

Package: SAFEPG (via r-universe)

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Type Package

Title A Novel SAFE Model for Predicting Climate-Related Extreme Losses

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Description The goal of 'SAFEPG' is to predict climate-related extreme losses by fitting a frequency-severity model. It improves predictive performance by introducing a sign-aligned regularization term, which ensures consistent signs for the coefficients across the frequency and severity components. This enhancement not only increases model accuracy but also enhances its interpretability, making it more suitable for practical applications in risk assessment.

Depends R (>= 3.5.0)

License GPL-2

Imports stats, Matrix, methods

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coef.safe	<i>Extract Model Coefficients from a 'safe' Object</i>
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Description

Retrieves the coefficients at specified values of 'lambda' from a fitted 'safe()' model.

Usage

```
## S3 method for class 'safe'
coef(object, s = NULL, ...)
```

Arguments

object	Fitted 'safe()' object.
s	Values of the penalty parameter 'lambda' for which coefficients are requested. Defaults to the entire sequence used during the model fit.
...	Not used.

Details

This function extracts coefficients for specified 'lambda' values from a 'safe()' object. If 's', the vector of 'lambda' values, contains values not originally used in the model fitting, the 'coef' function employs linear interpolation between the closest 'lambda' values from the original sequence to estimate coefficients at the new 'lambda' values.

Value

Returns a matrix or vector of coefficients corresponding to the specified 'lambda' values.

See Also

[safe](#), [predict.safe](#)

Examples

```

set.seed(1)
n <- 100
p <- 5
x <- matrix(rnorm(n * p), nrow = n, ncol = p)
beta_true <- rep(0.1, 5)
gamma_true <- c(rep(1, 3), -1, -1)
mu <- x %*% beta_true
k <- rpois(n, lambda = exp(mu))
alpha_val <- 1
theta <- exp(x %*% gamma_true) / alpha_val
y <- rgamma(n, shape = alpha_val, scale = theta)
lambda_val <- 1
fit <- safe(x, y, k, 1, ind_p = c(1, 1, 1, 0, 0))
coef(fit)

```

eccv.safe

Cross-validation for selecting the tuning parameter of the SAFE model

Description

Performs "electoral college" cross-validation for the [safe](#) function. Unlike a standard cross-validation approach, this method repeats the random partitioning into folds multiple times (controlled by `rep`), then selects the tuning parameter `lambda` via a majority vote across these repeated cross-validation runs. This function is largely similar [`glmnet::cv.glmnet()`].

Usage

```
eccv.safe(x, y, k, lambda, ind_p, rep = 24, nfolds = 5L, ...)
```

Arguments

<code>x</code>	A numeric matrix of dimensions $n \times p$, where n is the number of observations and p is the number of predictors.
<code>y</code>	A numeric vector of length n , representing the loss values (severity). These are assumed to follow a Gamma distribution with shape parameter <code>alpha</code> and scale parameter $\theta = \exp(x^\top \gamma) / \alpha$.
<code>k</code>	A numeric vector of length n , representing the number of claims (frequency), assumed to follow a Poisson distribution with mean $\mu = x^\top \beta$.
<code>lambda</code>	A user-supplied numeric vector of tuning parameters. The function will compute the solution for each value in <code>lambda</code> .
<code>ind_p</code>	A user-provided index or indicator specifying which predictors should share the same sign across frequency and severity.

rep	The number of repeated cross-validation cycles (default is 24). In each cycle, observations are randomly assigned to nolds folds, cross-validation is performed for all lambda values, and a "vote" is cast for the lambda with the lowest error in that cycle.
nolds	The number of folds in cross-validation. Default is 5.
...	Additional arguments passed to safe .

Details

The function computes the average cross-validation error and reports the best lambda that achieves the smallest cross-validation error.

Value

An object of class "safe", which is a list containing:

lambda.min	The tuning parameter lambda value that won the most votes (i.e., achieved the minimum cross-validation error most often).
lambda	The full sequence of candidate lambda values provided to the function.
ncvmat	A numeric matrix of dimension rep * the length of lambda, containing the cross-validation errors for each lambda across all rep runs.

Examples

```
set.seed(1)
n <- 100
p <- 5
x <- matrix(rnorm(n * p), nrow = n, ncol = p)
beta_true <- rep(0.1, 5)
gamma_true <- c(rep(1, 3), -1, -1)
mu <- x %*% beta_true
k <- rpois(n, lambda = exp(mu))
alpha_val <- 1
theta <- exp(x %*% gamma_true) / alpha_val
y <- rgamma(n, shape = alpha_val, scale = theta)
lambda_seq <- 10^seq(2, -8, length.out = 5)
fit <- eccv.safe(x, y, k, lambda=lambda_seq, ind_p = c(1, 1, 1, 0, 0))
```

predict.safe

Make Predictions from a 'safe' Object

Description

Produces fitted values for new predictor data using a fitted 'safe()' object.

Usage

```
## S3 method for class 'safe'
predict(object, newx, s = NULL, ...)
```

Arguments

object	Fitted 'safe()' object from which predictions are to be derived.
newx	Matrix of new predictor values for which predictions are desired. This must be a matrix and is a required argument.
s	Values of the penalty parameter 'lambda' for which predictions are requested. Defaults to the entire sequence used during the model fit.
...	Not used.

Details

This function generates predictions at specified 'lambda' values from a fitted 'safe()' object. It is essential to provide a new matrix of predictor values ('newx') at which these predictions are to be made.

Value

Returns the fitted values.

See Also

[safe](#), [coef.safe](#)

Examples

```
set.seed(1)
n <- 100
p <- 5
x <- matrix(rnorm(n * p), nrow = n, ncol = p)
beta_true <- rep(0.1, 5)
gamma_true <- c(rep(1, 3), -1, -1)
mu <- x %*% beta_true
k <- rpois(n, lambda = exp(mu))
alpha_val <- 1
theta <- exp(x %*% gamma_true) / alpha_val
y <- rgamma(n, shape = alpha_val, scale = theta)
lambda_val <- 1
fit <- safe(x, y, k, 1, ind_p = c(1, 1, 1, 0, 0))
set.seed(234)
newx <- matrix(rnorm(n * p), nrow = n, ncol = p)
predict(fit, newx)
```

 safe

Solve the sign-aligned frequency-severity (SAFE) model

Description

This function fits a Poisson-Gamma regression framework that aligns the signs of certain predictors in both the Poisson frequency component and the Gamma severity component. The solution path is computed at a sequence of tuning parameter values (λ).

Usage

```
safe(
  x,
  y,
  k,
  lambda,
  alpha = 1,
  eps = 1e-08,
  maxit = 1e+05,
  eps2 = 1e-08,
  ind_p,
  int_gam = NULL,
  int_beta = NULL
)
```

Arguments

x	A numeric matrix of dimensions $n \times p$, where n is the number of observations and p is the number of predictors.
y	A numeric vector of length n , representing the loss values (severity). These are assumed to follow a Gamma distribution with shape parameter α and scale parameter $\theta = \exp(x^\top \gamma) / \alpha$.
k	A numeric vector of length n , representing the number of claims (frequency), assumed to follow a Poisson distribution with mean $\mu = x^\top \beta$.
lambda	A user-supplied numeric vector of tuning parameters. The function will compute the solution for each value in <code>lambda</code> .
alpha	The shape parameter for the Gamma distribution (default is 1).
eps	Convergence tolerance for updating beta (default is $1e-8$).
maxit	An integer specifying the maximum number of iterations (default is $1e5$).
eps2	Convergence tolerance for updating gamma (default is $1e-8$).
ind_p	A user-provided vector specifying which predictors should share the same sign across frequency and severity.
int_gam	Optional numeric vector or matrix providing initial values for gamma; defaults to NULL.
int_beta	Optional numeric vector or matrix providing initial values for beta; defaults to NULL.

Details

The function uses an **accelerated proximal gradient descent** algorithm to simultaneously estimate β (for the Poisson frequency model) and γ (for the Gamma severity model) under a sign-alignment constraint for selected predictors (controlled by `ind_p`).

Value

An object of class `safe`, which is a list containing:

<code>beta</code>	A $p \times L$ matrix of frequency-model coefficients, where p is the number of predictors and L is the number of lambda values.
<code>gamma</code>	A $p \times L$ matrix of severity-model coefficients, with the same dimensions as <code>beta</code> .
<code>lambda</code>	The (sorted) sequence of lambda values used in the estimation.
<code>npass_beta</code>	Total number of iterations used to update <code>beta</code> across all lambda values.
<code>npass_gamma</code>	Total number of iterations used to update <code>gamma</code> across all lambda values.
<code>jerr</code>	An integer flag for warnings or errors; 0 indicates no issues encountered.

Examples

```
set.seed(1)
n <- 100
p <- 5
x <- matrix(rnorm(n * p), nrow = n, ncol = p)
beta_true <- rep(0.1, 5)
gamma_true <- c(rep(1, 3), -1, -1)
mu <- x %*% beta_true
k <- rpois(n, lambda = exp(mu))
alpha_val <- 1
theta <- exp(x %*% gamma_true) / alpha_val
y <- rgamma(n, shape = alpha_val, scale = theta)
lambda_val <- 1
fit <- safe(x, y, k, 1, ind_p = c(1, 1, 1, 0, 0))
```

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