

Package ‘BSGW’

May 6, 2026

Type Package

Title Bayesian Survival Model with Lasso Shrinkage Using Generalized Weibull Regression

Version 0.9.4

Date 2022-12-12

Author Alireza S. Mahani, Mansour T.A. Sharabiani

Maintainer Alireza S. Mahani <alireza.s.mahani@gmail.com>

Description Bayesian survival model using Weibull regression on both scale and shape parameters. Dependence of shape parameter on covariates permits deviation from proportional-hazard assumption, leading to dynamic - i.e. non-constant with time - hazard ratios between subjects. Bayesian Lasso shrinkage in the form of two Laplace priors - one for scale and one for shape coefficients - allows for many covariates to be included. Cross-validation helper functions can be used to tune the shrinkage parameters. Monte Carlo Markov Chain (MCMC) sampling using a Gibbs wrapper around Radford Neal's univariate slice sampler (R package MfUSampler) is used for coefficient estimation.

License GPL (>= 2)

Imports foreach, doParallel, survival, MfUSampler, methods

NeedsCompilation no

Repository CRAN

Date/Publication 2022-12-12 13:10:08 UTC

Contents

bsgw	2
bsgw.crossval	6
plot.bsgw	8
predict.bsgw	9
summary.bsgw	11

Index	13
--------------	-----------

Description

Bayesian survival model - with stratification and shrinkage - using Weibull regression on both scale and shape parameters, resulting in time-dependent (i.e. dynamic) hazard ratios.

Usage

```
bsgw(formula, data, formulas=formula, weights, subset, na.action=na.fail, init="survreg"
     , ordweib=FALSE, scale=0, control=bsgw.control(), print.level=2)
bsgw.control(scalex=TRUE, iter=1000, burnin=round(iter/2), sd.thresh=1e-4
     , lambda=0.0, lambdas=lambda, nskip=round(iter/10), alpha.min=0.1, alpha.max=10.0
     , beta.max=log(20), betas.max=5.0, memlim.gb=8)
## S3 method for class 'bsgw'
print(x, ...)
```

Arguments

formula	Survival formula expressing the time/status variables as well as covariates used in regression on scale parameter. Currently, only right and left censoring is supported. Must include intercept term.
data	Data frame containing the covariates and response variable.
formulas	Formula expressing the covariates used in regression on shape parameter. No left-hand side is necessary since the response variable information is extracted from formula. Default value is formula. Must include intercept term.
weights	Optional vector of case weights. *Not supported yet*
subset	Subset of the observations to be used in the fit. *Not supported yet*
na.action	Missing-data filter function. *Not supported yet (only na.fail behavior works)*
init	Initialization behavior. Currently, three options are supported: 1) If <code>init="survreg"</code> , an ordinary Weibull regression is performed and coefficients are used to initialize the bsgw MCMC run. 2) If <code>init</code> is a <code>survreg</code> object, e.g. from a previous Weibull regression fit, the object can be directly passed as parameter. 3) If <code>init</code> is any other value, or if <code>survreg</code> produces error or warning, we simply set all coefficients to zero.
ordweib	If <code>TRUE</code> , a Bayesian ordinary Weibull model is estimated, in which any covariates in <code>formulas</code> are stripped away, and the inverse-logit transformation in the shape-parameter regression is replaced with a simple exponential transformation. If shrinkage parameters are kept at 0, the result is a Bayesian equivalent of an ordinary Weibull regression.
scale	If <code>scale>0</code> , the value of the shape parameter is fixed, i.e. not estimated from data.
control	See <code>bsgw.control</code> for a description of the parameters inside the control list.

<code>print.level</code>	Controlling verbosity level.
<code>scalex</code>	If TRUE, each covariate vector is centered and scaled before model estimation. The scaling parameters are saved in return object, and used in subsequent calls to <code>predict</code> function. Users are strongly advised against turning this feature off, since the quality of Gibbs sampling MCMC is greatly enhanced by covariate centering and scaling.
<code>iter</code>	Number of MCMC samples to draw.
<code>burnin</code>	Number of initial MCMC samples to discard before calculating summary statistics.
<code>sd.thresh</code>	Threshold for standard deviation of a covariate (after possible centering/scaling). If below the threshold, the corresponding coefficient is removed from sampling, i.e. its value is clamped to zero.
<code>lambda</code>	Bayesian Lasso shrinkage parameter for scale-parameter coefficients.
<code>lambdas</code>	Bayesian Lasso shrinkage parameter for shape-parameter coefficients.
<code>nskip</code>	Controlling how often to print progress report during MCMC run. For example, if <code>nskip=10</code> , progress will be reported after 10,20,30,... samples.
<code>alpha.min</code>	Lower bound on the shape parameter.
<code>alpha.max</code>	Upper bound on the shape parameter.
<code>beta.max</code>	Upper bound on absolute value of coefficients of scale parameter (with the exception of the intercept).
<code>betas.max</code>	Upper bound on absolute value of coefficients of shape parameter (with the exception of the intercept).
<code>memlim.gb</code>	User-specified limit on total memory (in GB) available during prediction. Hazard, cumulative hazard, and survival prediction objects are all three-dimensional arrays which can quickly grow very large, depending on data length, number of MCMC samples collected, and number of time points along which prediction is made.
<code>x</code>	Object of class 'bsgw', usually the result of a call to the <code>bsgw</code> .
<code>...</code>	Arguments to be passed to/from other methods.

Value

The function `bsgw.control` returns a list with elements identical to the input parameters. The function `bsgw` returns an object of class `bsgw`, with the following components:

<code>call</code>	The matched call.
<code>formula</code>	Same as input.
<code>formulas</code>	Same as input.
<code>weights</code>	Same as input. *Not supported yet*
<code>subset</code>	Same as input. *Not supported yet*
<code>na.action</code>	Same as input. *Not supported yet* (current behavior is <code>na.fail</code>)
<code>init</code>	Initial values for scale and shape coefficients used in MCMC sampling, either by performing an ordinary Weibull regression or by extracting estimated coefficients from a previously-performed such regression.

<code>ordweib</code>	Same as input.
<code>survreg.scale.ref</code>	Value of scale parameter, estimated using ordinary Weibull regression by calling the <code>survreg</code> function in the <code>survival</code> package.
<code>ordreg</code>	The "survreg" object returned from calling the same function for initialization of coefficients.
<code>scale</code>	Same as input.
<code>control</code>	Same as input.
<code>X</code>	Model matrix used for regression on scale parameter, after potential centering and scaling. The corresponding vector of coefficients is called <code>beta</code> .
<code>Xs</code>	Model matrix used for regression on shape parameter, after potential centering and scaling. The corresponding vector of coefficients is called <code>betas</code> .
<code>y</code>	Survival response variable (time and status) used in the model.
<code>contrasts</code>	The contrasts used for scale-parameter coefficients.
<code>contrastss</code>	The contrasts used for shape-parameter coefficients.
<code>xlevels</code>	A record of the levels of the factors used in fitting for scale parameter regression.
<code>xlevelss</code>	A record of the levels of the factors used in fitting for shape parameter regression.
<code>terms</code>	The terms object used for scale parameter regression.
<code>termss</code>	The terms object used for shape parameter regression.
<code>colnamesX</code>	Names of columns for <code>X</code> , also names of scale coefficients.
<code>colnamesXs</code>	Names of columns for <code>Xs</code> , also names of shape coefficients.
<code>apply.scale.X</code>	Index of columns of <code>X</code> where scaling has been applied.
<code>apply.scale.Xs</code>	Index of columns of <code>Xs</code> where scaling has been applied.
<code>centerVec.X</code>	Vector of centering parameters for columns of <code>X</code> indicated by <code>apply.scale.X</code> .
<code>scaleVec.X</code>	Vector of scaling parameters for columns of <code>X</code> indicated by <code>apply.scale.X</code> .
<code>centerVec.Xs</code>	Vector of centering parameters for columns of <code>Xs</code> indicated by <code>apply.scale.Xs</code> .
<code>scaleVec.Xs</code>	Vector of scaling parameters for columns of <code>Xs</code> indicated by <code>apply.scale.Xs</code> .
<code>idx</code>	Vector of indexes into <code>X</code> for which sampling occurred. All columns of <code>X</code> whose standard deviation falls below <code>sd.thresh</code> are excluded from sampling and their corresponding coefficients are clamped to \emptyset .
<code>idxs</code>	Vector of indexes into <code>Xs</code> for which sampling occurred. All columns of <code>Xs</code> whose standard deviation falls below <code>sd.thresh</code> are excluded from sampling and their corresponding coefficients are clamped to \emptyset .
<code>median</code>	List of median values, with elements including <code>beta</code> (coefficients of scale regression), <code>betas</code> (coefficients of shape regression), <code>survreg.scale</code> (value of <code>surgreg</code> -style scale parameter for all training set observations).
<code>smp</code>	List of coefficient samples, with the following elements: 1) <code>beta</code> (scale parameter coefficients), 2) <code>betas</code> (shape parameter coefficients), 3) <code>lp</code> (vector of linear predictor for scale parameter, within-sample), 4) <code>loglike</code> (log-likelihood of

model), 5) logpost (log-posterior of mode, i.e. log-likelihood plus the shrinkage term). The last two entities are used during within-sample prediction of response, i.e. during a subsequent call to `predict`. Each parameter has `control$iter` samples.

<code>km.fit</code>	Kaplan-Meier fit to training data. Used in plot.bsgw method.
<code>tmax</code>	Maximum time value in training set. Used in predict.bsgw for automatic selection of the <code>tvec</code> parameter.

Author(s)

Alireza S. Mahani, Mansour T.A. Sharabiani

References

Mazucheli J., Louzada-Neto F. and Achnar J.A. (2002). Lifetime models with nonconstant shape parameters. *Confiabilidade. III Jornada Regional de Estatística e II Semana da Estatística, Maringá*.

Neal R.M. (2003). Slice Sampling. *Annals of Statistics*, **31**, 705-767.

Park T. and Casella G. (2008). The Bayesian Lasso. *Journal of the American Statistical Association*, **103**, 681-686.

See Also

For calculating median and lower/upper bounds on coefficients, use [summary.bsgw](#).

For prediction, use [predict.bsgw](#).

Examples

```
## model estimation using 800 samples, printing progress every 100 samples
library("survival")
data(ovarian)
est <- bsgw(Surv(futime, fustat) ~ ecog.ps + rx, ovarian
            , control=bsgw.control(iter=400, nskip=100))

## comparing shape of Weibull curves between ordinary Weibull and bsgw
## since in bsgw shape is dependent on covariates, only a population average is meaningful
## Note that survreg-style scale is inverse of bsgw shape parameter, see survreg help page
west <- survreg(Surv(futime, fustat) ~ ecog.ps + rx, ovarian)
cat("constant survreg-style scale parameter:", west$scale, "\n")
cat("population average of survreg-style scale parameter from bsgw model:"
    , mean(est$median$survreg.scale), "\n")
```

bsgw.crossval	<i>Convenience functions for cross-validation-based selection of shrinkage parameter in the bsgw model.</i>
---------------	---

Description

bsgw.crossval calculates cross-validation-based, out-of-sample log-likelihood of a bsgw model for a data set, given the supplied folds. bsgw.crossval.wrapper applies bsgw.crossval to a set of combinations of shrinkage parameters (lambda, lambdas) and produces the resulting vector of log-likelihood values as well as the specific combination of shrinkage parameters associated with the maximum log-likelihood. bsgw.generate.folds generates random partitions, while bsgw.generate.folds.eventbalanced generates random partitions with events evenly distributed across partitions. The latter feature is useful for cross-validation of small data sets with low event rates, since it prevents over-accumulation of events in one or two partitions, and lack of events altogether in other partitions.

Usage

```
bsgw.generate.folds(ntot, nfold=5)
bsgw.generate.folds.eventbalanced(formula, data, nfold=5)
bsgw.crossval(data, folds, all=FALSE, print.level=1
, control=bsgw.control(), ncores=1, ...)
bsgw.crossval.wrapper(data, folds, all=FALSE, print.level=1
, control=bsgw.control(), ncores=1
, lambda.vec=exp(seq(from=log(0.01), to=log(100), length.out = 10)), lambdas.vec=NULL
, lambda2=if (is.null(lambdas.vec)) cbind(lambda=lambda.vec, lambdas=lambda.vec)
else as.matrix(expand.grid(lambda=lambda.vec, lambdas=lambdas.vec))
, plot=TRUE, ...)
```

Arguments

ntot	Number of observations to create partitions for. It must typically be set to nrow(data).
nfold	Number of folds or partitions to generate.
formula	Survival formula, used to extract the binary status field from the data. Right-hand side of the formula is ignored, so a formula of the form Surv(time, status)~1 is sufficient.
data	Data frame used in model training and prediction.
folds	An integer vector of length nrow(data), defining fold/partition membership of each observation. For example, in 5-fold cross-validation for a data set of 200 observations, folds must be a 200-long vector with elements from the set {1, 2, 3, 4, 5}. Convenience functions bsgw.generate.folds and bsgw.generate.folds.eventbalanced can be used to generate the folds vector for a given survival data frame.
all	If TRUE, estimation objects from each cross-validation task is collected and returned for diagnostics purposes.

<code>print.level</code>	Verbosity of progress report.
<code>control</code>	List of control parameters, usually the output of <code>bsgw.control</code> .
<code>ncores</code>	Number of cores for parallel execution of cross-validation code.
<code>lambda.vec</code>	Vector of shrinkage parameters to be tested for scale-parameter coefficients.
<code>lambdas.vec</code>	Vector of shrinkage parameters to be tested for shape-parameter coefficients.
<code>lambda2</code>	A data frame that enumerates all combinations of <code>lambda</code> and <code>lambdas</code> to be tested. By default, it is constructed from forming all permutations of <code>lambda.vec</code> and <code>lambdas.vec</code> . If <code>lambdas.vec=NULL</code> , it will only try equal values of the two parameters in each combination.
<code>plot</code>	If TRUE, and if the <code>lambda</code> and <code>lambdas</code> entries in <code>lambda2</code> are identical, a plot of <code>loglike</code> as a function of either vector is produced.
<code>...</code>	Other arguments to be passed to <code>bsgw</code> .

Value

Functions `bsgw.generate.folds` and `bsgw.generate.folds.eventbalanced` produce integer vectors of length `ntot` or `nrow(data)` respectively. The output of these functions can be directly passed to `bsgw.crossval` or `bsgw.crossval.wrapper`. Function `bsgw.crossval` returns the log-likelihood of data under the assumed `bsgw` model, calculated using a cross-validation scheme with the supplied `fold` parameter. If `all=TRUE`, the estimation objects for each of the `nfold` estimation jobs will be returned as the "estobj" attribute of the returned value. Function `bsgw.crossval.wrapper` returns a list with elements `lambda` and `lambdas`, the optimal shrinkage parameters for scale and shape coefficients, respectively. Additionally, the following attributes are attached:

<code>loglike.vec</code>	Vector of log-likelihood values, one for each tested combination of <code>lambda</code> and <code>lambdas</code> .
<code>loglike.opt</code>	The maximum log-likelihood value from the <code>loglike.vec</code> .
<code>lambda2</code>	Data frame with columns <code>lambda</code> and <code>lambdas</code> . Each row of this data frame contains one combination of shrinkage parameters that are tested in the wrapper function.
<code>estobj</code>	If <code>all=TRUE</code> , a list of length <code>nrow(lambda2)</code> is returned, with each element being itself a list of <code>nfold</code> estimation objects associated with each call to the <code>bsgw</code> function. This object can be examined by the user for diagnostic purposes, e.g. by applying <code>plot</code> against each object.

Author(s)

Alireza S. Mahani, Mansour T.A. Sharabiani

Examples

```
library("survival")
data(ovarian)
folds <- bsgw.generate.folds.eventbalanced(Surv(futime, fustat) ~ 1, ovarian, 5)
cv <- bsgw.crossval(ovarian, folds, formula=Surv(futime, fustat) ~ ecog.ps + rx
, control=bsgw.control(iter=50, nskip=10), print.level = 3)
```

```
cv2 <- bsgw.crossval.wrapper(ovarian, folds, formula=Surv(futime, fustat) ~ ecog.ps + rx
, control=bsgw.control(iter=50, nskip=10)
, print.level=3, lambda.vec=exp(seq(from=log(0.1), to=log(1), length.out = 3)))
```

plot.bsgw

Plot diagnostics for a bsgw object

Description

Four sets of MCMC diagnostic plots are currently generated: 1) log-likelihood and log-posterior (including shrinkage effect) as a function of iteration number, 2) coefficient trace plots, 3) coefficient autocorrelation plots, 4) coefficient histograms.

Usage

```
## S3 method for class 'bsgw'
plot(x, pval=0.05, burnin=round(x$control$iter/2), nrow=2, ncol=3, ...)
```

Arguments

x	A bsgw object, typically the output of bsgw function.
pval	The P-value at which lower/upper bounds on coefficients are calculated and overlaid on trace plots and histograms.
burnin	Number of samples discarded from the beginning of an MCMC chain, after which parameter quantiles are calculated.
nrow	Number of rows of subplots within each figure, applied to plot sets 2-4.
ncol	Number of columns of subplots within each figure, applied to plot sets 2-4.
...	Further arguments to be passed to/from other methods.

Author(s)

Alireza S. Mahani, Mansour T.A. Sharabiani

Examples

```
library("survival")
data(ovarian)
est <- bsgw(Surv(futime, fustat) ~ ecog.ps + rx, ovarian
, control=bsgw.control(iter=400, nskip=100))
plot(est)
```

predict.bsgw	<i>Predict method for bsgw model fits</i>
--------------	---

Description

Calculates log-likelihood and hazard/cumulative hazard/survival functions over a user-supplied vector time values, based on BSGW model object.

Usage

```
## S3 method for class 'bsgw'
predict(object, newdata=NULL, tvec=NULL, burnin=object$control$burnin, ncores=1, ...)
## S3 method for class 'predict.bsgw'
summary(object, idx=1:length(object$median$survreg.scale), burnin=object$burnin, pval=0.05
, popmean=identical(idx,1:length(object$median$survreg.scale)), make.plot=TRUE, ...)
```

Arguments

object	For predict.bsgw, an object of class "bsgw", usually the result of a call to bsgw ; for summary.predict.bsgw, an object of class "predict.bsgw", usually the result of a call to predict.bsgw.
newdata	An optional data frame in which to look for variables with which to predict. If omitted, the fitted values (training set) are used.
tvec	An optional vector of time values, along which time-dependent entities (hazard, cumulative hazard, survival) will be predicted. If omitted, only the time-independent entities (currently only log-likelihood) will be calculated. If a single integer is provided for tvec, it is interpreted as number of time points, equally spaced from 0 to object\$tmax: <code>tvec <- seq(from=0.0, to=object\$tmax, length.out=tvec)</code> .
burnin	Number of samples to discard from the beginning of each MCMC chain before calculating median value(s) for time-independent entities.
ncores	Number of cores to use for parallel prediction.
...	Further arguments to be passed to/from other methods.
idx	Index of observations (rows of newdata or training data) for which to generate summary statistics. Default is the entire data.
pval	Desired p-value, based on which lower/upper bounds will be calculated. Default is 0.05.
popmean	Whether population averages must be calculated or not. By default, population averages are only calculated when the entire data is included in prediction.
make.plot	Whether population mean and other plots must be created or not.

Details

The time-dependent predicted objects (except `loglike`) are three-dimensional arrays of size (`nsmp` x `nt` x `nobs`), where `nsmp` = number of MCMC samples, `nt` = number of time values in `tvec`, and `nobs` = number of rows in `newdata`. Therefore, even for modest data sizes, these objects can occupy large chunks of memory. For example, for `nsmp=1000`, `nt=100`, `nobs=1000`, the three objects `h`, `H`, `S` have a total size of 2.2GB. Since applying `quantile` to these arrays is time-consuming (as needed for calculation of median and lower/upper bounds), we have left such summaries out of the scope of `predict` function. Users can instead apply `summary` to the prediction object to obtain summary statistics. During cross-validation-based selection of shrinkage parameter `lambda`, there is no need to supply `tvec` since we only the log-likelihood value. This significantly speeds up the parameter-tuning process. The function `summary.predict.bsgw` allows the user to calculate summary statistics for a subset (or all of) data, if desired. This approach is in line with the overall philosophy of delaying the data summarization until necessary, to avoid unnecessary loss in accuracy due to premature blending of information contained in individual samples.

Value

The function `predict.bsgw` returns as object of class "predict.bsgw" with the following fields:

<code>tvec</code>	Actual vector of time values (if any) used for prediction.
<code>burnin</code>	Same as input.
<code>median</code>	List of median values for predicted entities. Currently, only <code>loglike</code> and <code>survreg.scale</code> median is produced. See 'Details' for explanation.
<code>smp</code>	List of MCMC samples for predicted entities. Elements include <code>h</code> (hazard function), <code>H</code> (cumulative hazard function), <code>S</code> (survival function), <code>survreg.scale</code> (inverse of shape parameter in <code>rweibull</code>), and <code>loglike</code> (model log-likelihood). All functions are evaluated over time values specified in <code>tvec</code> .
<code>km.fit</code>	Kaplan-Meyer fit of the data used for prediction (if data contains response fields).

The function `summary.predict.bsgw` returns a list with the following fields:

<code>lower</code>	A list of lower-bound values for <code>h</code> , <code>H</code> , <code>S</code> , <code>hr</code> (hazard ratio of <code>idx[2]</code> to <code>idx[1]</code> observation), and <code>S.diff</code> (survival probability of <code>idx[2]</code> minus <code>idx[1]</code>). The last two are only included if <code>length(idx)==2</code> .
<code>median</code>	List of median values for same entities described in <code>lower</code> .
<code>upper</code>	List of upper-bound values for same entities described in <code>lower</code> .
<code>popmean</code>	Lower-bound/median/upper-bound values for population average of survival probability.
<code>km.fit</code>	Kaplan-Meyer fit associated with the prediction object (if available).

Author(s)

Alireza S. Mahani, Mansour T.A. Sharabiani

Examples

```
library("survival")
data(ovarian)
est <- bsgw(Surv(futime, fustat) ~ ecog.ps + rx, ovarian
            , control=bsgw.control(iter=400, nskip=100))
pred <- predict(est, tvec=100)
predsumm <- summary(pred, idx=1:10)
```

summary.bsgw	<i>Summarizing Bayesian Survival Generalized Weibull (BSGW) model fits</i>
--------------	--

Description

summary method for class "bsgw".

Usage

```
## S3 method for class 'bsgw'
summary(object, pval = 0.05, burnin = object$control$burnin, ...)
## S3 method for class 'summary.bsgw'
print(x, ...)
```

Arguments

object	An object of class "bsgw", usually the result of a call to <code>bsgw</code> .
x	An object of class "summary.bsgw", usually the result of a call to <code>summary.bsgw</code> .
pval	Desired p-value, based on which lower/upper bounds will be calculated. Default is 0.05.
burnin	Number of samples to discard from the beginning of each MCMC chain before calculating median and lower/upper bounds.
...	Further arguments to be passed to/from other methods.

Value

The function `summary.bsgw` calculates median as well as lower/upper bounds for all model coefficients, given the supplied p-value. It also calculates the p-value for coefficients being significant smaller/larger than zero. It contains returns an object of class "summary.bsgw" with the following elements:

call	The matched call.
pval	Same as input.
burnin	Same as input.
coefficients	A $p \times 4$ matrix with columns for the estimated coefficient median, its lower and upper bounds given the user-supplied p-value, and the p-value for being smaller/larger than zero.
survreg.scale	List of lower, median, and upper values of the survreg-style scale parameter (i.e. inverse of shape parameter in <code>rweibull</code>) for the training-set population.

Author(s)

Alireza S. Mahani, Mansour T.A. Sharabiani

See Also

See [summary](#) for a description of the generic method.

The model fitting function is [bsgw](#).

Examples

```
library("survival")
data(ovarian)
est <- bsgw(Surv(futime, fustat) ~ ecog.ps + rx, ovarian
           , control=bsgw.control(iter=400, nskip=100))
summary(est, pval=0.1)
```

Index

BSGW (bsgw), [2](#)
bsgw, [2](#), [7–9](#), [11](#), [12](#)
bsgw.control, [7](#)
bsgw.crossval, [6](#)
bsgw.generate.folds (bsgw.crossval), [6](#)

plot.bsgw, [5](#), [8](#)
predict.bsgw, [5](#), [9](#)
print.bsgw (bsgw), [2](#)
print.summary.bsgw (summary.bsgw), [11](#)

summary, [12](#)
summary.bsgw, [5](#), [11](#)
summary.predict.bsgw (predict.bsgw), [9](#)