of EM iterations before stopping; default is 1000
n.trees	Integer specifying the total number of trees to fit. This is equivalent to the number of iterations and the number of basis functions in the additive expansion. Default is 100.
interaction.depth	Integer specifying the maximum depth of each tree (i.e., the highest level of variable interactions allowed). A value of 1 implies an additive model, a value of 2 implies a model with up to 2-way interactions, etc. Default is 1.
shrinkage	a shrinkage parameter applied to each tree in the expansion. Also known as the learning rate or step-size reduction; 0.001 to 0.1 usually work, but a smaller learning rate typically requires more trees. Default is 0.1.
bag.fraction	the fraction of the training set observations randomly selected to propose the next tree in the expansion. This introduces randomness into the model fit. If bag.fraction < 1 then running the same model twice will result in similar but different fits. Default is 1 (for a deterministic prediction).

## Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous \*latent\* data and (2) connecting the latent data to the observed data via a \*transformation and rounding\* operation. The Gaussian model in this case is a GBM.

## Value

a list with the following elements:

- fitted.values: the fitted values at the MLEs (training)
- fitted.values.test: the fitted values at the MLEs (testing)
- g.hat a function containing the (known or estimated) transformation
- sigma.hat the MLE of the standard deviation
- mu.hat the MLE of the conditional mean (on the transformed scale)
- z.hat the estimated latent data (on the transformed scale) at the MLEs
- residuals the Dunn-Smyth residuals (randomized)

- residuals\_rep the Dunn-Smyth residuals (randomized) for 10 replicates
- logLik the log-likelihood at the MLEs
- logLik0 the log-likelihood at the MLEs for the *\*unrounded\** initialization
- lambda the Box-Cox nonlinear parameter
- gbmObj: the object returned by gbm() at the MLEs
- and other parameters that (1) track the parameters across EM iterations and (2) record the model specifications

### Note

Infinite latent data values may occur when the transformed Gaussian model is highly inadequate. In that case, the function returns the *\*indices\** of the data points with infinite latent values, which are significant outliers under the model. Deletion of these indices and re-running the model is one option, but care must be taken to ensure that (i) it is appropriate to treat these observations as outliers and (ii) the model is adequate for the remaining data points.

### References

Kowal, D. R., & Wu, B. (2021). Semiparametric count data regression for self-reported mental health. *Biometrics*. doi:10.1111/biom.13617

### Examples

```
# Simulate data with count-valued response y:
sim_dat = simulate_nb_friedman(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X

# EM algorithm for STAR (using the log-link)
library(gbm)
fit_em = gbm_star(y = y, X = X,
                 transformation = 'log')

# Evaluate convergence:
plot(fit_em$logLik_all, type='l', main = 'GBM-STAR-log', xlab = 'Iteration', ylab = 'log-lik')

# Fitted values:
y_hat = fitted(fit_em)
plot(y_hat, y);

# Residuals:
plot(residuals(fit_em))
qqnorm(residuals(fit_em)); qqline(residuals(fit_em))

# Log-likelihood at MLEs:
fit_em$logLik
```

genEM\_star

*Generalized EM estimation for STAR***Description**

Compute MLEs and log-likelihood for a generalized STAR model. The STAR model requires a *transformation* and an *estimation function* for the conditional mean given observed data. The transformation can be known (e.g., log or sqrt) or unknown (Box-Cox or estimated nonparametrically) for greater flexibility. The estimator can be any least squares estimator, including nonlinear models. Standard function calls including `coefficients()`, `fitted()`, and `residuals()` apply.

**Usage**

```
genEM_star(
  y,
  estimator,
  transformation = "np",
  y_max = Inf,
  sd_init = 10,
  tol = 10^-10,
  max_iters = 1000
)
```

**Arguments**

<code>y</code>	<code>n x 1</code> vector of observed counts
<code>estimator</code>	a function that inputs data <code>y</code> and outputs a list with two elements: <ol style="list-style-type: none"> <li>1. The fitted values <code>fitted.values</code></li> <li>2. The parameter estimates <code>coefficients</code></li> </ol>
<code>transformation</code>	transformation to use for the latent data; must be one of <ul style="list-style-type: none"> <li>• "identity" (identity transformation)</li> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>• "box-cox" (box-cox transformation with learned parameter)</li> </ul>
<code>y_max</code>	a fixed and known upper bound for all observations; default is <code>Inf</code>
<code>sd_init</code>	add random noise for EM algorithm initialization scaled by <code>sd_init</code> times the Gaussian MLE standard deviation; default is <code>10</code>
<code>tol</code>	tolerance for stopping the EM algorithm; default is <code>10^-10</code> ;
<code>max_iters</code>	maximum number of EM iterations before stopping; default is <code>1000</code>

## Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous \*latent\* data and (2) connecting the latent data to the observed data via a \*transformation and rounding\* operation.

The expectation-maximization (EM) algorithm is used to produce maximum likelihood estimators (MLEs) for the parameters defined in the estimator function, such as linear regression coefficients, which define the Gaussian model for the continuous latent data. Fitted values (point predictions), residuals, and log-likelihood values are also available. Inference for the estimators proceeds via classical maximum likelihood. Initialization of the EM algorithm can be randomized to monitor convergence. However, the log-likelihood is concave for all transformations (except 'box-cox'), so global convergence is guaranteed.

There are several options for the transformation. First, the transformation can belong to the \*Box-Cox\* family, which includes the known transformations 'identity', 'log', and 'sqrt', as well as a version in which the Box-Cox parameter is estimated within the EM algorithm ('box-cox'). Second, the transformation can be estimated (before model fitting) using the empirical distribution of the data  $y$ . Options in this case include the empirical cumulative distribution function (CDF), which is fully nonparametric ('np'), or the parametric alternatives based on Poisson ('pois') or Negative-Binomial ('neg-bin') distributions. For the parametric distributions, the parameters of the distribution are estimated using moments (means and variances) of  $y$ .

## Value

a list with the following elements:

- `coefficients` the MLEs of the coefficients
- `fitted.values` the fitted values at the MLEs
- `g.hat` a function containing the (known or estimated) transformation
- `sigma.hat` the MLE of the standard deviation
- `mu.hat` the MLE of the conditional mean (on the transformed scale)
- `z.hat` the estimated latent data (on the transformed scale) at the MLEs
- `residuals` the Dunn-Smyth residuals (randomized)
- `residuals_rep` the Dunn-Smyth residuals (randomized) for 10 replicates
- `logLik` the log-likelihood at the MLEs
- `logLik0` the log-likelihood at the MLEs for the \*unrounded\* initialization
- `lambda` the Box-Cox nonlinear parameter
- and other parameters that (1) track the parameters across EM iterations and (2) record the model specifications

## Note

Infinite latent data values may occur when the transformed Gaussian model is highly inadequate. In that case, the function returns the \*indices\* of the data points with infinite latent values, which are significant outliers under the model. Deletion of these indices and re-running the model is one option, but care must be taken to ensure that (i) it is appropriate to treat these observations as outliers and (ii) the model is adequate for the remaining data points.

## References

Kowal, D. R., & Wu, B. (2021). Semiparametric count data regression for self-reported mental health. *Biometrics*. doi:10.1111/biom.13617

## Examples

```
# Simulate data with count-valued response y:
sim_dat = simulate_nb_friedman(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X

# Select a transformation:
transformation = 'np'

# Example using GAM as underlying estimator (for illustration purposes only)
if(require("mgcv")){
  fit_em = genEM_star(y = y,
                    estimator = function(y) gam(y ~ s(X1)+s(X2),
                    data=data.frame(y,X)),
                    transformation = transformation)
}

# Fitted coefficients:
coef(fit_em)

# Fitted values:
y_hat = fitted(fit_em)
plot(y_hat, y);

# Log-likelihood at MLEs:
fit_em$logLik
```

## Description

Run the MCMC algorithm for STAR given

1. a function to initialize model parameters; and
2. a function to sample (i.e., update) model parameters.

The transformation can be known (e.g., log or sqrt) or unknown (Box-Cox or estimated nonparametrically) for greater flexibility.

**Usage**

```
genMCMC_star(
  y,
  sample_params,
  init_params,
  transformation = "np",
  y_max = Inf,
  nsave = 1000,
  nburn = 1000,
  nskip = 0,
  save_y_hat = FALSE,
  verbose = TRUE
)
```

**Arguments**

<code>y</code>	<code>n x 1</code> vector of observed counts
<code>sample_params</code>	a function that inputs data <code>y</code> and a named list <code>params</code> containing <ol style="list-style-type: none"> <li><code>mu</code>: the <code>n x 1</code> vector of conditional means (fitted values)</li> <li><code>sigma</code>: the conditional standard deviation</li> <li><code>coefficients</code>: a named list of parameters that determine <code>mu</code></li> </ol> and outputs an updated list <code>params</code> of samples from the full conditional posterior distribution of coefficients and <code>sigma</code> (and updates <code>mu</code> )
<code>init_params</code>	an initializing function that inputs data <code>y</code> and initializes the named list <code>params</code> of <code>mu</code> , <code>sigma</code> , and coefficients
<code>transformation</code>	transformation to use for the latent data; must be one of <ul style="list-style-type: none"> <li>"identity" (identity transformation)</li> <li>"log" (log transformation)</li> <li>"sqrt" (square root transformation)</li> <li>"np" (nonparametric transformation estimated from empirical CDF)</li> <li>"pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>"neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>"box-cox" (box-cox transformation with learned parameter)</li> </ul>
<code>y_max</code>	a fixed and known upper bound for all observations; default is <code>Inf</code>
<code>nsave</code>	number of MCMC iterations to save
<code>nburn</code>	number of MCMC iterations to discard
<code>nskip</code>	number of MCMC iterations to skip between saving iterations, i.e., save every <code>(nskip + 1)</code> th draw
<code>save_y_hat</code>	logical; if <code>TRUE</code> , compute and save the posterior draws of the expected counts, $E(y)$ , which may be slow to compute
<code>verbose</code>	logical; if <code>TRUE</code> , print time remaining

## Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous \*latent\* data and (2) connecting the latent data to the observed data via a \*transformation and rounding\* operation.

Posterior and predictive inference is obtained via a Gibbs sampler that combines (i) a latent data augmentation step (like in probit regression) and (ii) an existing sampler for a continuous data model.

There are several options for the transformation. First, the transformation can belong to the \*Box-Cox\* family, which includes the known transformations 'identity', 'log', and 'sqrt', as well as a version in which the Box-Cox parameter is inferred within the MCMC sampler ('box-cox'). Second, the transformation can be estimated (before model fitting) using the empirical distribution of the data  $y$ . Options in this case include the empirical cumulative distribution function (CDF), which is fully nonparametric ('np'), or the parametric alternatives based on Poisson ('pois') or Negative-Binomial ('neg-bin') distributions. For the parametric distributions, the parameters of the distribution are estimated using moments (means and variances) of  $y$ .

For this generic function, the Bayesian nonparametric transformation(s) are not available.

## Value

a list with at least the following elements:

- `post.pred`: draws from the posterior predictive distribution of  $y$
- `post.sigma`: draws from the posterior distribution of  $\sigma$
- `post.log.like.point`: draws of the log-likelihood for each of the  $n$  observations
- `WAIC`: Widely-Applicable/Watanabe-Akaike Information Criterion
- `p_waic`: Effective number of parameters based on WAIC
- `post.lambda`: draws from the posterior distribution of  $\lambda$  (NULL unless `transformation='box-cox'`)
- `fitted.values`: the posterior mean of the conditional expectation of the counts  $y$  (NULL if `save_y_hat=FALSE`)
- `post.fitted.values`: posterior draws of the conditional mean of the counts  $y$  (NULL if `save_y_hat=FALSE`)

If the coefficients list from `init_params` and `sample_params` contains a named element `beta`, e.g. for linear regression, then the function output contains

- `coefficients`: the posterior mean of the  $\beta$  coefficients
- `post.beta`: draws from the posterior distribution of  $\beta$
- `post.othercoefs`: draws from the posterior distribution of any other sampled coefficients, e.g. variance terms

If no `beta` exists in the parameter coefficients, then the output list just contains

- `coefficients`: the posterior mean of all coefficients
- `post.beta`: draws from the posterior distribution of all coefficients

Additionally, if `init_params` and `sample_params` have output `mu_test`, then the sampler will output `post.predtest`, which contains draws from the posterior predictive distribution at test points.

**Examples**

```

# Simulate data with count-valued response y:
sim_dat = simulate_nb_lm(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X

# STAR: log-transformation:
fit_log = genMCMC_star(y = y,
                      sample_params = function(y, X, params) sample_lm_gprior(y, X, params),
                      init_params = function(y) init_lm_gprior(y, X),
                      transformation = 'log')

# What is included:
names(fit_log)

# Posterior mean of each coefficient:
coef(fit_log)

# WAIC for STAR-log:
fit_log$WAIC

# MCMC diagnostics:
plot(as.ts(fit_log$post.beta[,1:3]))

# Posterior predictive check:
hist(apply(fit_log$post.pred, 1,
          function(x) mean(x==0)), main = 'Proportion of Zeros', xlab='');
abline(v = mean(y==0), lwd=4, col = 'blue')

```

---

getEffSize

*Summarize of effective sample size*


---

**Description**

Compute the summary statistics for the effective sample size (ESS) across posterior samples for possibly many variables

**Usage**

```
getEffSize(postX)
```

**Arguments**

postX            An array of arbitrary dimension (nsims x ... x ...), where nsims is the number of posterior samples

**Value**

Table of summary statistics using the function summary().

**Examples**

```
# ESS for iid simulations:
library(coda)
rand_iid = rnorm(n = 10^4)
getEffSize(rand_iid)

# ESS for several AR(1) simulations with coefficients 0.1, 0.2,...,0.9:
rand_ar1 = sapply(seq(0.1, 0.9, by = 0.1), function(x) arima.sim(n = 10^4, list(ar = x)))
getEffSize(rand_ar1)
```

---

g\_bc

*Box-Cox transformation*

---

**Description**

Evaluate the Box-Cox transformation, which is a scaled power transformation to preserve continuity in the index lambda at zero. Negative values are permitted.

**Usage**

```
g_bc(t, lambda)
```

**Arguments**

t	argument(s) at which to evaluate the function
lambda	Box-Cox parameter

**Value**

The evaluation(s) of the Box-Cox function at the given input(s) t.

**Note**

Special cases include the identity transformation (lambda = 1), the square-root transformation (lambda = 1/2), and the log transformation (lambda = 0).

**Examples**

```
# Log-transformation:
g_bc(1:5, lambda = 0); log(1:5)

# Square-root transformation: note the shift and scaling
g_bc(1:5, lambda = 1/2); sqrt(1:5)
```

---

`g_cdf`*Cumulative distribution function (CDF)-based transformation*

---

**Description**

Compute a CDF-based transformation using the observed count data. The CDF can be estimated nonparametrically or parametrically based on the Poisson or Negative Binomial distributions. In the parametric case, the parameters are determined based on the moments of  $y$ . Note that this is a fixed quantity and does not come with uncertainty quantification.

**Usage**

```
g_cdf(y, distribution = "np")
```

**Arguments**

<code>y</code>	<code>n x 1</code> vector of observed counts
<code>distribution</code>	the distribution used for the CDF; must be one of <ul style="list-style-type: none"><li>• "np" (empirical CDF)</li><li>• "pois" (moment-matched marginal Poisson CDF)</li><li>• "neg-bin" (moment-matched marginal Negative Binomial CDF)</li></ul>

**Value**

A smooth monotone function which can be used for evaluations of the transformation.

**Examples**

```
# Sample some data:
y = rpois(n = 500, lambda = 5)

# Empirical CDF version:
g_np = g_cdf(y, distribution = 'np')

# Poisson version:
g_pois = g_cdf(y, distribution = 'pois')

# Negative binomial version:
g_negbin = g_cdf(y, distribution = 'neg-bin')

# Plot together:
t = 1:max(y) # grid
plot(t, g_np(t), type='l')
lines(t, g_pois(t), lty = 2)
lines(t, g_negbin(t), lty = 3)
```

---

`g_inv` *Inverse transformation*

---

**Description**

Compute the inverse transformation on a vector of real-valued inputs based on the marginal CDF of  $z$  and the marginal CDF of  $y$ .

**Usage**

```
g_inv(Fz_eval, Fy_eval, y_grid)
```

**Arguments**

<code>Fz_eval</code>	the marginal CDF of $z$ evaluated on some inputs
<code>Fy_eval</code>	the marginal CDF of $y$ evaluated on <code>y_grid</code>
<code>y_grid</code>	a grid of non-negative integers

**Value**

The inverse transformation function evaluated on the same inputs as `Fz_eval`.

**Note**

The inputs for `Fz_eval` do not need to be known to compute the inverse transformation, so they are not required for this function.

The function will return NA if `Fz_eval` is greater than all `Fy_eval` values. When `y_max < Inf`, this implies that the inverse is `y_max`; otherwise, it means that `y_grid` needs to use larger values. Both are handled externally.

---

`g_inv_approx` *Approximate inverse transformation*

---

**Description**

Compute the inverse function of a transformation  $g$  based on a grid search.

**Usage**

```
g_inv_approx(g, t_grid)
```

**Arguments**

<code>g</code>	the transformation function
<code>t_grid</code>	grid of arguments at which to evaluate the transformation function

**Value**

A function which can be used for evaluations of the (approximate) inverse transformation function.

**Examples**

```
# Sample some data:
y = rpois(n = 500, lambda = 5)

# Empirical CDF transformation:
g_np = g_cdf(y, distribution = 'np')

# Grid for approximation:
t_grid = seq(1, max(y), length.out = 100)

# Approximate inverse:
g_inv = g_inv_approx(g = g_np, t_grid = t_grid)

# Check the approximation:
plot(t_grid, g_inv(g_np(t_grid)), type='p')
lines(t_grid, t_grid)
```

---

g\_inv\_bc

*Inverse Box-Cox transformation*


---

**Description**

Evaluate the inverse Box-Cox transformation. Negative values are permitted.

**Usage**

```
g_inv_bc(s, lambda)
```

**Arguments**

s	argument(s) at which to evaluate the function
lambda	Box-Cox parameter

**Value**

The evaluation(s) of the inverse Box-Cox function at the given input(s) s.

**Note**

Special cases include the identity transformation ( $\lambda = 1$ ), the square-root transformation ( $\lambda = 1/2$ ), and the log transformation ( $\lambda = 0$ ).

#' @examples # (Inverse) log-transformation: g\_inv\_bc(1:5, lambda = 0); exp(1:5)

# (Inverse) square-root transformation: note the shift and scaling g\_inv\_bc(1:5, lambda = 1/2); (1:5)^2

---

HPDregion	<i>Compute highest posterior density (HPD) regions</i>
-----------	--

---

**Description**

Given a vector of draws from the posterior (predictive) distribution, compute a prob their cumulative probability exceeds prob.

**Usage**

```
HPDregion(post_pred, prob = 0.95, merge_gaps = FALSE)
```

**Arguments**

post_pred	vector of draws from the posterior (predictive) distribution
prob	numeric scalar in (0,1) giving the target probability content of the region
merge_gaps	logical; if TRUE, add singleton points that are skipped over

**Value**

An ordered vector of values comprising the HPD region

**Note**

The HPD region is not necessarily contiguous. This function is primarily designed for discrete distributions, so that the unique values in post\_pred accumulate multiple realizations. merge\_gaps allows some smoothing over "skipped" points, e.g., {0, 1, 3} becomes {0, 1, 2, 3}.

---

init_lm_gprior	<i>Initialize linear regression parameters assuming a g-prior</i>
----------------	---

---

**Description**

Initialize the parameters for a linear regression model assuming a g-prior for the coefficients.

**Usage**

```
init_lm_gprior(y, X, X_test = NULL)
```

**Arguments**

y	n x 1 vector of data
X	n x p matrix of predictors
X_test	n_test x p matrix of predictors at test points (default is NULL)

**Value**

a named list params containing at least

1. mu: vector of conditional means (fitted values)
2. sigma: the conditional standard deviation
3. coefficients: a named list of parameters that determine mu

Additionally, if X\_test is not NULL, then the list includes an element mu\_test, the vector of conditional means at the test points

**Note**

The parameters in coefficients are:

- beta: the  $p \times 1$  vector of regression coefficients components of beta

**Examples**

```
# Simulate data for illustration:
sim_dat = simulate_nb_lm(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X

# Initialize:
params = init_lm_gprior(y = y, X = X)
names(params)
names(params$coefficients)
```

---

lm\_star

*Fitting frequentist STAR linear model via EM algorithm*


---

**Description**

Compute the MLEs and log-likelihood for the STAR linear model. The regression coefficients are estimated using least squares within an EM algorithm.

**Usage**

```
lm_star(
  formula,
  data = NULL,
  transformation = "np",
  y_max = Inf,
  sd_init = 10,
  tol = 10^-10,
  max_iters = 1000
)
```

**Arguments**

formula	an object of class "formula" (see <code>lm</code> for details on model specification)
data	an optional data frame, list or environment (or object coercible by <code>as.data.frame</code> to a data frame) containing the variables in the model; like <code>lm</code> , if not found in data, the variables are taken from <code>environment(formula)</code>
transformation	transformation to use for the latent data; must be one of <ul style="list-style-type: none"> <li>• "identity" (identity transformation)</li> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>• "box-cox" (box-cox transformation with learned parameter)</li> </ul>
y_max	a fixed and known upper bound for all observations; default is Inf
sd_init	add random noise for EM algorithm initialization scaled by <code>sd_init</code> times the Gaussian MLE standard deviation; default is 10
tol	tolerance for stopping the EM algorithm; default is $10^{-10}$ ;
max_iters	maximum number of EM iterations before stopping; default is 1000

**Details**

Standard function calls including `coefficients`, `fitted`, and `residuals` apply. Fitted values are the expectation at the MLEs, and as such are not necessarily count-valued.

**Value**

an object of class "lmstar", which is a list with the following elements:

- `coefficients` the MLEs of the coefficients
- `fitted.values` the fitted values at the MLEs
- `g.hat` a function containing the (known or estimated) transformation
- `ginv.hat` a function containing the inverse of the transformation
- `sigma.hat` the MLE of the standard deviation
- `mu.hat` the MLE of the conditional mean (on the transformed scale)
- `z.hat` the estimated latent data (on the transformed scale) at the MLEs
- `residuals` the Dunn-Smyth residuals (randomized)
- `residuals_rep` the Dunn-Smyth residuals (randomized) for 10 replicates
- `logLik` the log-likelihood at the MLEs
- `logLik0` the log-likelihood at the MLEs for the \*unrounded\* initialization
- `lambda` the Box-Cox nonlinear parameter
- and other parameters that (1) track the parameters across EM iterations and (2) record the model specifications

**Note**

Infinite latent data values may occur when the transformed Gaussian model is highly inadequate. In that case, the function returns the *\*indices\** of the data points with infinite latent values, which are significant outliers under the model. Deletion of these indices and re-running the model is one option, but care must be taken to ensure that (i) it is appropriate to treat these observations as outliers and (ii) the model is adequate for the remaining data points.

**References**

Kowal, D. R., & Wu, B. (2021). Semiparametric count data regression for self-reported mental health. *Biometrics*. doi:10.1111/biom.13617

**Examples**

```
# Simulate data with count-valued response y:
sim_dat = simulate_nb_lm(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X[,-1] # remove intercept

# Fit model
fit_em = lm_star(y ~ X)

# Fitted coefficients:
coef(fit_em)

# Fitted values:
y_hat = fitted(fit_em)
plot(y_hat, y);

# Residuals:
plot(residuals(fit_em))
qqnorm(residuals(fit_em)); qqline(residuals(fit_em))
```

---

plot\_coef

---

*Plot the estimated regression coefficients and credible intervals*


---

**Description**

Plot the estimated regression coefficients and credible intervals for the linear effects in up to two models.

**Usage**

```
plot_coef(
  post_coefficients_1,
  post_coefficients_2 = NULL,
  alpha = 0.05,
  labels = NULL
)
```

**Arguments**

post_coefficients_1	Nsims x p matrix of simulations from the posterior distribution of the p coefficients, where Nsims is the number of simulations
post_coefficients_2	Nsims x p matrix of simulations from the posterior distribution of the p coefficients from another model
alpha	confidence level for the credible intervals
labels	p dimensional string of labels for the coefficient names

**Value**

A plot of regression coefficients and credible intervals for 1-2 models

---

plot_fitted	<i>Plot the fitted values and the data</i>
-------------	--

---

**Description**

Plot the fitted values, plus pointwise credible intervals, against the data. For simulations, one may use the true values in place of the data.

**Usage**

```
plot_fitted(y, post_y, y_hat = NULL, alpha = 0.05, ...)
```

**Arguments**

y	n x 1 vector of data
post_y	Nsims x n matrix of simulated fitted values, where Nsims is the number of simulations
y_hat	n x 1 vector of fitted values; if NULL, use the pointwise sample mean colMeans(post_y)
alpha	confidence level for the credible intervals
...	other arguments for plotting

**Value**

A plot with the fitted values and the credible intervals against the data

---

plot\_pmf

*Plot the empirical and model-based probability mass functions*


---

**Description**

Plot the empirical probability mass function, i.e., the proportion of data values  $y$  that equal  $j$  for each  $j=0, 1, \dots$ , together with the model-based estimate of the probability mass function based on the posterior predictive distribution.

**Usage**

```
plot_pmf(y, post.pred, error.bars = FALSE, alpha = 0.05)
```

**Arguments**

<code>y</code>	<code>n x 1</code> vector of data
<code>post.pred</code>	nsave draws from the posterior predictive distribution of $y$
<code>error.bars</code>	logical; if TRUE, include errors bars on the model-based PMF
<code>alpha</code>	confidence level for the credible intervals

**Value**

A plot of the empirical PMF of  $y$  along with a PMF estimate from the model posterior predictive distribution

---

predict.lmstar

*Predict method for response in STAR linear model*


---

**Description**

Outputs predicted values based on an lmstar fit and optionally prediction intervals based on the the (plug-in) predictive distribution for the STAR linear model

**Usage**

```
## S3 method for class 'lmstar'
predict(object, newdata = NULL, interval = FALSE, level = 0.95, N = 1000, ...)
```

**Arguments**

object	Object of class "lmstar" as output by <code>lm_star</code>
newdata	An optional matrix/data frame in which to look for variables with which to predict. If omitted, the fitted values are used.
interval	logical; whether or not to include prediction intervals (default FALSE)
level	Level for prediction intervals
N	number of Monte Carlo samples from the posterior predictive distribution used to approximate intervals; default is 1000
...	Ignored

**Details**

If `interval=TRUE`, then `predict.lmstar` uses a Monte Carlo approach to estimating the (plug-in) predictive distribution for the STAR linear model. The algorithm iteratively samples (i) the latent data given the observed data, (ii) the latent predictive data given the latent data from (i), and (iii) (inverse) transforms and rounds the latent predictive data to obtain a draw from the integer-valued predictive distribution.

The appropriate quantiles of these Monte Carlo draws are computed and reported as the prediction interval.

**Value**

Either a vector of predictions (if `interval=FALSE`) or a matrix of predictions and bounds with column names `fit`, `lwr`, and `upr`

**Note**

The "plug-in" predictive distribution is a crude approximation. Better approaches are available using the Bayesian models, e.g. `blm_star`, which provide samples from the posterior predictive distribution.

For highly skewed responses, prediction intervals especially at lower levels may not include the predicted value itself, since the mean is often much larger than the median.

**Examples**

```
# Simulate data with count-valued response y:
x = seq(0, 1, length.out = 100)
y = rpois(n = length(x), lambda = exp(1.5 + 5*(x -.5)^2))

# Estimate model--assume a quadratic effect (better for illustration purposes)
fit = lm_star(y~x+I(x^2), transformation = 'sqrt')

#Compute the predictive draws for the test points (same as observed points here)
#Also compute intervals using plug-in predictive distribution
y_pred = predict(fit, interval=TRUE)

# Plot the results
plot(x, y, ylim = range(y, y_pred), main = 'STAR: Predictions and 95% PI')
```

```
lines(x,y_pred[,"fit"], col='black', type='s', lwd=4)
lines(x, y_pred[,"lwr"], col='darkgray', type='s', lwd=4)
lines(x, y_pred[,"upr"], col='darkgray', type='s', lwd=4)
```

---

pvals	<i>Compute coefficient p-values for STAR linear regression using likelihood ratio test</i>
-------	--

---

### Description

For a linear regression model within the STAR framework, compute p-values for regression coefficients using a likelihood ratio test. It also computes a p-value for excluding all predictors, akin to a (partial) F test.

### Usage

```
pvals(object)
```

### Arguments

object            Object of class "lmstar" as output by [lm\\_star](#)

### Value

a list of p+1 p-values, one for each predictor as well as the joint p-value excluding all predictors

### Examples

```
# Simulate data with count-valued response y:
sim_dat = simulate_nb_lm(n = 100, p = 2)
y = sim_dat$y; X = sim_dat$X[,-1] # remove intercept

# Select a transformation:
transformation = 'np'

#Estimate model
fit = lm_star(y~X, transformation = transformation)

#Compute p-values
pvals(fit)
```

---

randomForest\_star      *Fit Random Forest STAR with EM algorithm*

---

## Description

Compute the MLEs and log-likelihood for the Random Forest STAR model. The STAR model requires a *transformation* and an *estimation function* for the conditional mean given observed data. The transformation can be known (e.g., log or sqrt) or unknown (Box-Cox or estimated nonparametrically) for greater flexibility. The estimator in this case is a random forest. Standard function calls including [fitted](#) and [residuals](#) apply.

## Usage

```
randomForest_star(
  y,
  X,
  X.test = NULL,
  transformation = "np",
  y_max = Inf,
  sd_init = 10,
  tol = 10^-3,
  max_iters = 1000,
  ntree = 500,
  mtry = max(floor(ncol(X)/3), 1),
  nodesize = 5
)
```

## Arguments

y	n x 1 vector of observed counts
X	n x p matrix of predictors
X.test	m x p matrix of out-of-sample predictors
transformation	transformation to use for the latent data; must be one of <ul style="list-style-type: none"> <li>• "identity" (identity transformation)</li> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>• "box-cox" (box-cox transformation with learned parameter)</li> </ul>
y_max	a fixed and known upper bound for all observations; default is Inf
sd_init	add random noise for EM algorithm initialization scaled by sd_init times the Gaussian MLE standard deviation; default is 10

tol	tolerance for stopping the EM algorithm;
max_iters	maximum number of EM iterations before stopping; default is 1000
nree	Number of trees to grow. This should not be set to too small a number, to ensure that every input row gets predicted at least a few times. Default is 500.
mtry	Number of variables randomly sampled as candidates at each split. Default is $p/3$ .
nodesize	Minimum size of terminal nodes. Setting this number larger causes smaller trees to be grown (and thus take less time). Default is 5.

### Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous *\*latent\** data and (2) connecting the latent data to the observed data via a *\*transformation and rounding\** operation.

The expectation-maximization (EM) algorithm is used to produce maximum likelihood estimators (MLEs) for the parameters defined in the The fitted values are computed using out-of-bag samples. As a result, the log-likelihood is based on out-of-bag prediction, and it is similarly straightforward to compute out-of-bag squared and absolute errors.

### Value

a list with the following elements:

- `fitted.values`: the fitted values at the MLEs based on out-of-bag samples (training)
- `fitted.values.test`: the fitted values at the MLEs (testing)
- `g.hat` a function containing the (known or estimated) transformation
- `sigma.hat` the MLE of the standard deviation
- `mu.hat` the MLE of the conditional mean (on the transformed scale)
- `z.hat` the estimated latent data (on the transformed scale) at the MLEs
- `residuals` the Dunn-Smyth residuals (randomized)
- `residuals_rep` the Dunn-Smyth residuals (randomized) for 10 replicates
- `logLik` the log-likelihood at the MLEs
- `logLik0` the log-likelihood at the MLEs for the *\*unrounded\** initialization
- `lambda` the Box-Cox nonlinear parameter
- `rfObj`: the object returned by `randomForest()` at the MLEs
- and other parameters that (1) track the parameters across EM iterations and (2) record the model specifications

### Note

Since the random forest produces random predictions, the EM algorithm will never converge exactly.

Infinite latent data values may occur when the transformed Gaussian model is highly inadequate. In that case, the function returns the *\*indices\** of the data points with infinite latent values, which are significant outliers under the model. Deletion of these indices and re-running the model is one option, but care must be taken to ensure that (i) it is appropriate to treat these observations as outliers and (ii) the model is adequate for the remaining data points.

## References

Kowal, D. R., & Wu, B. (2021). Semiparametric count data regression for self-reported mental health. *Biometrics*. doi:10.1111/biom.13617

## Examples

```
# Simulate data with count-valued response y:
sim_dat = simulate_nb_friedman(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X

# EM algorithm for STAR (using the log-link)
library(randomForest)
fit_em = randomForest_star(y = y, X = X,
                           transformation = 'log')

# Fitted values (out-of-bag)
y_hat = fitted(fit_em)
plot(y_hat, y);

# Residuals:
plot(residuals(fit_em))
qqnorm(residuals(fit_em)); qqline(residuals(fit_em))

# Log-likelihood at MLEs (out-of-bag):
fit_em$logLik
```

---

rdir

*Dirichlet sampler*

---

## Description

Compute one Monte Carlo draw from a Dirichlet distribution

## Usage

```
rdir(params)
```

## Arguments

params            the vector of (nonnegative) Dirichlet parameters

## Details

This function computes one draw from a Dirichlet distribution. The output is on a simplex (non-negative, sum to one) and has the same length as the input params.

**Value**

the vector of the Dirichlet draw on the simplex

**Examples**

```
# Example draw:  
rdir(params = c(1,2,3))
```

---

roaches	<i>Data on the efficacy of a pest management system at reducing the number of roaches in urban apartments.</i>
---------	--

---

**Description**

Data on the efficacy of a pest management system at reducing the number of roaches in urban apartments.

**Usage**

```
roaches
```

**Format**

```
## 'roaches' A data frame with 262 obs. of 6 variables:
```

**y** Number of roaches caught

**roach1** Pretreatment number of roaches

**treatment** Treatment indicator

**senior** Indicator for only elderly residents in building

**exposure2** Number of days for which the roach traps were used

**Source**

Gelman and Hill (2007); package 'rstanarm'

---

round_floor	<i>Rounding function</i>
-------------	--------------------------

---

**Description**

Define the rounding operator associated with the floor function. The function also returns zero whenever the input is negative and caps the value at `y_max`, where `y_max` is a known upper bound on the data `y` (if specified).

**Usage**

```
round_floor(z, y_max = Inf)
```

**Arguments**

<code>z</code>	the real-valued input(s)
<code>y_max</code>	a fixed and known upper bound for all observations; default is <code>Inf</code>

**Value**

The count-valued output(s) from the rounding function.

**Examples**

```
# Floor function:
round_floor(1.5)
round_floor(0.5)

# Special treatment of negative numbers:
round_floor(-1)
```

---

sample_lm_gprior	<i>Sample the linear regression parameters assuming a g-prior</i>
------------------	---

---

**Description**

Sample the parameters for a linear regression model assuming a g-prior for the coefficients. The coefficients and error standard deviation are sampled jointly.

**Usage**

```
sample_lm_gprior(y, X, params, psi = length(y), chXtX = NULL, X_test = NULL)
```

**Arguments**

<code>y</code>	<code>n x 1</code> vector of data
<code>X</code>	<code>n x p</code> matrix of predictors
<code>params</code>	the named list of parameters containing <ol style="list-style-type: none"> <li>1. <code>mu</code>: vector of conditional means (fitted values)</li> <li>2. <code>sigma</code>: the conditional standard deviation</li> <li>3. <code>coefficients</code>: a named list of parameters that determine <code>mu</code></li> </ol>
<code>psi</code>	the prior variance for the g-prior
<code>cholXtX</code>	the <code>p x p</code> matrix of <code>chol(crossprod(X))</code> (one-time cost); if <code>NULL</code> , compute within the function
<code>X_test</code>	matrix of predictors at test points (default is <code>NULL</code> )

**Value**

The updated named list `params` with draws from the full conditional distributions of `sigma` and `coefficients` (along with updated `mu` and `mu_test` if applicable).

**Note**

The parameters in `coefficients` are:

- `beta`: the `p x 1` vector of regression coefficients components of `beta`

**Examples**

```
# Simulate data for illustration:
sim_dat = simulate_nb_lm(n = 100, p = 5)
y = sim_dat$y; X = sim_dat$X
# Initialize:
params = init_lm_gprior(y = y, X = X)
# Sample:
params = sample_lm_gprior(y = y, X = X, params = params)
names(params)
names(params$coefficients)
```

---

 simBaS

---

*Compute Simultaneous Band Scores (SimBaS)*


---

**Description**

Compute simultaneous band scores (SimBaS) from Meyer et al. (2015, Biometrics). SimBaS uses MC(MC) simulations of a function of interest to compute the minimum `alpha` such that the joint credible bands at the `alpha` level do not include zero. This quantity is computed for each grid point (or observation point) in the domain of the function.

**Usage**

```
simBaS(sampFuns)
```

**Arguments**

sampFuns          Nsims x m matrix of Nsims MCMC samples and m points along the curve

**Value**

m x 1 vector of simBaS

**Note**

The input needs not be curves: the simBaS may be computed for vectors to achieve a multiplicity adjustment.

The minimum of the returned value, PsimBaS\_t, over the domain t is the Global Bayesian P-Value (GBPv) for testing whether the function is zero everywhere.

---

simulate\_nb\_friedman    *Simulate count data from Friedman's nonlinear regression*

---

**Description**

Simulate data from a negative-binomial distribution with nonlinear mean function.

**Usage**

```
simulate_nb_friedman(
  n = 100,
  p = 10,
  r_nb = 1,
  b_int = log(1.5),
  b_sig = log(5),
  sigma_true = sqrt(2 * log(1)),
  seed = NULL
)
```

**Arguments**

n                    number of observations

p                    number of predictors

r\_nb                the dispersion parameter of the Negative Binomial dispersion; smaller values imply greater overdispersion, while larger values approximate the Poisson distribution.

b\_int                intercept; default is log(1.5).

b\_sig                regression coefficients for true signals; default is log(5.0).

sigma\_true standard deviation of the Gaussian innovation; default is zero.  
 seed optional integer to set the seed for reproducible simulation; default is NULL which results in a different dataset after each run

### Details

The log-expected counts are modeled using the Friedman (1991) nonlinear function with interactions, possibly with additional Gaussian noise (on the log-scale). We assume that half of the predictors are associated with the response, i.e., true signals. For sufficiently large dispersion parameter  $r_{nb}$ , the distribution will approximate a Poisson distribution. Here, the predictor variables are simulated from independent uniform distributions.

### Value

A named list with the simulated count response  $y$ , the simulated design matrix  $X$ , and the true expected counts  $Ey$ .

### Note

Specifying  $\text{sigma\_true} = \sqrt{2 \cdot \log(1 + a)}$  implies that the expected counts are inflated by  $100 \cdot a\%$  (relative to  $\exp(X \cdot \beta)$ ), in addition to providing additional overdispersion.

### Examples

```
# Simulate and plot the count data:
sim_dat = simulate_nb_friedman(n = 100, p = 10);
plot(sim_dat$y)
```

---

simulate_nb_lm	<i>Simulate count data from a linear regression</i>
----------------	---

---

### Description

Simulate data from a negative-binomial distribution with linear mean function.

### Usage

```
simulate_nb_lm(
  n = 100,
  p = 10,
  r_nb = 1,
  b_int = log(1.5),
  b_sig = log(2),
  sigma_true = sqrt(2 * log(1)),
  ar1 = 0,
  seed = NULL
)
```

**Arguments**

n	number of observations
p	number of predictors (including the intercept)
r_nb	the dispersion parameter of the Negative Binomial dispersion; smaller values imply greater overdispersion, while larger values approximate the Poisson distribution.
b_int	intercept; default is $\log(1.5)$ , which implies the expected count is 1.5 when all predictors are zero
b_sig	regression coefficients for true signals; default is $\log(2.0)$ , which implies a twofold increase in the expected counts for a one unit increase in x
sigma_true	standard deviation of the Gaussian innovation; default is zero.
ar1	the autoregressive coefficient among the columns of the X matrix; default is zero.
seed	optional integer to set the seed for reproducible simulation; default is NULL which results in a different dataset after each run

**Details**

The log-expected counts are modeled as a linear function of covariates, possibly with additional Gaussian noise (on the log-scale). We assume that half of the predictors are associated with the response, i.e., true signals. For sufficiently large dispersion parameter `r_nb`, the distribution will approximate a Poisson distribution. Here, the predictor variables are simulated from independent standard normal distributions.

**Value**

A named list with the simulated count response `y`, the simulated design matrix `X` (including an intercept), the true expected counts `Ey`, and the true regression coefficients `beta_true`.

**Note**

Specifying `sigma_true = sqrt(2*log(1 + a))` implies that the expected counts are inflated by `100*a%` (relative to  $\exp(X*\beta)$ ), in addition to providing additional overdispersion.

**Examples**

```
# Simulate and plot the count data:
sim_dat = simulate_nb_lm(n = 100, p = 10);
plot(sim_dat$y)
```

---

spline\_star

*Posterior and predictive inference for Bayesian STAR splines*


---

### Description

Compute samples from the posterior and predictive distributions of a STAR spline regression model using either a Gibbs sampling approach or exact Monte Carlo sampling. Cubic B-splines are used with a prior that penalizes roughness.

### Usage

```
spline_star(
  y,
  x = NULL,
  x_test = NULL,
  transformation = "bnp",
  y_max = Inf,
  psi = NULL,
  nbasis = NULL,
  use_MCMC = TRUE,
  nsave = 1000,
  nburn = 1000,
  nskip = 0,
  alpha = 1,
  F0 = NULL,
  verbose = TRUE
)
```

### Arguments

y	n x 1 vector of observed counts
x	n x 1 vector of observation points; if NULL, assume equally-spaced on [0,1]
x_test	n_test x 1 vector of testing points; default is x
transformation	transformation to use for the latent data; must be one of <ul style="list-style-type: none"> <li>• "identity" (identity transformation)</li> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> <li>• "box-cox" (box-cox transformation with learned parameter)</li> <li>• "bnp" (Bayesian nonparametric transformation)</li> </ul>
y_max	a fixed and known upper bound for all observations; default is Inf

psi	prior variance (1/smoothing parameter); if NULL, sample this parameter
nbasis	number of spline basis functions; if NULL, use the default from spikeSlabGAM: : sm
use_MCMC	logical; whether to run Gibbs sampler or Monte Carlo (default is TRUE)
nsave	number of MC(MC) iterations to save
nburn	number of MCMC iterations to discard
nskip	number of MCMC iterations to skip between saving iterations, i.e., save every (nskip + 1)th draw
alpha	prior precision for the Dirichlet Process prior ('bnp' transformation only); default is one
F0	function to evaluate the base measure CDF supported on $\{\theta, \dots, y_{\max}\}$ ('bnp' transformation only)
verbose	logical; if TRUE, print time remaining

## Details

STAR defines a count-valued probability model by (1) specifying a Gaussian model for continuous *latent* data and (2) connecting the latent data to the observed data via a *transformation and rounding* operation. Here, the continuous latent data model is a spline regression.

There are several options for the transformation. First, the transformation can belong to the *Box-Cox* family, which includes the known transformations 'identity', 'log', and 'sqrt', as well as a version in which the Box-Cox parameter is inferred within the MCMC sampler ('box-cox').

Second, the transformation can be estimated (before model fitting) using the data  $y$ . Options in this case include the empirical cumulative distribution function (ECDF), which is fully nonparametric ('np'), or the parametric alternatives based on Poisson ('pois') or Negative-Binomial ('neg-bin') distributions. For the parametric distributions, the parameters of the distribution are estimated using moments (means and variances) of  $y$ .

Lastly, the transformation can be modeled nonparametrically using Bayesian nonparametrics via Dirichlet processes ('bnp'). The 'bnp' option is the default because it is highly flexible, accounts for uncertainty when the transformation is unknown, and is computationally efficient.

The Monte Carlo sampler (use\_MCMC=FALSE) produces direct, joint draws from the posterior predictive distribution. When  $n$  is moderate to large, MCMC sampling (use\_MCMC=TRUE) is much faster and more convenient.

## Value

a list with the following elements:

- `coefficients`: the posterior mean of the spline coefficients
- `fitted.values` the posterior predictive mean at the test points `x_test`
- `post.beta`: `nsave`  $\times$  `p` samples from the posterior distribution of the regression coefficients
- `post.pred`: `nsave`  $\times$  `n_test` samples from the posterior predictive distribution at `x_test`
- `post.psi`: `nsave` draws from the prior variance (inverse smoothing) `psi`
- `marg_like`: the marginal likelihood (only if use\_MCMC=FALSE; otherwise NULL)

**Examples**

```

# Simulate some data:
n = 100
x = seq(0,1, length.out = n)
y = round_floor(exp(1 + rnorm(n)/4 + poly(x, 4)%*%rnorm(n=4, sd = 4:1)))

# Sample from the predictive distribution of a STAR spline model:
fit = spline_star(y = y, x = x)

# Compute 90% prediction intervals:
pi_y = t(apply(fit$post.pred, 2, quantile, c(0.05, .95)))

# Plot the results: intervals, median, and smoothed mean
plot(x, y, ylim = range(pi_y, y))
polygon(c(x, rev(x)),c(pi_y[,2], rev(pi_y[,1])),col='gray', border=NA)
lines(x, apply(fit$post.pred, 2, median), lwd=5, col='black')
lines(x, smooth.spline(x, apply(fit$post.pred, 2, mean))$y, lwd=5, col='blue')
lines(x, y, type='p')

```

warpDLM

*Posterior Inference for warpDLM model with latent structural DLM***Description**

This function outputs posterior quantities and forecasts from a univariate warpDLM model. Currently two latent DLM specifications are supported: local level and the local linear trend.

**Usage**

```

warpDLM(
  y,
  type = c("level", "trend"),
  transformation = c("np", "identity", "log", "sqrt", "pois", "neg-bin"),
  y_max = Inf,
  R0 = 10,
  nsave = 5000,
  nburn = 5000,
  nskip = 1,
  n.ahead = 1
)

```

**Arguments**

**y** the count-valued time series

**type** the type of latent DLM (must be either level or trend)

**transformation** transformation to use for the latent process (default is np); must be one of

	<ul style="list-style-type: none"> <li>• "identity" (identity transformation)</li> <li>• "log" (log transformation)</li> <li>• "sqrt" (square root transformation)</li> <li>• "np" (nonparametric transformation estimated from empirical CDF)</li> <li>• "pois" (transformation for moment-matched marginal Poisson CDF)</li> <li>• "neg-bin" (transformation for moment-matched marginal Negative Binomial CDF)</li> </ul>
y_max	a fixed and known upper bound for all observations; default is Inf
R0	the variance for the initial state $\theta_0$ ; default is 10
nsave	number of MCMC iterations to save
nburn	number of MCMC iterations to discard
nskip	number of MCMC iterations to skip between saving iterations, i.e., save every (nskip + 1)th draw
n.ahead	number of steps to forecast ahead

**Value**

A list with the following elements:

- V\_post: posterior draws of the observation variance
- W\_post: posterior draws of the state update variance(s)
- fc\_post: draws from the forecast distribution (of length n.ahead)
- post\_pred: draws from the posterior predictive distribution of y
- g\_func: transformation function
- g\_inv\_func: inverse transformation function
- KFAS\_mod: the final KFAS model representing the latent DLM

# Index

- \* **datasets**
  - roaches, [39](#)
- a\_j, [3](#)
- bam\_star, [4](#)
- bart\_star, [6](#)
- blm\_star, [10, 34](#)
- coefficients, [30](#)
- confint.lmstar, [13](#)
- credBands, [14](#)
- ergMean, [14](#)
- fitted, [15, 30, 36](#)
- formula, [30](#)
- g\_bc, [24](#)
- g\_cdf, [25](#)
- g\_inv, [26](#)
- g\_inv\_approx, [26](#)
- g\_inv\_bc, [27](#)
- gbm\_star, [15](#)
- genEM\_star, [18](#)
- genMCMC\_star, [20](#)
- getEffSize, [23](#)
- HPDregion, [28](#)
- init\_lm\_gprior, [28](#)
- lm, [30](#)
- lm\_star, [13, 29, 34, 35](#)
- plot\_coef, [31](#)
- plot\_fitted, [32](#)
- plot\_pmf, [33](#)
- predict.lmstar, [33](#)
- pvals, [35](#)
- randomForest\_star, [36](#)
- rdir, [38](#)
- residuals, [15, 30, 36](#)
- roaches, [39](#)
- round\_floor, [40](#)
- sample\_lm\_gprior, [40](#)
- simBaS, [41](#)
- simulate\_nb\_friedman, [42](#)
- simulate\_nb\_lm, [43](#)
- spline\_star, [45](#)
- warpDLM, [47](#)