

Package ‘GPLTR’

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Description Combining a generalized linear model with an additional tree part on the same scale. A four-step procedure is proposed to fit the model and test the joint effect of the selected tree part while adjusting on confounding factors. We also proposed an ensemble procedure based on the bagging to improve prediction accuracy and computed several scores of importance for variable selection. See 'Cyprien Mbogning et al.'(2014)<[doi:10.1186/2043-9113-4-6](https://doi.org/10.1186/2043-9113-4-6)> and 'Cyprien Mbogning et al.'(2015)<[doi:10.1159/000380850](https://doi.org/10.1159/000380850)> for an overview of all the methods implemented in this package.

License GPL (>= 2.0)

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GPLTR-package	<i>Fit a generalized partially linear tree-based regression model</i>
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Description

Combining a generalized linear model with an additional tree part on the same scale. A four-step procedure is proposed to fit the model and test the joint effect of the selected tree part while adjusting on confounding factors. We also proposed an ensemble procedure based on the bagging to improve prediction accuracy and computed several scores of importance for variable selection. See 'Cyprien Mbogning et al.'(2014)<doi:10.1186/2043-9113-4-6>, 'Cyprien Mbogning et al.'(2015)<doi:10.1159/000380850> for an overview of all the methods implemented in this package.

Details

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Author(s)

Cyprien Mbogning and Wilson Toussile

Maintainer: Cyprien Mbogning <cyprien.mbogning@gmail.com>

References

Mbogning, C., Perdry, H., Broet, P.: A Bagged partially linear tree-based regression procedure for prediction and variable selection. *Human Heredity*, 79(3-4):1 82-93 (2015)

Mbogning, C., Perdry, H., Toussile, W., Broet, P.: A novel tree-based procedure for deciphering the genomic spectrum of clinical disease entities. *Journal of Clinical Bioinformatics* 4:6, (2014)

Terry M. Therneau, Elizabeth J. Atkinson (2013) An Introduction to Recursive Partitioning Using the RPART Routines. Mayo Foundation.

Chen, J., Yu, K., Hsing, A., Therneau, T.M.: A partially linear tree-based regression model for assessing complex joint gene-gene and gene-environment effects. *Genetic Epidemiology* 31, 238-251 (2007)

Examples

```
#####
## Example on a public dataset: the burn data
#####
## The burn data are also displayed in the KMSurv package
#####
## Not run:
data(burn)

## Build the rpart tree with all the variables

rpart.burn <- rpart(D2 ~ Z1 + Z2 + Z3 + Z4 + Z5 + Z6 + Z7 + Z8 + Z9
                  + Z10 + Z11, data = burn, method = "class")

plot(rpart.burn, main = 'rpart tree')
text(rpart.burn, xpd = TRUE, cex = .6, use.n = TRUE)

## fit the PLTR model after adjusting on gender (Z2) using the proposed method

args.rpart <- list(minbucket = 10, maxdepth = 4, cp = 0, maxcompete = 0,
                  maxsurrogate = 0)
family <- "binomial"
X.names = "Z2"
Y.name = "D2"
G.names = c('Z1', 'Z3', 'Z4', 'Z5', 'Z6', 'Z7', 'Z8', 'Z9', 'Z10', 'Z11')

pltr.burn <- pltr.glm(burn, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 4, iterMin = 3, verbose = FALSE)

## Pruned back the maximal tree using either the BIC or the AIC criterion

pltr.burn_prun <- best.tree.BIC.AIC(xtree = pltr.burn$tree, burn, Y.name,
                                  X.names, family = family)

## plot the BIC selected tree

plot(pltr.burn_prun$tree$BIC, main = 'BIC selected tree')
text(pltr.burn_prun$tree$BIC, xpd = TRUE, cex = .6, col = 'blue')

## Summary of the selected tree by a BIC criterion

summary(pltr.burn_prun$tree$BIC)

## Summary of the final selected pltr model
```

```

summary(pltr.burn_prun$fit_glm$BIC)

## fit the PLTR model after adjusting on gender (Z2) using the parametric
## bootstrap method

## set numWorkers = 1 on a windows platform

args.parallel = list(numWorkers = 10)

best_bootstrap <- best.tree.bootstrap(pltr.burn$tree, burn, Y.name, X.names,
  G.names, B = 2000, BB = 2000, args.rpart = args.rpart, epsi = 0.008,
  iterMax = 6, iterMin = 5, family = family, LEVEL = 0.05, LB = FALSE,
  args.parallel = args.parallel, verbose = FALSE)

plot(best_bootstrap$selected_model$tree, main = 'original method')
text(best_bootstrap$selected_model$tree, xpd = TRUE)

## Bagging a set of basic unpruned pltr predictors
# ?bagging.pltr

Bag.burn <- bagging.pltr(burn, Y.name, X.names, G.names, family,
  args.rpart, epsi = 0.01, iterMax = 4, iterMin = 3,
  Bag = 10, verbose = FALSE, doprune = FALSE)

## The threshold values used

Bag.burn$CUT

## The set of PLTR models in the bagging procedure

PLTR_BAG.burn <- Bag.burn$Glm_BAG

## The set of trees in the bagging procedure

TREE_BAG.burn <- Bag.burn$Tree_BAG

## Use the bagging procedure to predict new features
# ?predict_bagg.pltr

Pred_Bag.burn <- predict_bagg.pltr(Bag.burn, Y.name, newdata = burn,
  type = "response", threshold = seq(0, 1, by = 0.1))

## The confusion matrix for each threshold value using the majority vote

Pred_Bag.burn$CONF1

## The prediction error for each threshold value

Pred_Bag.burn$PRED_ERROR1

## Compute the variable importances using the bagging procedure

```

```

Var_Imp_BAG.burn <- VIMPBAG(Bag.burn, burn, Y.name)

## Importance score using the permutaion method for each thresshold value

Var_Imp_BAG.burn$PIS

## Shadow plot of three proposed scores

par(mfrow=c(1,3))
barplot(Var_Imp_BAG.burn$PIS$CUT5, main = 'PIS', horiz = TRUE, las = 1,
        cex.names = .8, col = 'lightblue')
barplot(Var_Imp_BAG.burn$DIS, main = 'DIS', horiz = TRUE, las = 1,
        cex.names = .8, col = 'grey')
barplot(Var_Imp_BAG.burn$DDIS, main = 'DDIS', horiz = TRUE, las = 1,
        cex.names = .8, col = 'purple')

## End(Not run)

```

bag.aucoob

AUC on the Out Of Bag samples

Description

Compute the AUC on the OOB samples of the bagging procedure for the binomial family. The true and false positive rates are also returned and could be helpfull for plotting the ROC curves.

Usage

```
bag.aucoob(bag_pltr, xdata, Y.name)
```

Arguments

bag_pltr	The output of the function bagging.pltr
xdata	The learning dataset containing the dependent variable, the confounding variables and the predictors variables
Y.name	The name of the binary dependent variable

Details

The thresshold values used for computing the AUC are defined when building the bagging predictor. see [bagging.pltr](#) for the convenient parameterization.

Value

A list of 4 elements

AUCOOB	the AUC computed on OOB samples of the Bagging procedure
TPR	the true positive rate for several thresshold values
FPR	the false positive rate for several thresshold values
OOB	the Out Of Bag error for each thresshold value

Note

The plot of the ROC curve is straightforward using the TPR and FPR obtained with the function `bag.aucoob`

Author(s)

Cyprien Mbogning

References

Mbogning, C., Perdry, H., Broet, P.: A Bagged partially linear tree-based regression procedure for prediction and variable selection (submitted 2014)

Examples

```
##
```

<code>bagging.pltr</code>	<i>bagging pltr models</i>
---------------------------	----------------------------

Description

bagging procedure to aggregate several PLTR models for accurate prediction and variable selection

Usage

```
bagging.pltr(xdata, Y.name, X.names, G.names, family = "binomial",
  args.rpart, epsi = 0.001, iterMax = 5, iterMin = 3, LB = FALSE,
  args.parallel = list(numWorkers = 1),
  Bag = 20, Pred_Data = data.frame(), verbose = TRUE, doprune = FALSE
  , threshold = seq(0, 1, by = 0.1))
```

Arguments

<code>xdata</code>	the learning data frame
<code>Y.name</code>	the name of the binary dependent variable
<code>X.names</code>	the names of independent variables to consider in the linear part of the glm and as offset in the tree part
<code>G.names</code>	the names of independent variables to consider in the tree part of the hybrid glm.
<code>family</code>	the glm family considered depending on the type of the dependent variable (only the binomial family works in this function for the moment) .
<code>args.rpart</code>	a list of options that control details of the rpart algorithm. <code>minbucket</code> : the minimum number of observations in any terminal <leaf> node; <code>cp</code> : complexity parameter (Any split that does not decrease the overall lack of fit by a factor of <code>cp</code> is not attempted); <code>maxdepth</code> : the maximum depth of any node of the final tree, with the root node counted as depth 0. ... See rpart.control for further details

epsi	a threshold value to check the convergence of the algorithm
iterMax	the maximal number of iteration to consider
iterMin	the minimum number of iteration to consider
LB	a binary indicator with values TRUE or FALSE indicating whether the loading is balanced or not in the parallel computing. It is nevertheless useless on a windows platform. See mclapply
args.parallel	a list of two elements containing the number of workers and the type of parallelization to achieve see mclapply .
Bag	The number of Bagging samples to consider
Pred_Data	An optional data frame to validate the bagging procedure (the test dataset)
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging)
doprune	a binary indicator with values TRUE or FALSE indicating whether the set of trees in the bagging procedure are pruned (by a BIC procedure) or not
threshold	a vector of numerical values between 0 and 1 used as threshold values for the computation of the OOB error rate

Details

For the Bagging procedure, it is mandatory to set `maxcompete = 0` and `maxsurrogate = 0` within the `rpart` arguments. This will ensure the correct calculation of the importance of variables.

Value

A list with eleven elements

IND_OOB	A list of length <code>Bag</code> containing the Out Of Bag (OOB) individuals for each PLTR model.
EOOB	The vector of OOB errors of the bagging procedure for each threshold value.
OOB_ERRORS_PBP	A matrix with <code>Bag</code> columns and <code>threshold</code> rows containing OOB error of each PLTR model in the bagging sequence for each threshold value.
OOB_ERROR_PBP	A vector containing the mean of <code>OOB_ERRORS_PBP</code> for each threshold value.
Tree_BAG	A list of length <code>Bag</code> containing the bagging trees
Glm_BAG	A list of length <code>Bag</code> containing the bagging <code>pltr</code> model; could be helpful for prediction on new features.
LOST	The 0, 1 lost matrix for OOB observations at each threshold value
TEST	A value of NULL if <code>Pred_Data</code> is not available. A list of three elements otherwise: <code>PRED_ERROR</code> : the estimated error of the Bagging procedure on the test sample for each threshold value; <code>PRED_IND</code> : A list of length the length of the threshold vector, each element containing a matrix with the prediction of the testing data individuals using each PLTR model of the bagging sequence (column by column); <code>FINAL_PRED_IND</code> : A list containing the final prediction of each individual of the testing data by the bagging procedure (the modal prediction) for each threshold value.

Var_IMP	A numeric vector containing the relative variable importance of the bagging procedure
Timediff	The execution time of the bagging procedure
CUT	The threshold value used inside the bagging procedure

Author(s)

Cyprien Mbogning

References

Mbogning, C., Perdry, H., Broet, P.: A Bagged partially linear tree-based regression procedure for prediction and variable selection. *Human Heredity* (To appear) (2015)

Leo Breiman: Bagging Predictors. *Machine Learning*, 24, 123-140 (1996)

See Also

[predict_bagg.pltr](#)

Examples

```
## Not run:
##load the data set

data(burn)

## set the parameters

args.rpart <- list(minbucket = 10, maxdepth = 4, cp = 0, maxsurrogate = 0)
family <- "binomial"
Y.name <- "D2"
X.names <- "Z2"
G.names <- c('Z1','Z3','Z4','Z5','Z6','Z7','Z8','Z9','Z10','Z11')
args.parallel = list(numWorkers = 1)

## Bagging a set of basic unpruned pltr predictors

Bag.burn <- bagging.pltr(burn, Y.name, X.names, G.names, family,
  args.rpart,epsi = 0.01, iterMax = 4, iterMin = 3,
  Bag = 20, verbose = FALSE, doprune = FALSE)

## End(Not run)
```

best.tree.BIC.AIC *Prunning the Maximal tree*

Description

this function is set to prune back the maximal tree by using the BIC or the AIC criterion.

Usage

```
best.tree.BIC.AIC(xtree, xdata, Y.name, X.names,  
                  family = "binomial", verbose = TRUE)
```

Arguments

xtree	a tree to prune
xdata	the dataset used to build the tree
Y.name	the name of the dependent variable
X.names	the names of independent confounding variables to consider in the linear part of the glm
family	the glm family considered depending on the type of the dependent variable.
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging)

Value

a list of four elements:

best_index	The size of the selected trees by BIC and AIC
tree	The selected trees by BIC and AIC
fit_glm	The fitted pltr models selected with BIC, and AIC
Timediff	The execution time of the selection procedure

Author(s)

Cyprien Mbogning and Wilson Toussile

References

Mbogning, C., Perdry, H., Toussile, W., Broet, P.: A novel tree-based procedure for deciphering the genomic spectrum of clinical disease entities. *Journal of Clinical Bioinformatics* 4:6, (2014)

Akaike, H.: A new look at the statistical model identification. *IEEE Trans. Automat. Control* AC-19, 716-723 (1974)

Schwarz, G.: Estimating the dimension of a model. *The Annals of Statistics* 6, 461-464 (1978)

See Also

[best.tree.CV](#), [pltr.glm](#)

Examples

```

data(burn)

args.rpart <- list(minbucket = 10, maxdepth = 4, cp = 0, maxcompete = 0,
                  maxsurrogate = 0)
family <- "binomial"
X.names = "Z2"
Y.name = "D2"
G.names = c('Z1', 'Z3', 'Z4', 'Z5', 'Z6', 'Z7', 'Z8', 'Z9', 'Z10', 'Z11')

pltr.burn <- pltr.glm(burn, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 4, iterMin = 3, verbose = FALSE)

## Pruned back the maximal tree using either the BIC or the AIC criterion

pltr.burn_prun <- best.tree.BIC.AIC(xtree = pltr.burn$tree, burn, Y.name,
                                   X.names, family = family)

## plot the BIC selected tree

plot(pltr.burn_prun$tree$BIC, main = 'BIC selected tree')
text(pltr.burn_prun$tree$BIC, xpd = TRUE, cex = .6, col = 'blue')

## Not run:
##load the data set

data(data_pltr)

## Set the parameters

args.rpart <- list(minbucket = 40, maxdepth = 10, cp = 0)
family <- "binomial"
Y.name <- "Y"
X.names <- "G1"
G.names <- paste("G", 2:15, sep="")

## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 5, iterMin = 3)

##pruned back the maximal tree by BIC or AIC criterion

tree_select <- best.tree.BIC.AIC(xtree = fit_pltr$tree, data_pltr, Y.name,
                                X.names, family = family)

plot(tree_select$tree$BIC, main = 'BIC TREE')
text(tree_select$tree$BIC, minlength = 0L, xpd = TRUE, cex = .6)

```

```
## End(Not run)
```

```
best.tree.bootstrap    parametric bootstrap on a pltr model
```

Description

a parametric bootstrap procedure to select and test at the same time the selected tree

Usage

```
best.tree.bootstrap(xtree, xdata, Y.name, X.names, G.names, B = 10, BB = 10,
  args.rpart = list(cp = 0, minbucket = 20, maxdepth = 10), epsi = 0.001,
  iterMax = 5, iterMin = 3, family = "binomial", LEVEL = 0.05, LB = FALSE,
  args.parallel = list(numWorkers = 1), verbose = TRUE)
```

Arguments

<code>xtree</code>	the maximal tree obtained by the function <code>pltr.glm</code>
<code>xdata</code>	the data frame used to build <code>xtree</code>
<code>Y.name</code>	the name of the dependent variable
<code>X.names</code>	the names of independent variables to consider in the linear part of the glm
<code>G.names</code>	the names of independent variables to consider in the tree part of the hybrid glm.
<code>B</code>	the size of the bootstrap sample
<code>BB</code>	the size of the bootstrap sample to compute the adjusted p-value
<code>args.rpart</code>	a list of options that control details of the <code>rpart</code> algorithm. <code>minbucket</code> : the minimum number of observations in any terminal <leaf> node; <code>cp</code> : complexity parameter (Any split that does not decrease the overall lack of fit by a factor of <code>cp</code> is not attempted); <code>maxdepth</code> : the maximum depth of any node of the final tree, with the root node counted as depth 0. ... See rpart.control for further details
<code>epsi</code>	a threshold value to check the convergence of the algorithm
<code>iterMax</code>	the maximal number of iteration to consider
<code>iterMin</code>	the minimum number of iteration to consider
<code>family</code>	the glm family considered depending on the type of the dependent variable.
<code>LEVEL</code>	the level of the test
<code>LB</code>	a binary indicator with values <code>TRUE</code> or <code>FALSE</code> indicating whether the loading is balanced or not in the parallel computing. It is useless on a windows platform.
<code>args.parallel</code>	parameters of the parallelization. See mclapply for more details
<code>verbose</code>	Logical; <code>TRUE</code> for printing progress during the computation (helpful for debugging)

Value

a list with six elements

selected_model

a list with the fit of the selected pltr model `fit_glm`, the selected tree `tree`, the p-value of the selected tree `p.value`, the adjusted p-value of the selected tree `adj_p.value` and an indicator `Tree_Selected` to assess whether the test is significant or not.

`fit_glm` the fitted pltr model under the null hypothesis if the test is not significant

`Timediff` The execution time of the parametric bootstrap procedure

`comp_p_values` The P-values of the competing trees

`Badj` The number of samples used in the inner level of the procedure

`BBadj` The number of samples used in the outer level of the procedure

Author(s)

Cyprien Mbogning and Wilson Toussile

References

Chen, J., Yu, K., Hsing, A., Therneau, T.M.: A partially linear tree-based regression model for assessing complex joint gene-gene and gene-environment effects. *Genetic Epidemiology* 31, 238-251 (2007)

See Also

[p.val.tree](#)

Examples

```
#load the data set
data(data_pltr)
args.rpart <- list(minbucket = 40, maxdepth = 10, cp = 0)
family <- "binomial"
Y.name <- "Y"
X.names <- "G1"
G.names <- paste("G", 2:15, sep="")
## Not run:
## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names,
  args.rpart = args.rpart, family = family, iterMax = 5, iterMin = 3)

## select an test the selected tree by a parametric bootstrap procedure
args.parallel = list(numWorkers = 1, type = "PSOCK")

best_bootstrap <- best.tree.bootstrap(fit_pltr$tree, data_pltr, Y.name, X.names,
  G.names, B = 10, BB = 10, args.rpart = args.rpart, epsi = 0.001,
  iterMax = 5, iterMin = 3, family = family, LEVEL = 0.05, LB = FALSE,
  args.parallel = args.parallel)
```

```
## End(Not run)
```

best.tree.CV	<i>Prunning the Maximal tree</i>
--------------	----------------------------------

Description

this function is set to prune back the maximal tree by using a K-fold cross-validation procedure.

Usage

```
best.tree.CV(xtree, xdata, Y.name, X.names, G.names, family = "binomial",
args.rpart = list(cp = 0, minbucket = 20, maxdepth = 10), epsi = 0.001,
iterMax = 5, iterMin = 3, ncv = 10, verbose = TRUE)
```

Arguments

xtree	a tree to prune
xdata	the dataset used to build the tree
Y.name	the name of the dependent variable
X.names	the names of independent variables to consider in the linear part of the glm
G.names	the names of independent variables to consider in the tree part of the hybrid glm.
family	the glm family considered depending on the type of the dependent variable.
args.rpart	a list of options that control details of the rpart algorithm. minbucket: the minimum number of observations in any terminal <leaf> node; cp: complexity parameter (Any split that does not decrease the overall lack of fit by a factor of cp is not attempted); maxdepth: the maximum depth of any node of the final tree, with the root node counted as depth 0. ... See rpart.control for further details
epsi	a treshold value to check the convergence of the algorithm
iterMax	the maximal number of iteration to consider
iterMin	the minimum number of iteration to consider
ncv	The number of folds to consider for the cross-validation
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging)

Value

a list of five elements:

best_index	The size of the selected tree by the cross-validation procedure
tree	The selected tree by CV
fit_glm	The fitted gpltr models selected with CV

CV_ERRORS	A list of two elements containing the cross-validation error of the selected tree by the CV procedure and a vector of cross-validation errors of all the competing models
Timediff	The execution time of the Cross-Validation procedure

Author(s)

Cyprien Mbogning

References

Mbogning, C., Perdry, H., Toussile, W., Broet, P.: A novel tree-based procedure for deciphering the genomic spectrum of clinical disease entities. *Journal of Clinical Bioinformatics* 4:6, (2014)

See Also

[best.tree.BIC.AIC, pltr.glm](#)

Examples

```
## Not run:
##load the data set

data(data_pltr)

## set the parameters

args.rpart <- list(minbucket = 40, maxdepth = 10, cp = 0)
family <- "binomial"
Y.name <- "Y"
X.names <- "G1"
G.names <- paste("G", 2:15, sep="")

## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 5, iterMin = 3)

##pruned back the maximal tree by a cross-validation procedure

tree_selected <- best.tree.CV(fit_pltr$tree, data_pltr, Y.name, X.names, G.names,
                              family = family, args.rpart = args.rpart, epsi = 0.001, iterMax = 5,
                              iterMin = 3, ncv = 10)

plot(tree_selected$tree, main = 'CV TREE')
text(tree_selected$tree, minlength = 0L, xpd = TRUE, cex = .6)

## End(Not run)
```

best.tree.permute *permutation test on a pltr model*

Description

a unified permutation test procedure to select and test at the same time the selected tree

Usage

```
best.tree.permute(xtree, xdata, Y.name, X.names, G.names, B = 10,
  args.rpart = list(cp = 0, minbucket = 20, maxdepth = 10), epsi = 0.001,
  iterMax = 5, iterMin = 3, family = "binomial", LEVEL = 0.05,
  LB = FALSE, args.parallel = list(numWorkers = 1, type = "PSOCK"), verbose = TRUE)
```

Arguments

xtree	the maximal tree obtained by the function <code>pltr.glm</code>
xdata	the data frame used to build xtree
Y.name	the name of the dependent variable
X.names	the names of independent variables to consider in the linear part of the glm. For this function, only a binary variable is supported.
G.names	the names of independent variables to consider in the tree part of the hybrid glm.
B	the size of the bootstrap sample
args.rpart	a list of options that control details of the rpart algorithm. <code>minbucket</code> : the minimum number of observations in any terminal <leaf> node; <code>cp</code> : complexity parameter (Any split that does not decrease the overall lack of fit by a factor of <code>cp</code> is not attempted); <code>maxdepth</code> : the maximum depth of any node of the final tree, with the root node counted as depth 0. ... See rpart.control for further details
epsi	a threshold value to check the convergence of the algorithm
iterMax	the maximal number of iteration to consider
iterMin	the minimum number of iteration to consider
family	the binomial family.
LEVEL	the level of the test
LB	a binary indicator with values TRUE or FALSE indicating whether the loading is balanced or not in the parallel computing. It is useless on a windows platform.
args.parallel	parameters of the parallelization. See mclapply for more details.
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging)


```
## End(Not run)
```

burn	<i>burn dataset</i>
------	---------------------

Description

The burn data frame has 154 rows and 17 columns.

Usage

```
data(burn)
```

Format

A data frame with 154 observations on the following 17 variables.

Obs Observation number

Z1 Treatment: 0=routine bathing 1=Body cleansing

Z2 Gender (0=male 1=female)

Z3 Race: 0=nonwhite 1=white

Z4 Percentage of total surface area burned

Z5 Burn site indicator: head 1=yes, 0=no

Z6 Burn site indicator: buttock 1=yes, 0=no

Z7 Burn site indicator: trunk 1=yes, 0=no

Z8 Burn site indicator: upper leg 1=yes, 0=no

Z9 Burn site indicator: lower leg 1=yes, 0=no

Z10 Burn site indicator: respiratory tract 1=yes, 0=no

Z11 Type of burn: 1=chemical, 2=scald, 3=electric, 4=flame

T1 Time to excision or on study time

D1 Excision indicator: 1=yes 0=no

T2 Time to prophylactic antibiotic treatment or on study time

D2 Prophylactic antibiotic treatment: 1=yes 0=no

T3 Time to straphylocous aureaus infection or on study time

D3 Straphylocous aureaus infection: 1=yes 0=no

Source

Klein and Moeschberger (1997) Survival Analysis Techniques for Censored and truncated data, Springer.

Ichida et al. Stat. Med. 12 (1993): 301-310.

Examples

```
data(burn)
## maybe str(burn) ;
```

data_pltr	<i>gpltr data example</i>
-----------	---------------------------

Description

A data frame to test the functions of the package

Usage

```
data(data_pltr)
```

Format

A data frame with 3000 observations on the following 16 variables.

G1 a numeric vector
G2 a factor with levels 0 1
G3 a factor with levels 0 1
G4 a factor with levels 0 1
G5 a factor with levels 0 1
G6 a binary numeric vector
G7 a binary numeric vector
G8 a binary numeric vector
G9 a binary numeric vector
G10 a binary numeric vector
G11 a binary numeric vector
G12 a binary numeric vector
G13 a binary numeric vector
G14 a binary numeric vector
G15 a binary numeric vector
Y a binary numeric vector

Details

The numeric variable G1 is considered as offset in the simulated PLTR model; the variables G2,...,G5 are used to simulate the tree part, while G6,...,G15 are noise variables.

Examples

```
data(data_pltr)
## maybe str(data_pltr) ...
```

nested.trees	<i>compute the nested trees</i>
--------------	---------------------------------

Description

Compute a sequence of nested competing trees for the pruning step

Usage

```
nested.trees(xtree, xdata, Y.name, X.names, MaxTreeSize = NULL,
family = "binomial", verbose = TRUE)
```

Arguments

xtree	a tree inheriting to the rpart method
xdata	the dataset used to build the tree
Y.name	the name of the dependent variable in the tree model
X.names	the names of independent variables considered as offset in the tree model
MaxTreeSize	The maximal size of the competing trees
family	the glm family considered depending on the type of the dependent variable.
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging)

Value

a list with 4 elements:

leaves	a list of leaves of the competing trees to consider for the optimal tree
null_deviance	the deviance of the null model (linear part of the glm)
deviances	a vector of deviances of the competing PLTR models
diff_deviances	a vector of the deviance differences between the competing PLTR models and the null model

Author(s)

Cyprien Mbogning and Wilson Toussile

Examples

```
## Not run:
## load the data set

data(data_pltr)
args.rpart <- list(minbucket = 40, maxdepth = 10, cp = 0)
family <- "binomial"
Y.name <- "Y"
```

```

X.names <- "G1"
G.names <- paste("G", 2:15, sep="")

## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 5, iterMin = 3)

## compute the competing trees

nested_trees <- nested.trees(fit_pltr$tree, data_pltr, Y.name, X.names,
                              MaxTreeSize = 10, family = family)

## End(Not run)

```

p.val.tree

Compute the p-value

Description

Test whether the selected tree by either BIC, AIC or CV procedure is significantly associated to the dependent variable or not, while adjusting for a confounding effect.

Usage

```

p.val.tree(xtree, xdata, Y.name, X.names, G.names, B = 10, args.rpart =
list(minbucket = 40, maxdepth = 10, cp = 0), epsi = 0.001, iterMax = 5,
iterMin = 3, family = "binomial", LB = FALSE,
args.parallel = list(numWorkers = 1), index = 4, verbose = TRUE)

```

Arguments

xtree	the maximal tree obtained by the function <code>pltr.glm</code>
xdata	the data frame used to build xtree
Y.name	the name of the dependent variable
X.names	the names of independent confounding variables to consider in the linear part of the <code>glm</code>
G.names	the names of independent variables to consider in the tree part of the hybrid <code>glm</code> .
B	the resampling size of the deviance difference
args.rpart	a list of options that control details of the <code>rpart</code> algorithm. <code>minbucket</code> : the minimum number of observations in any terminal <leaf> node; <code>cp</code> : complexity parameter (Any split that does not decrease the overall lack of fit by a factor of <code>cp</code> is not attempted); <code>maxdepth</code> : the maximum depth of any node of the final tree, with the root node counted as depth 0. ... See rpart.control for further details
epsi	a threshold value to check the convergence of the algorithm

iterMax	the maximal number of iteration to consider
iterMin	the minimum number of iteration to consider
family	the glm family considered depending on the type of the dependent variable.
LB	a binary indicator with values TRUE or FALSE indicating whether the loadings are balanced or not in the parallel computing
args.parallel	parameters of the parallelization. See mclapply for more details.
index	the size of the selected tree (by the functions best.tree.BIC.AIC or best.tree.CV) using one of the proposed criteria
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging)

Value

A list of three elements:

p.value	The P-value of the selected tree
Timediff	The execution time of the test procedure
Badj	The number of samples used inside the the procedure

Author(s)

Cyprien Mbogning

References

- Mbogning, C., Perdry, H., Toussile, W., Broet, P.: A novel tree-based procedure for deciphering the genomic spectrum of clinical disease entities. *Journal of Clinical Bioinformatics* 4:6, (2014)
- Fan, J., Zhang, C., Zhang, J.: Generalized likelihood ratio statistics and WILKS phenomenon. *Annals of Statistics* 29(1), 153-193 (2001)

See Also

[best.tree.bootstrap](#), [best.tree.permute](#)

Examples

```
## Not run:
## load the data set

data(data_pltr)

## set the parameters

args.rpart <- list(minbucket = 40, maxdepth = 10, cp = 0)
family <- "binomial"
Y.name <- "Y"
X.names <- "G1"
G.names <- paste("G", 2:15, sep="")
```

```

## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 5, iterMin = 3)

##pruned back the maximal tree by BIC or AIC criterion

tree_select <- best.tree.BIC.AIC(xtree = fit_pltr$tree, data_pltr, Y.name,
                                X.names, family = family)

## Compute the p-value of the selected tree by BIC

args.parallel = list(numWorkers = 10, type = "PSOCK")
index = tree_select$best_index[[1]]
p_value <- p.val.tree(xtree = fit_pltr$tree, data_pltr, Y.name, X.names, G.names,
                     B = 100, args.rpart = args.rpart, epsi = 1e-3,
                     iterMax = 5, iterMin = 3, family = family, LB = FALSE,
                     args.parallel = args.parallel, index = index)

## End(Not run)

```

pltr.glm

Partially tree-based regression model function

Description

The `pltr.glm` function is designed to fit an hybrid glm model with an additive tree part on a glm scale.

Usage

```

pltr.glm(data, Y.name, X.names, G.names, family = "binomial",
         args.rpart = list(cp = 0, minbucket = 20, maxdepth = 10),
         epsi = 0.001, iterMax = 5, iterMin = 3, verbose = TRUE)

```

Arguments

<code>data</code>	a data frame containing the variables in the model
<code>Y.name</code>	the name of the dependent variable
<code>X.names</code>	the names of independent variables to consider in the linear part of the glm
<code>G.names</code>	the names of independent variables to consider in the tree part of the hybrid glm.
<code>family</code>	the glm family considered depending on the type of the dependent variable.
<code>args.rpart</code>	a list of options that control details of the <code>rpart</code> algorithm. <code>minbucket</code> : the minimum number of observations in any terminal <leaf> node; <code>cp</code> : complexity parameter (Any split that does not decrease the overall lack of fit by a factor of <code>cp</code> is not attempted); <code>maxdepth</code> : the maximum depth of any node of the final tree, with the root node counted as depth 0. ... See rpart.control for further details

epsi	a treshold value to check the convergence of the algorithm
iterMax	the maximal number of iteration to consider
iterMin	the minimum number of iteration to consider
verbose	Logical; TRUE for printing progress during the computation (helpful for debugging)

Details

The `pltr.glm` function use an iterative procedure to fit the linear part of the `glm` and the tree part. The tree obtained at the convergence of the procedure is a maximal tree which overfits the data. It's then mandatory to pruned back this tree by using one of the proposed criteria (BIC, AIC and CV).

Value

A list with four elements:

fit	the <code>glm</code> fitted on the confounding factors at the end of the iterative algorithm
tree	the maximal tree obtained at the end of the algorithm
nber_iter	the number of iterations used by the algorithm
Timediff	The execution time of the iterative procedure

Note

The tree obtained at the end of these iterative procedure usually overfits the data. It's therefore mendatory to use either `best.tree.BIC.AIC` or `best.tree.CV` to prunne back the tree.

Author(s)

Cyprien Mbogning and Wilson Toussile

References

- Mbogning, C., Perdry, H., Toussile, W., Broet, P.: A novel tree-based procedure for deciphering the genomic spectrum of clinical disease entities. *Journal of Clinical Bioinformatics* 4:6, (2014)
- Terry M. Therneau, Elizabeth J. Atkinson (2013) An Introduction to Recursive Partitioning Using the RPART Routines. Mayo Foundation.
- Chen, J., Yu, K., Hsing, A., Therneau, T.M.: A partially linear tree-based regression model for assessing complex joint gene-gene and gene-environment effects. *Genetic Epidemiology* 31, 238-251 (2007)

See Also

[rpart](#)

Examples

```

data(burn)

args.rpart <- list(minbucket = 10, maxdepth = 4, cp = 0, maxcompete = 0,
                  maxsurrogate = 0)
family <- "binomial"
X.names = "Z2"
Y.name = "D2"
G.names = c('Z1', 'Z3', 'Z4', 'Z5', 'Z6', 'Z7', 'Z8', 'Z9', 'Z10', 'Z11')

pltr.burn <- pltr.glm(burn, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 4, iterMin = 3, verbose = FALSE)

## Not run:
## load the data set

data(data_pltr)

## set the parameters

args.rpart <- list(minbucket = 40, maxdepth = 10, cp = 0)
family <- "binomial"
Y.name <- "Y"
X.names <- "G1"
G.names <- paste("G", 2:15, sep="")

## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names, args.rpart = args.rpart,
                   family = family, iterMax = 5, iterMin = 3)

plot(fit_pltr$tree, main = 'MAXIMAL TREE')
text(fit_pltr$tree, minlength = 0L, xpd = TRUE, cex = .6)

## End(Not run)

```

predict_bagg.pltr *prediction on new features*

Description

Prediction on new features using a set of bagging pltr models

Usage

```

predict_bagg.pltr(bag_pltr, Y.name, newdata, type = "response",
                 thresshold = seq(0, 1, by = 0.1))

```


Arguments

bag_pltr	the bagging result obtained with the function bagging.pltr
Y.name	the name of the binary dependent variable
newdata	a data frame in which to look for predictors and the dependant variable.
type	the type of prediction required. type = "response" is the default; It gives the predicted probabilities. At this stage of the package, only this type is take into account. Other types such as "link" and "terms" are useless.
threshold	a vector of cutoff values for binary prediction. Could be helpfull for computing the AUC on the test sample.

Value

A list with 8 elements

FINAL_PRED_IND1	A list of size the length of the threshold vector, containing the final prediction of each individual of the testing data by the bagging procedure using the majority rule (the modal prediction).
FINAL_PRED_IND2	A list of size the length of the threshold vector, containing the final prediction of each individual of the testing data by the bagging procedure using the mean estimated probability.
PRED_ERROR1	A vector of estimated errors of the Bagging procedure on the test sample for each threshold value using FINAL_PRED_IND1.
PRED_ERROR2	A vector of estimated errors of the Bagging procedure on the test sample for each threshold value using FINAL_PRED_IND2.
CONF1	A list of confusion matrix using FINAL_PRED_IND1
CONF2	A list of confusion matrix using FINAL_PRED_IND2
PRED_ERRORS_PBP	A list of size the length of the threshold vector. Each element representing the prediction error obtained via each predictor in the bagging sequence for each threshold value
PRED_ERROR_PBP	A vector containing the mean of PRED_ERRORS_PBP for each threshold value

Author(s)

Cyprien Mbogning

References

Mbogning, C., Perdry, H., Broet, P.: A Bagged partially linear tree-based regression procedure for prediction and variable selection. *Human Heredity* (To appear), (2015)

See Also

[bagging.pltr](#), [predict.glm](#)

Examples

```

## Not run:
## load the data set

data(burn)

## set the parameters

args.rpart <- list(minbucket = 10, maxdepth = 4, cp = 0, maxsurrogate = 0)
family <- "binomial"
Y.name <- "D2"
X.names <- "Z2"
G.names <- c('Z1', 'Z3', 'Z4', 'Z5', 'Z6', 'Z7', 'Z8', 'Z9', 'Z10', 'Z11')
args.parallel = list(numWorkers = 1)

## Bagging a set of basic unpruned pltr predictors

Bag.burn <- bagging.pltr(burn, Y.name, X.names, G.names, family,
                        args.rpart, epsi = 0.01, iterMax = 4, iterMin = 3,
                        Bag = 20, verbose = FALSE, doprune = FALSE)

## Use the bagging procedure to predict new features

# ?predict_bagg.pltr

Pred_Bag.burn <- predict_bagg.pltr(Bag.burn, Y.name, newdata = burn,
                                   type = "response", threshold = seq(0, 1, by = 0.1))

## The confusion matrix for each threshold value using the majority vote

Pred_Bag.burn$CONF1

## End(Not run)

```

predict_pltr

prediction

Description

prediction on new features using a pltr tree and the name of the confounding variable

Usage

```

predict_pltr(xtree, xdata, Y.name, X.names, newdata, type = "response",
            family = 'binomial', threshold = seq(0.1, 0.9, by = 0.1))

```

Arguments

xtree a tree obtained with the pltr procedure

xdata	the dataframe used to learn the pltr model
Y.name	the name of the main variable
X.names	the names of the confounding variables
newdata	the new data with all the predictors and the dependent variable
type	the type of prediction
family	the glm family considered
threshold	the threshold value to consider for binary prediction. It could be a vector, helping to compute the AUC

Value

A list of two element

predict_glm	the predicted vector, depending on the family used. For the binomial family with a vector of threshold, a matrix with each column corresponding to a threshold value
ERR_PRED	either the prediction error of the pltr procedure on the test set or a vector of prediction error when the family is binomial with a vector of threshold values

Author(s)

Cyprien Mbogning

References

Mbogning, C., Perdry, H., Toussile, W., Broet, P.: A novel tree-based procedure for deciphering the genomic spectrum of clinical disease entities. *Journal of Clinical Bioinformatics* 4:6, (2014)

See Also

[pltr.glm](#), [predict.glm](#)

Examples

```
##
```

tree2glm	<i>tree to GLM</i>
----------	--------------------

Description

fit the PLTR model for a given tree. The tree is coerced into dummy covariates.

Usage

```
tree2glm(xtree, xdata, Y.name, X.names, family = "binomial")
```

Arguments

xtree	a tree inherits from the rpart method
xdata	a data frame containing the variables in the model
Y.name	the name of the dependent variable
X.names	the names of independent variables to consider in the linear part of the glm
family	the glm family considered depending on the type of the dependent variable.

Value

the pltr fitted model (fit)

Author(s)

Cyprien Mbogning and Wilson Toussile

Examples

```
## Not run:
##load the data set

data(data_pltr)

## set the parameters

args.rpart <- list(minbucket = 40, cp = 0)
family <- "binomial"
Y.name <- "Y"
X.names <- "G1"
G.names <- paste("G", 2:15, sep="")

## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names, args.rpart = args.rpart,
                   family = family, iterMax = 5, iterMin = 3)

## Coerce a tree into a glm model using the confounding factor

fit_glm <- tree2glm(fit_pltr$tree, data_pltr, Y.name, X.names,
                   family = family)

summary(fit_glm)

## End(Not run)
```

tree2indicators	<i>From a tree to indicators (or dummy variables)</i>
-----------------	---

Description

Coerces a given tree structure to binary covariates.

Usage

```
tree2indicators(fit)
```

Arguments

`fit` a tree structure inheriting to the `rpart` method

Value

a list of indicators

Author(s)

Cyprien Mbogning and Wilson Toussile

Examples

```
## Not run:
## load the data set

data(data_pltr)

## set the parameters

args.rpart <- list(minbucket = 40, xval = 10, cp = 0)
family <- "binomial"
Y.name <- "Y"
X.names <- "G1"
G.names <- paste("G", 2:15, sep="")

## build a maximal tree

fit_pltr <- pltr.glm(data_pltr, Y.name, X.names, G.names, args.rpart = args.rpart,
                    family = family, iterMax = 5, iterMin = 3)

## Compute a list of indicator from the leaves of the tree fitted tree

tree2indicators(fit_pltr$tree)

## End(Not run)
```

VIMPBAG *score of importance for variables*

Description

Several variable importance scores are computed: the deviance importance score (DIS), the permutation importance score (PIS), the depth deviance importance score (DDIS), the minimal depth importance score (MinDepth) and the occurrence score (OCCUR).

Usage

```
VIMPBAG(BAGGRES, data, Y.name)
```

Arguments

BAGGRES	The output of the bagging procedure (bagging.pltr)
data	The learning dataframe used within the bagging procedure
Y.name	The name of the binary dependant variable used in the bagging procedure

Details

several choices for variable selection using the bagging procedure are proposed. A discussion about the scores of importance PIS, DIS, and DDIS is available in Mbogning et al. 2015

Value

A list with 9 elements

PIS	A list of length the length of the threshold value used in the bagging procedure, containing the permutation importance score displayed in decreasing order for each threshold value
StdPIS	The standard error of the PIS
OCCUR	The occurrence number for each variable in the bagging sequence displayed in decreasing order
DIS	The deviance importance score displayed in decreasing order
DDIS	The depth deviance importance score displayed in decreasing order
MinDepth	The minimal depth score for each variable, displayed in increasing order
dimtrees	A vector containing the dimensions of trees within the bagging sequence
E00B	A vector containing the OOB error of the bagging procedure for each threshold value
Bagfinal	The number of Bagging iterations used

Author(s)

Cyprien Mbogning

References

Mbogning, C., Perdry, H., Broet, P.: A Bagged partially linear tree-based regression procedure for prediction and variable selection. *Human Heredity (To appear)*, (2015)

See Also

[bagging.pltr](#)

Examples

```
## Not run:
## load the data set

data(burn)

## set the parameters

args.rpart <- list(minbucket = 10, maxdepth = 4, cp = 0, maxsurrogate = 0)
family <- "binomial"
Y.name <- "D2"
X.names <- "Z2"
G.names <- c('Z1', 'Z3', 'Z4', 'Z5', 'Z6', 'Z7', 'Z8', 'Z9', 'Z10', 'Z11')
args.parallel = list(numWorkers = 1)

## Bagging a set of basic unpruned pltr predictors

Bag.burn <- bagging.pltr(burn, Y.name, X.names, G.names, family,
                        args.rpart, epsi = 0.01, iterMax = 4, iterMin = 3,
                        Bag = 20, verbose = FALSE, doprune = FALSE)

## Several importance scores for variables, using the bagging procedure

Var_Imp_BAG.burn <- VIMPBAG(Bag.burn, burn, Y.name)

## Importance score using the permutaion method for each thresshold value

Var_Imp_BAG.burn$PIS

## Importance score using the deviance criterion

Var_Imp_BAG.burn$DIS

## End(Not run)
```

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