

Package ‘SuperLearner’

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Author Eric Polley, Mark van der Laan

Maintainer Eric Polley <eric.polley@nih.gov>

Description This package implements the super learner prediction method and contains a library of prediction algorithms to be used in the super learner.

License GPL-3

URL <https://github.com/ecpolley/SuperLearner>

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CV.SuperLearner	<i>Function to get V-fold cross-validated risk estimate for super learner</i>
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Description

Function to get V-fold cross-validated risk estimate for super learner. This function simply splits the data into V folds and then calls SuperLearner. Most of the arguments are passed directly to SuperLearner.

Usage

```
CV.SuperLearner(Y, X, V = 20, family = gaussian(), SL.library,
  method = "method.NNLS", id = NULL, verbose = FALSE,
  control = list(saveFitLibrary = FALSE), cvControl = list(),
  obsWeights = NULL, saveAll = TRUE, parallel = "seq")
```

Arguments

Y	The outcome.
X	The covariates.
V	The number of folds for CV.SuperLearner. This is not the number of folds for SuperLearner. The number of folds for SuperLearner is controlled with cvControl.
family	Currently allows gaussian or binomial to describe the error distribution. Link function information will be ignored and should be contained in the method argument below.
SL.library	Either a character vector of prediction algorithms or a list containing character vectors. See details below for examples on the structure. A list of functions included in the SuperLearner package can be found with listWrappers().
method	A list (or a function to create a list) containing details on estimating the coefficients for the super learner and the model to combine the individual algorithms in the library. See ?method.template for details. Currently, the built in options are either "method.NNLS" (the default), "method.NNLS2", "method.NNloglik",

"method.CC_LS", "method.CC_nloglik", or "method.AUC". NNLS and NNLS2 are non-negative least squares based on the Lawson-Hanson algorithm and the dual method of Goldfarb and Idnani, respectively. NNLS and NNLS2 will work for both gaussian and binomial outcomes. NNloglik is a non-negative binomial likelihood maximization using the BFGS quasi-Newton optimization method. NN* methods are normalized so weights sum to one. CC_LS uses Goldfarb and Idnani's quadratic programming algorithm to calculate the best convex combination of weights to minimize the squared error loss. CC_nloglik calculates the convex combination of weights that minimize the negative binomial log likelihood on the logistic scale using the sequential quadratic programming algorithm. AUC, which only works for binary outcomes, uses the Nelder-Mead method via the optim function to minimize rank loss (equivalent to maximizing AUC).

id	Optional cluster identification variable. For the cross-validation splits, id forces observations in the same cluster to be in the same validation fold. id is passed to the prediction and screening algorithms in SL.library, but be sure to check the individual wrappers as many of them ignore the information.
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging).
control	A list of parameters to control the estimation process. Parameters include saveFitLibrary and trimLogit. See SuperLearner.control for details.
cvControl	A list of parameters to control the cross-validation process. Parameters include V, stratifyCV, shuffle and validRows. See SuperLearner.CV.control for details.
obsWeights	Optional observation weights variable. As with id above, obsWeights is passed to the prediction and screening algorithms, but many of the built in wrappers ignore (or can't use) the information. If you are using observation weights, make sure the library you specify uses the information.
saveAll	Logical; Should the entire SuperLearner object be saved for each fold?
parallel	Options for parallel computation of the V-fold step. Use "seq" (the default) for sequential computation. parallel = 'multicore' to use mclapply for the V-fold step (but note that SuperLearner() will still be sequential). Or parallel can be the name of a snow cluster and will use parLapply for the V-fold step. For both multicore and snow, the inner SuperLearner calls will be sequential.

Details

The SuperLearner function builds a estimator, but does not contain an estimate on the performance of the estimator. Various methods exist for estimator performance evaluation. If you are familiar with the super learner algorithm, it should be no surprise we recommend using cross-validation to evaluate the honest performance of the super learner estimator. The function CV.SuperLearner computes the usual V-fold cross-validated risk estimate for the super learner (and all algorithms in SL.library for comparison).

Value

An object of class CV.SuperLearner (a list) with components:

<code>call</code>	The matched call.
<code>AllSL</code>	If <code>saveAll = TRUE</code> , a list with output from each call to <code>SuperLearner</code> , otherwise <code>NULL</code> .
<code>SL.predict</code>	The predicted values from the super learner when each particular row was part of the validation fold.
<code>discreteSL.predict</code>	The traditional cross-validated selector. Picks the algorithm with the smallest cross-validated risk (in super learner terms, gives that algorithm coefficient 1 and all others 0).
<code>whichDiscreteSL</code>	A list of length <code>V</code> . The elements in the list are the algorithm that had the smallest cross-validated risk estimate for that fold.
<code>library.predict</code>	A matrix with the predicted values from each algorithm in <code>SL.library</code> . The columns are the algorithms in <code>SL.library</code> and the rows represent the predicted values when that particular row was in the validation fold (i.e. not used to fit that estimator).
<code>coef</code>	A matrix with the coefficients for the super learner on each fold. The columns are the algorithms in <code>SL.library</code> the rows are the folds.
<code>folds</code>	A list containing the row numbers for each validation fold.
<code>V</code>	Number of folds for <code>CV.SuperLearner</code> .
<code>libraryNames</code>	A character vector with the names of the algorithms in the library. The format is 'predictionAlgorithm_screeningAlgorithm' with '_All' used to denote the prediction algorithm run on all variables in <code>X</code> .
<code>SL.library</code>	Returns <code>SL.library</code> in the same format as the argument with the same name above.
<code>method</code>	A list with the method functions.
<code>Y</code>	The outcome

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[SuperLearner](#)

Examples

```
set.seed(23432)
## training set
n <- 500
p <- 50
X <- matrix(rnorm(n*p), nrow = n, ncol = p)
colnames(X) <- paste("X", 1:p, sep="")
X <- data.frame(X)
Y <- X[, 1] + sqrt(abs(X[, 2] * X[, 3])) + X[, 2] - X[, 3] + rnorm(n)
```

```
# build Library and run Super Learner
SL.library <- c("SL.glm", "SL.randomForest", "SL.gam", "SL.polymars", "SL.mean")
## Not run:
test <- CV.SuperLearner(Y = Y, X = X, V = 10, SL.library = SL.library,
  verbose = TRUE, method = "method.NNLS")
test
summary(test)
# Look at the coefficients across folds
coef(test)

## End(Not run)
```

CVFolds

Generate list of row numbers for each fold in the cross-validation

Description

Generate list of row numbers for each fold in the cross-validation. CVFolds is used in the SuperLearner to create the cross-validation splits.

Usage

```
CVFolds(N, id, Y, cvControl)
```

Arguments

N	Sample size
id	Optional cluster id variable. If present, all observations in the same cluster will always be in the same split.
Y	outcome
cvControl	Control parameters for the cross-validation step. See SuperLearner.CV.control for details.

Value

validRows	A list of length V where each element in the list is a vector with the row numbers of the corresponding validation sample.
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Author(s)

Eric C Polley <eric.polley@nih.gov>

listWrappers	<i>list all wrapper functions in SuperLearner</i>
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Description

List all wrapper functions in [SuperLearner](#) package

Usage

```
listWrappers(what = "both")
```

Arguments

what	What list to return. Can be both for both prediction algorithms and screening algorithms, SL for the prediction algorithms, screen for the screening algorithms, method for the estimation method details, or anything else will return a list of all (exported) functions in the SuperLearner package. Additional wrapper functions are available at https://github.com/ecpolley/SuperLearnerExtra .
------	---

Value

Invisible character vector with all exported functions in the SuperLearner package

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[SuperLearner](#)

Examples

```
listWrappers(what = "SL")
listWrappers(what = "screen")
```

plot.CV.SuperLearner	<i>Graphical display of the V-fold CV risk estimates</i>
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Description

The function plots the V-fold cross-validated risk estimates for the super learner, the discrete super learner and each algorithm in the library. By default the estimates will be sorted and include an asymptotic 95% confidence interval.

Usage

```
## S3 method for class 'CV.SuperLearner'  
plot(x, package = "ggplot2", constant = qnorm(0.975), sort = TRUE, ...)
```

Arguments

x	The output from CV.SuperLearner.
package	Either "ggplot2" or "lattice". The package selected must be available.
constant	A numeric value. The confidence interval is defined as $p \pm \text{constant} * \text{se}$, where p is the point estimate and se is the standard error. The default is the quantile of the standard normal corresponding to a 95% CI.
sort	Logical. Should the rows in the plot be sorted from the smallest to the largest point estimate. If FALSE, then the order is super learner, discrete super learner, then the estimators in SL.library.
...	Additional arguments for summary.CV.SuperLearner

Details

see [summary.CV.SuperLearner](#) for details on how the estimates are computed

Value

Returns the plot (either a ggplot2 object (class ggplot) or a lattice object (class trellis))

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[summary.CV.SuperLearner](#) and [CV.SuperLearner](#)

predict.SuperLearner *Predict method for SuperLearner object*

Description

Obtains predictions on a new data set from a SuperLearner fit. May require the original data if one of the library algorithms uses the original data in its predict method.

Usage

```
## S3 method for class 'SuperLearner'  
predict(object, newdata, X = NULL, Y = NULL, onlySL = FALSE, ...)
```

Arguments

object	Fitted object from SuperLearner
newdata	New X values for prediction
X	Original data set used to fit object
Y	Original outcome used to fit object
onlySL	Logical. If TRUE, only compute predictions for algorithms with non-zero coefficients in the super learner object. Default is FALSE (computes predictions for all algorithms in library).
...	Additional arguments passed to the <code>predict.SL.*</code> functions

Details

If `newdata` is omitted the predicted values from `object` are returned. Each algorithm in the Super Learner library needs to have a corresponding prediction function with the “predict.” prefixed onto the algorithm name (e.g. `predict.SL.glm` for `SL.glm`).

Value

<code>pred</code>	Predicted values from Super Learner fit
<code>library.predict</code>	Predicted values for each algorithm in library

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[SuperLearner](#)

SampleSplitSuperLearner

Super Learner Prediction Function

Description

A Prediction Function for the Super Learner. The `SuperLearner` function takes a training set pair (X,Y) and returns the predicted values based on a validation set. `SampleSplitSuperLearner` uses sample split validation whereas `SuperLearner` uses V-fold cross-validation.

Usage

```
SampleSplitSuperLearner(Y, X, newX = NULL, family = gaussian(), SL.library,
  method = "method.NNLS", id = NULL, verbose = FALSE,
  control = list(), split = 0.8, obsWeights = NULL)
```

Arguments

Y	The outcome in the training data set. Must be a numeric vector.
X	The predictor variables in the training data set, usually a data.frame.
newX	The predictor variables in the validation data set. The structure should match X. If missing, uses X for newX.
SL.library	Either a character vector of prediction algorithms or a list containing character vectors. See details below for examples on the structure. A list of functions included in the SuperLearner package can be found with <code>listWrappers()</code> .
verbose	logical; TRUE for printing progress during the computation (helpful for debugging).
family	Currently allows gaussian or binomial to describe the error distribution. Link function information will be ignored and should be contained in the method argument below.
method	A list (or a function to create a list) containing details on estimating the coefficients for the super learner and the model to combine the individual algorithms in the library. See <code>?method.template</code> for details. Currently, the built in options are either "method.NNLS" (the default), "method.NNLS2", "method.NNloglik", "method.CC_LS", or "method.CC_nloglik". NNLS and NNLS2 are non-negative least squares based on the Lawson-Hanson algorithm and the dual method of Goldfarb and Idnani, respectively. NNLS and NNLS2 will work for both gaussian and binomial outcomes. NNloglik is a non-negative binomial likelihood maximization using the BFGS quasi-Newton optimization method. NN* methods are normalized so weights sum to one. CC_LS uses Goldfarb and Idnani's quadratic programming algorithm to calculate the best convex combination of weights to minimize the squared error loss. CC_nloglik calculates the convex combination of weights that minimize the negative binomial log likelihood on the logistic scale using the sequential quadratic programming algorithm.
id	Optional cluster identification variable. For the cross-validation splits, id forces observations in the same cluster to be in the same validation fold. id is passed to the prediction and screening algorithms in SL.library, but be sure to check the individual wrappers as many of them ignore the information.
obsWeights	Optional observation weights variable. As with id above, obsWeights is passed to the prediction and screening algorithms, but many of the built in wrappers ignore (or can't use) the information. If you are using observation weights, make sure the library you specify uses the information.
control	A list of parameters to control the estimation process. Parameters include <code>saveFitLibrary</code> and <code>trimLogit</code> . See SuperLearner.control for details.
split	Either a single value between 0 and 1 indicating the fraction of the samples for the training split. A value of 0.8 will randomly assign 80 percent of the samples to the training split and the other 20 percent to the validation split. Alternatively, split can be a numeric vector with the row numbers of X corresponding to the validation split. All other rows not in the vector will be considered in the training split.

Details

SuperLearner fits the super learner prediction algorithm. The weights for each algorithm in `SL.library` is estimated, along with the fit of each algorithm.

The prescreen algorithms. These algorithms first rank the variables in X based on either a univariate regression p-value of the `randomForest` variable importance. A subset of the variables in X is selected based on a pre-defined cut-off. With this subset of the X variables, the algorithms in `SL.library` are then fit.

The SuperLearner package contains a few prediction and screening algorithm wrappers. The full list of wrappers can be viewed with `listWrappers()`. The design of the SuperLearner package is such that the user can easily add their own wrappers. We also maintain a website with additional examples of wrapper functions at <https://github.com/ecpolley/SuperLearnerExtra>.

Value

<code>call</code>	The matched call.
<code>libraryNames</code>	A character vector with the names of the algorithms in the library. The format is 'predictionAlgorithm_screeningAlgorithm' with '_All' used to denote the prediction algorithm run on all variables in X .
<code>SL.library</code>	Returns <code>SL.library</code> in the same format as the argument with the same name above.
<code>SL.predict</code>	The predicted values from the super learner for the rows in <code>newX</code> .
<code>coef</code>	Coefficients for the super learner.
<code>library.predict</code>	A matrix with the predicted values from each algorithm in <code>SL.library</code> for the rows in <code>newX</code> .
<code>Z</code>	The Z matrix (the cross-validated predicted values for each algorithm in <code>SL.library</code>).
<code>cvRisk</code>	A numeric vector with the V-fold cross-validated risk estimate for each algorithm in <code>SL.library</code> . Note that this does not contain the CV risk estimate for the SuperLearner, only the individual algorithms in the library.
<code>family</code>	Returns the family value from above
<code>fitLibrary</code>	A list with the fitted objects for each algorithm in <code>SL.library</code> on the full training data set.
<code>varNames</code>	A character vector with the names of the variables in X .
<code>validRows</code>	A list containing the row numbers for the V-fold cross-validation step.
<code>method</code>	A list with the method functions.
<code>whichScreen</code>	A logical matrix indicating which variables passed each screening algorithm.
<code>control</code>	The control list.
<code>split</code>	The split value.
<code>errorsInCVLibrary</code>	A logical vector indicating if any algorithms experienced an error within the CV step.
<code>errorsInLibrary</code>	A logical vector indicating if any algorithms experienced an error on the full data.

Author(s)

Eric C Polley <eric.polley@nih.gov>

References

van der Laan, M. J., Polley, E. C. and Hubbard, A. E. (2008) Super Learner, *Statistical Applications of Genetics and Molecular Biology*, **6**, article 25. <http://www.bepress.com/sagmb/vol6/iss1/art25>

Examples

```
## Not run:
## simulate data
set.seed(23432)
## training set
n <- 500
p <- 50
X <- matrix(rnorm(n*p), nrow = n, ncol = p)
colnames(X) <- paste("X", 1:p, sep="")
X <- data.frame(X)
Y <- X[, 1] + sqrt(abs(X[, 2] * X[, 3])) + X[, 2] - X[, 3] + rnorm(n)

## test set
m <- 1000
newX <- matrix(rnorm(m*p), nrow = m, ncol = p)
colnames(newX) <- paste("X", 1:p, sep="")
newX <- data.frame(newX)
newY <- newX[, 1] + sqrt(abs(newX[, 2] * newX[, 3])) + newX[, 2] -
  newX[, 3] + rnorm(m)

# generate Library and run Super Learner
SL.library <- c("SL.glm", "SL.randomForest", "SL.gam",
  "SL.polymars", "SL.mean")
test <- SampleSplitSuperLearner(Y = Y, X = X, newX = newX, SL.library = SL.library,
  verbose = TRUE, method = "method.NNLS")
test

# library with screening
SL.library <- list(c("SL.glmnet", "All"), c("SL.glm", "screen.randomForest",
  "All", "screen.SIS"), "SL.randomForest", c("SL.polymars", "All"), "SL.mean")
test <- SuperLearner(Y = Y, X = X, newX = newX, SL.library = SL.library,
  verbose = TRUE, method = "method.NNLS")
test

# binary outcome
set.seed(1)
N <- 200
X <- matrix(rnorm(N*10), N, 10)
X <- as.data.frame(X)
Y <- rbinom(N, 1, plogis(.2*X[, 1] + .1*X[, 2] - .2*X[, 3] +
  .1*X[, 3]*X[, 4] - .2*abs(X[, 4])))
```

```

SL.library <- c("SL.glmnet", "SL.glm", "SL.knn", "SL.gam", "SL.mean")

# least squares loss function
test.NNLS <- SampleSplitSuperLearner(Y = Y, X = X, SL.library = SL.library,
  verbose = TRUE, method = "method.NNLS", family = binomial())
test.NNLS

## End(Not run)

```

```
summary.CV.SuperLearner
```

Summary Function for Cross-Validated Super Learner

Description

summary method for the CV.SuperLearner function

Usage

```

## S3 method for class 'CV.SuperLearner'
summary(object, obsWeights = NULL, ...)

## S3 method for class 'summary.CV.SuperLearner'
print(x, digits, ...)

```

Arguments

object	An object of class "CV.SuperLearner", the result of a call to CV.SuperLearner.
x	An object of class "summary.CV.SuperLearner", the result of a call to summary.CV.SuperLearner.
obsWeights	Optional vector for observation weights.
digits	The number of significant digits to use when printing.
...	additional arguments ...

Details

Summary method for CV.SuperLearner. Calculates the V-fold cross-validated estimate of either the mean squared error or the $-2 \cdot \log(L)$ depending on the loss function used.

Value

summary.CV.SuperLearner returns a list with components

call	The function call from CV.SuperLearner
method	Describes the loss function used. Currently either least squares or negative log Likelihood.
V	Number of folds

Risk.SL	Risk estimate for the super learner
Risk.dSL	Risk estimate for the discrete super learner (the cross-validation selector)
Risk.library	A matrix with the risk estimates for each algorithm in the library
Table	A table with the mean risk estimate and standard deviation across the folds for the super learner and all algorithms in the library

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[CV.SuperLearner](#)

SuperLearner

Super Learner Prediction Function

Description

A Prediction Function for the Super Learner. The SuperLearner function takes a training set pair (X,Y) and returns the predicted values based on a validation set.

Usage

```
SuperLearner(Y, X, newX = NULL, family = gaussian(), SL.library,
  method = "method.NNLS", id = NULL, verbose = FALSE,
  control = list(), cvControl = list(), obsWeights = NULL)
```

Arguments

Y	The outcome in the training data set. Must be a numeric vector.
X	The predictor variables in the training data set, usually a data.frame.
newX	The predictor variables in the validation data set. The structure should match X. If missing, uses X for newX.
SL.library	Either a character vector of prediction algorithms or a list containing character vectors. See details below for examples on the structure. A list of functions included in the SuperLearner package can be found with listWrappers().
verbose	logical; TRUE for printing progress during the computation (helpful for debugging).
family	Currently allows gaussian or binomial to describe the error distribution. Link function information will be ignored and should be contained in the method argument below.

method	A list (or a function to create a list) containing details on estimating the coefficients for the super learner and the model to combine the individual algorithms in the library. See <code>?method.template</code> for details. Currently, the built in options are either "method.NNLS" (the default), "method.NNLS2", "method.NNloglik", "method.CC_LS", "method.CC_nloglik", or "method.AUC". NNLS and NNLS2 are non-negative least squares based on the Lawson-Hanson algorithm and the dual method of Goldfarb and Idnani, respectively. NNLS and NNLS2 will work for both gaussian and binomial outcomes. NNloglik is a non-negative binomial likelihood maximization using the BFGS quasi-Newton optimization method. NN* methods are normalized so weights sum to one. CC_LS uses Goldfarb and Idnani's quadratic programming algorithm to calculate the best convex combination of weights to minimize the squared error loss. CC_nloglik calculates the convex combination of weights that minimize the negative binomial log likelihood on the logistic scale using the sequential quadratic programming algorithm. AUC, which only works for binary outcomes, uses the Nelder-Mead method via the <code>optim</code> function to minimize rank loss (equivalent to maximizing AUC).
id	Optional cluster identification variable. For the cross-validation splits, <code>id</code> forces observations in the same cluster to be in the same validation fold. <code>id</code> is passed to the prediction and screening algorithms in <code>SL.library</code> , but be sure to check the individual wrappers as many of them ignore the information.
obsWeights	Optional observation weights variable. As with <code>id</code> above, <code>obsWeights</code> is passed to the prediction and screening algorithms, but many of the built in wrappers ignore (or can't use) the information. If you are using observation weights, make sure the library you specify uses the information.
control	A list of parameters to control the estimation process. Parameters include <code>saveFitLibrary</code> and <code>trimLogit</code> . See SuperLearner.control for details.
cvControl	A list of parameters to control the cross-validation process. Parameters include <code>V</code> , <code>stratifyCV</code> , <code>shuffle</code> and <code>validRows</code> . See SuperLearner.CV.control for details.

Details

`SuperLearner` fits the super learner prediction algorithm. The weights for each algorithm in `SL.library` is estimated, along with the fit of each algorithm.

The prescreen algorithms. These algorithms first rank the variables in X based on either a univariate regression p-value of the `randomForest` variable importance. A subset of the variables in X is selected based on a pre-defined cut-off. With this subset of the X variables, the algorithms in `SL.library` are then fit.

The `SuperLearner` package contains a few prediction and screening algorithm wrappers. The full list of wrappers can be viewed with `listWrappers()`. The design of the `SuperLearner` package is such that the user can easily add their own wrappers. We also maintain a website with additional examples of wrapper functions at <https://github.com/ecpolley/SuperLearnerExtra>.

Value

`call` The matched call.

libraryNames	A character vector with the names of the algorithms in the library. The format is 'predictionAlgorithm_screeningAlgorithm' with '_All' used to denote the prediction algorithm run on all variables in X.
SL.library	Returns SL.library in the same format as the argument with the same name above.
SL.predict	The predicted values from the super learner for the rows in newX.
coef	Coefficients for the super learner.
library.predict	A matrix with the predicted values from each algorithm in SL.library for the rows in newX.
Z	The Z matrix (the cross-validated predicted values for each algorithm in SL.library).
cvRisk	A numeric vector with the V-fold cross-validated risk estimate for each algorithm in SL.library. Note that this does not contain the CV risk estimate for the SuperLearner, only the individual algorithms in the library.
family	Returns the family value from above
fitLibrary	A list with the fitted objects for each algorithm in SL.library on the full training data set.
varNames	A character vector with the names of the variables in X.
validRows	A list containing the row numbers for the V-fold cross-validation step.
method	A list with the method functions.
whichScreen	A logical matrix indicating which variables passed each screening algorithm.
control	The control list.
cvControl	The cvControl list.
errorsInCVLibrary	A logical vector indicating if any algorithms experienced an error within the CV step.
errorsInLibrary	A logical vector indicating if any algorithms experienced an error on the full data.

Author(s)

Eric C Polley <eric.polley@nih.gov>

References

van der Laan, M. J., Polley, E. C. and Hubbard, A. E. (2008) Super Learner, *Statistical Applications of Genetics and Molecular Biology*, **6**, article 25. <http://www.bepress.com/sagmb/vol6/iss1/art25>

Examples

```

## Not run:
## simulate data
set.seed(23432)
## training set
n <- 500
p <- 50
X <- matrix(rnorm(n*p), nrow = n, ncol = p)
colnames(X) <- paste("X", 1:p, sep="")
X <- data.frame(X)
Y <- X[, 1] + sqrt(abs(X[, 2] * X[, 3])) + X[, 2] - X[, 3] + rnorm(n)

## test set
m <- 1000
newX <- matrix(rnorm(m*p), nrow = m, ncol = p)
colnames(newX) <- paste("X", 1:p, sep="")
newX <- data.frame(newX)
newY <- newX[, 1] + sqrt(abs(newX[, 2] * newX[, 3])) + newX[, 2] -
  newX[, 3] + rnorm(m)

# generate Library and run Super Learner
SL.library <- c("SL.glm", "SL.randomForest", "SL.gam",
  "SL.polymars", "SL.mean")
test <- SuperLearner(Y = Y, X = X, newX = newX, SL.library = SL.library,
  verbose = TRUE, method = "method.NNLS")
test

# library with screening
SL.library <- list(c("SL.glmnet", "All"), c("SL.glm", "screen.randomForest",
  "All", "screen.SIS"), "SL.randomForest", c("SL.polymars", "All"), "SL.mean")
test <- SuperLearner(Y = Y, X = X, newX = newX, SL.library = SL.library,
  verbose = TRUE, method = "method.NNLS")
test

# binary outcome
set.seed(1)
N <- 200
X <- matrix(rnorm(N*10), N, 10)
X <- as.data.frame(X)
Y <- rbinom(N, 1, plogis(.2*X[, 1] + .1*X[, 2] - .2*X[, 3] +
  .1*X[, 3]*X[, 4] - .2*abs(X[, 4])))

SL.library <- c("SL.glmnet", "SL.glm", "SL.knn", "SL.gam", "SL.mean")

# least squares loss function
test.NNLS <- SuperLearner(Y = Y, X = X, SL.library = SL.library,
  verbose = TRUE, method = "method.NNLS", family = binomial())
test.NNLS

# negative log binomial likelihood loss function
test.NNloglik <- SuperLearner(Y = Y, X = X, SL.library = SL.library,
  verbose = TRUE, method = "method.NNloglik", family = binomial())

```

```

test.NNloglik

# 1 - AUC loss function
test.AUC <- SuperLearner(Y = Y, X = X, SL.library = SL.library,
  verbose = TRUE, method = "method.AUC", family = binomial())
test.AUC

# 2
# adapted from library(SIS)
set.seed(1)
# training
b <- c(2, 2, 2, -3*sqrt(2))
n <- 150
p <- 200
truerho <- 0.5
corrmat <- diag(rep(1-truerho, p)) + matrix(truerho, p, p)
corrmat[, 4] = sqrt(truerho)
corrmat[4, ] = sqrt(truerho)
corrmat[4, 4] = 1
cholmat <- chol(corrmat)
x <- matrix(rnorm(n*p, mean=0, sd=1), n, p)
x <- x
feta <- x[, 1:4]
fprob <- exp(feta) / (1 + exp(feta))
y <- rbinom(n, 1, fprob)

# test
m <- 10000
newx <- matrix(rnorm(m*p, mean=0, sd=1), m, p)
newx <- newx
newfeta <- newx[, 1:4]
newfprob <- exp(newfeta) / (1 + exp(newfeta))
newy <- rbinom(m, 1, newfprob)

DATA2 <- data.frame(Y = y, X = x)
newDATA2 <- data.frame(Y = newy, X=newx)

create.SL.knn <- function(k = c(20, 30)) {
  for(mm in seq(length(k))){
    eval(parse(text = paste('SL.knn.', k[mm], '<- function(..., k = ', k[mm],
      ') SL.knn(..., k = k)', sep = '')), envir = .GlobalEnv)
  }
  invisible(TRUE)
}
create.SL.knn(c(20, 30, 40, 50, 60, 70))

# library with screening
SL.library <- list(c("SL.glmnet", "All"), c("SL.glm", "screen.randomForest"),
  "SL.randomForest", "SL.knn", "SL.knn.20", "SL.knn.30", "SL.knn.40",
  "SL.knn.50", "SL.knn.60", "SL.knn.70",
  c("SL.polymars", "screen.randomForest"))
test <- SuperLearner(Y = DATA2$Y, X = DATA2[, -1], newX = newDATA2[, -1],
  SL.library = SL.library, verbose = TRUE, family = binomial())

```

```

test

## examples with multicore
set.seed(23432)
## training set
n <- 500
p <- 50
X <- matrix(rnorm(n*p), nrow = n, ncol = p)
colnames(X) <- paste("X", 1:p, sep="")
X <- data.frame(X)
Y <- X[, 1] + sqrt(abs(X[, 2] * X[, 3])) + X[, 2] - X[, 3] + rnorm(n)

## test set
m <- 1000
newX <- matrix(rnorm(m*p), nrow = m, ncol = p)
colnames(newX) <- paste("X", 1:p, sep="")
newX <- data.frame(newX)
newY <- newX[, 1] + sqrt(abs(newX[, 2] * newX[, 3])) + newX[, 2] - newX[, 3] + rnorm(m)

# generate Library and run Super Learner
SL.library <- c("SL.glm", "SL.randomForest", "SL.gam",
               "SL.polymars", "SL.mean")

testMC <- mcSuperLearner(Y = Y, X = X, newX = newX, SL.library = SL.library,
                        method = "method.NNLS")
testMC

## examples with snow
library(parallel)
cl <- makeCluster(2, type = "PSOCK") # can use different types here
clusterSetRNGStream(cl, iseed = 2343)
testSNOW <- snowSuperLearner(cluster = cl, Y = Y, X = X, newX = newX,
                             SL.library = SL.library, method = "method.NNLS")
testSNOW
stopCluster(cl)

## snow example with user-generated wrappers
# If you write your own wrappers and are using snowSuperLearner()
# These new wrappers need to be added to the SuperLearner namespace and exported to the clusters
# Using a simple example here, but can define any new SuperLearner wrapper
my.SL.wrapper <- function(...) SL.glm(...)
# assign function into SuperLearner namespace
environment(my.SL.wrapper) <- asNamespace("SuperLearner")

cl <- makeCluster(2, type = "PSOCK") # can use different types here
clusterSetRNGStream(cl, iseed = 2343)
clusterExport(cl, c("my.SL.wrapper")) # copy the function to all clusters
testSNOW <- snowSuperLearner(cluster = cl, Y = Y, X = X, newX = newX,
                             SL.library = c("SL.glm", "SL.mean", "my.SL.wrapper"), method = "method.NNLS")
testSNOW
stopCluster(cl)

## timing

```

```

replicate(5, system.time(SuperLearner(Y = Y, X = X, newX = newX,
  SL.library = SL.library, method = "method.NNLS")))

replicate(5, system.time(mcSuperLearner(Y = Y, X = X, newX = newX,
  SL.library = SL.library, method = "method.NNLS")))

cl <- makeCluster(2, type = 'PSOCK')
replicate(5, system.time(snowSuperLearner(cl, Y = Y, X = X, newX = newX,
  SL.library = SL.library, method = "method.NNLS")))
stopCluster(cl)

## End(Not run)

```

SuperLearner.control *Control parameters for the SuperLearner*

Description

Control parameters for the SuperLearner

Usage

```
SuperLearner.control(saveFitLibrary = TRUE, trimLogit = 0.001)
```

Arguments

`saveFitLibrary` Logical. Should the fit for each algorithm be saved in the output from SuperLearner.
`trimLogit` number between 0.0 and 0.5. What level to truncate the logit transformation to maintain a bounded loss function when using the NNloglik method.

Value

A list containing the control parameters.

SuperLearner.CV.control *Control parameters for the cross validation steps in SuperLearner*

Description

Control parameters for the cross validation steps in SuperLearner

Usage

```
SuperLearner.CV.control(V = 10L, stratifyCV = FALSE, shuffle = TRUE,
  validRows = NULL)
```

Arguments

<code>V</code>	Integer. Number of splits for the V-fold cross-validation step. The default is 10. In most cases, between 10 and 20 splits works well.
<code>stratifyCV</code>	Logical. Should the data splits be stratified by a binary response? Attempts to maintain the same ratio in each training and validation sample.
<code>shuffle</code>	Logical. Should the rows of X be shuffled before creating the splits.
<code>validRows</code>	A List. Use this to pass pre-specified rows for the sample splits. The length of the list should be V and each entry in the list should contain a vector with the row numbers of the corresponding validation sample.

Value

A list containing the control parameters

SuperLearnerNews	<i>Show the NEWS file for the SuperLearner package</i>
------------------	--

Description

Show the NEWS file of the SuperLearner package. The function is simply a wrapper for the RShowDoc function

Usage

```
SuperLearnerNews(...)  
SuperLearnerDocs(what = 'SuperLearnerR.pdf', ...)
```

Arguments

<code>...</code>	additional arguments passed to RShowDoc
<code>what</code>	specify what document to open. Currently supports the NEWS file and the PDF files 'SuperLearner.pdf' and 'SuperLearnerR.pdf'.

Value

A invisible character string given the path to the SuperLearner NEWS file

trimLogit	<i>truncated-probabilities logit transformation</i>
-----------	---

Description

computes the logit transformation on the truncated probabilities

Usage

```
trimLogit(x, trim = 1e-05)
```

Arguments

x	vector of probabilities.
trim	value to truncate probabilities at. Currently symmetric truncation (trim and 1-trim).

Value

logit transformed values

Examples

```
x <- c(0.00000001, 0.0001, 0.001, 0.01, 0.1, 0.3, 0.7, 0.9, 0.99,
       0.999, 0.9999, 0.99999999)
trimLogit(x, trim = 0.001)
data.frame(Prob = x, Logit = qlogis(x), trimLogit = trimLogit(x, 0.001))
```

write.method.template *Method to estimate the coefficients for the super learner*

Description

These functions contain the information on the loss function and the model to combine algorithms

Usage

```
write.method.template(file = "", ...)
```

a few built in options:

```
method.NNLS()
method.NNLS2()
method.NNloglik()
method.CC_LS()
method.CC_nloglik()
method.AUC(optim_method = "Nelder-Mead")
```

Arguments

file	A connection, or a character string naming a file to print to. Passed to <code>cat</code> .
optim_method	Passed to the <code>optim</code> call method. See <code>optim</code> for details.
...	Additional arguments passed to <code>cat</code> .

Details

A SuperLearner method must be a list (or a function to create a list) with exactly 3 elements. The 3 elements must be named `require`, `computeCoef` and `computePred`.

Value

A list containing 3 elements:

require	A character vector listing any required packages. Use NULL if no additional packages are required
computeCoef	A function. The arguments are: <code>Z</code> , <code>Y</code> , <code>libraryNames</code> , <code>obsWeights</code> , <code>control</code> , <code>verbose</code> . The value is a list with two items: <code>cvRisk</code> and <code>coef</code> . This function computes the coefficients of the super learner. As the super learner minimizes the cross-validated risk, the loss function information is contained in this function as well as the model to combine the algorithms in <code>SL.library</code> .
computePred	A function. The arguments are: <code>predY</code> , <code>coef</code> , <code>control</code> . The value is a numeric vector with the super learner predicted values.

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[SuperLearner](#)

Examples

```
write.method.template(file = '')
```

`write.screen.template` *screening algorithms for SuperLearner*

Description

Screening algorithms for SuperLearner to be used with `SL.library`.

Usage

```
write.screen.template(file = "", ...)
```

Arguments

file A connection, or a character string naming a file to print to. Passed to [cat](#).
 ... Additional arguments passed to [cat](#)

Details

Explain structure of a screening algorithm here:

Value

whichVariable A logical vector with the length equal to the number of columns in X. TRUE indicates the variable (column of X) should be included.

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[SuperLearner](#)

Examples

```
write.screen.template(file = '')
```

write.SL.template *Wrapper functions for prediction algorithms in SuperLearner*

Description

Template function for SuperLearner prediction wrappers and built in options.

Usage

```
write.SL.template(file = "", ...)
```

Arguments

file A connection, or a character string naming a file to print to. Passed to [cat](#).
 ... Additional arguments passed to [cat](#)

Details

Describe SL.* structure here

Value

A list with two elements:

<code>pred</code>	The predicted values for the rows in <code>newX</code> .
<code>fit</code>	A list. Contains all objects necessary to get predictions for new observations from specific algorithm.

Author(s)

Eric C Polley <eric.polley@nih.gov>

See Also

[SuperLearner](#)

Examples

```
write.SL.template(file = '')
```

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