Package ‘ranger’

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Type Package
Title A Fast Implementation of Random Forests
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Description A fast implementation of Random Forests, particularly suited for high
dimensional data. Ensembles of classification, regression, survival and
probability prediction trees are supported. Data from genome-wide association
studies can be analyzed efficiently. In addition to data frames, datasets of
class ‘gwaa.data’ (R package ‘GenABEL’) and ‘dgCMatrix’ (R package ‘Matrix’) can
be directly analyzed.
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R topics documented:

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Case-specific random forests.

Description
In case-specific random forests (CSRF), random forests are built specific to the cases of interest. Instead of using equal probabilities, the cases are weighted according to their difference to the case of interest.

Usage
```r
csrp(formula, training_data, test_data, params1 = list(), params2 = list())
```

Arguments
- **formula**: Object of class `formula` or character describing the model to fit.
- **training_data**: Training data of class `data.frame`.
- **test_data**: Test data of class `data.frame`.
- **params1**: Parameters for the proximity random forest grown in the first step.
- **params2**: Parameters for the prediction random forests grown in the second step.

Details
The algorithm consists of 3 steps:
1. Grow a random forest on the training data
2. For each observation of interest (test data), the weights of all training observations are computed by counting the number of trees in which both observations are in the same terminal node.
3. For each test observation, grow a weighted random forest on the training data, using the weights obtained in step 2. Predict the outcome of the test observation as usual.

In total, n+1 random forests are grown, where n is the number observations in the test dataset. For details, see Xu et al. (2014).

Value

Predictions for the test dataset.

Author(s)

Marvin N. Wright

References


Examples

```r
## Split in training and test data
train.idx <- sample(nrow(iris), 2/3 * nrow(iris))
iris.train <- iris[train.idx, ]
iris.test <- iris[-train.idx, ]

## Run case-specific RF
csr.rf(Species ~ ., training_data = iris.train, test_data = iris.test,
      params1 = list(num.trees = 50, mtry = 4),
      params2 = list(num.trees = 5))
```

---

getTerminalNodeIDs

Get terminal node IDs (deprecated)

Description

This function is deprecated. Please use predict() with type = "terminalNodes" instead. This function calls predict() now.

Usage

getTerminalNodeIDs(rf, dat)

Arguments

rf

ranger object.

dat

New dataset. Terminal node IDs for this dataset are obtained.
Value

Matrix with terminal nodeIDs for all observations in dataset and trees.

Examples

```r
library(ranger)
rf <- ranger(Species ~ ., data = iris, num.trees = 5, write.forest = TRUE)
getTerminalNodeIDs(rf, iris)
```

Description

Grow two random forests on two cross-validation folds. Instead of out-of-bag data, the other fold is used to compute permutation importance. Related to the novel permutation variable importance by Janitza et al. (2015).

Usage

```r
holdoutRF()
```

Arguments

... All arguments are passed to `ranger()` (except importance, case.weights, replace and holdout).

Value

Hold-out random forests with variable importance.

Author(s)

Marvin N. Wright

References


See Also

`ranger`
importance.ranger

importance.ranger  
* ranger variable importance

---

**Description**

Extract variable importance of ranger object.

**Usage**

```r
## S3 method for class 'ranger'
importance(x, ...)
```

**Arguments**

- `x` ranger object.
- `...` Further arguments passed to or from other methods.

**Value**

Variable importance measures.

**Author(s)**

Marvin N. Wright

**See Also**

- `ranger`

---

importance_pvalues  
* ranger variable importance p-values

---

**Description**

Compute variable importance with p-values. For high dimensional data, the fast method of Janitza et al. (2016) can be used. The permutation approach of Altmann et al. (2010) is computationally intensive but can be used with all kinds of data. See below for details.

**Usage**

```r
importance_pvalues(x, method = c("janitza", "altmann"),
                   num.permutations = 100, formula = NULL, data = NULL, ...)
```
importance_pvalues

Arguments

x  ranger or holdoutRF object.
method  Method to compute p-values. Use "janitza" for the method by Janitza et al. (2016) or "altmann" for the non-parametric method by Altmann et al. (2010).
num_permutations  Number of permutations. Used in the "altmann" method only.
formula  Object of class formula or character describing the model to fit. Used in the "altmann" method only.
data  Training data of class data.frame or matrix. Used in the "altmann" method only.
...  Further arguments passed to ranger(). Used in the "altmann" method only.

Details

The method of Janitza et al. (2016) uses a clever trick: With an unbiased variable importance measure, the importance values of non-associated variables vary randomly around zero. Thus, all non-positive importance values are assumed to correspond to these non-associated variables and they are used to construct a distribution of the importance under the null hypothesis of no association to the response. Since only the non-positive values of this distribution can be observed, the positive values are created by mirroring the negative distribution. See Janitza et al. (2016) for details.

The method of Altmann et al. (2010) uses a simple permutation test: The distribution of the importance under the null hypothesis of no association to the response is created by several replications of permuting the response, growing an RF and computing the variable importance. The authors recommend 50-100 permutations. However, much larger numbers have to be used to estimate more precise p-values. We add 1 to the numerator and denominator to avoid zero p-values.

Value

Variable importance and p-value for each variable.

Author(s)

Marvin N. Wright

References


See Also

ranger
**Examples**

```r
require(ranger)

## Janitza's p-values with corrected Gini importance
n <- 50
p <- 400
dat <- data.frame(y = factor(rbinom(n, 1, .5)), replicate(p, runif(n)))
rf.sim <- ranger(y ~ ., dat, importance = "impurity_corrected")
importance_pvalues(rf.sim, method = "janitza")

## Permutation p-values
## Not run:
rf.iris <- ranger(Species ~ ., data = iris, importance = 'permutation')
importance_pvalues(rf.iris, method = "altmann", formula = Species ~ ., data = iris)

## End(Not run)
```

---

**parse.formula**

**Parse formula**

**Description**

Parse formula and return dataset containing selected columns. Interactions are supported for numerical columns only. An interaction column is the product of all interacting columns.

**Usage**

```r
parse.formula(formula, data, env = parent.frame())
```

**Arguments**

- `formula`: Object of class formula or character describing the model to fit.
- `data`: Training data of class data.frame.
- `env`: The environment in which the left hand side of `formula` is evaluated.

**Value**

Dataset including selected columns and interactions.
predict.ranger  

**Ranger prediction**

**Description**

Prediction with new data and a saved forest from Ranger.

**Usage**

```r
## S3 method for class 'ranger'
predict(object, data = NULL, predict.all = FALSE,
        num.trees = object$num.trees, type = "response",
        se.method = "infjack", quantiles = c(0.1, 0.5, 0.9), seed = NULL,
        num.threads = NULL, verbose = TRUE, ...)  
```

**Arguments**

- `object`  
  Ranger ranger object.
- `data`  
  New test data of class `data.frame` or `gwaa.data` (GenABEL).
- `predict.all`  
  Return individual predictions for each tree instead of aggregated predictions for all trees. Return a matrix (sample x tree) for classification and regression, a 3d array for probability estimation (sample x class x tree) and survival (sample x time x tree).
- `num.trees`  
  Number of trees used for prediction. The first `num.trees` in the forest are used.
- `type`  
  Type of prediction. One of 'response', 'se', 'terminalNodes', 'quantiles' with default 'response'. See below for details.
- `se.method`  
  Method to compute standard errors. One of 'jack', 'infjack' with default 'infjack'. Only applicable if `type = 'se'`. See below for details.
- `quantiles`  
  Vector of quantiles for quantile prediction. Set `type = 'quantiles'` to use.
- `seed`  
  Random seed. Default is NULL, which generates the seed from R. Set to 0 to ignore the R seed. The seed is used in case of ties in classification mode.
- `num.threads`  
  Number of threads. Default is number of CPUs available.
- `verbose`  
  Verbose output on or off.
- `...`  
  further arguments passed to or from other methods.

**Details**

For `type = 'response'` (the default), the predicted classes (classification), predicted numeric values (regression), predicted probabilities (probability estimation) or survival probabilities (survival) are returned. For `type = 'se'`, the standard error of the predictions are returned (regression only). The jackknife-after-bootstrap or infinitesimal jackknife for bagging is used to estimate the standard errors based on out-of-bag predictions. See Wager et al. (2014) for details. For `type = 'terminalNodes'`, the IDs of the terminal node in each tree for each observation in the
given dataset are returned. For type = 'quantiles', the selected quantiles for each observation are estimated. See Meinshausen (2006) for details.

If type = 'se' is selected, the method to estimate the variances can be chosen with se.method. Set se.method = 'jack' for jackknife-after-bootstrap and se.method = 'infjack' for the infinitesimal jackknife for bagging.

For classification and predict.all = TRUE, a factor levels are returned as numerics. To retrieve the corresponding factor levels, use rf$forest$levels, if rf is the ranger object.

Value

Object of class ranger.prediction with elements

- predictions: Predicted classes/values (only for classification and regression)
- unique.death.times: Unique death times (only for survival).
- chf: Estimated cumulative hazard function for each sample (only for survival).
- survival: Estimated survival function for each sample (only for survival).
- num.trees: Number of trees.
- num.independent.variables: Number of independent variables.
- tree.type: Type of forest/tree. Classification, regression or survival.
- num.samples: Number of samples.

Author(s)

Marvin N. Wright

References


See Also

- ranger

Description

Prediction with new data and a saved forest from Ranger.
Usage

## S3 method for class 'ranger.forest'

predict(object, data, predict.all = FALSE, 
num.trees = object$num.trees, type = "response", 
se.method = "infjack", seed = NULL, num.threads = NULL, 
verbose = TRUE, inbag.counts = NULL, ...)

Arguments

- **object**: Ranger ranger.forest object.
- **data**: New test data of class data.frame or gwaa.data (GenABEL).
- **predict.all**: Return individual predictions for each tree instead of aggregated predictions for all trees. Return a matrix (sample x tree) for classification and regression, a 3d array for probability estimation (sample x class x tree) and survival (sample x time x tree).
- **num.trees**: Number of trees used for prediction. The first num.trees in the forest are used.
- **type**: Type of prediction. One of 'response', 'se', 'terminalNodes', 'quantiles' with default 'response'. See below for details.
- **se.method**: Method to compute standard errors. One of 'jack', 'infjack' with default 'infjack'. Only applicable if type = 'se'. See below for details.
- **seed**: Random seed. Default is NULL, which generates the seed from R. Set to 0 to ignore the R seed. The seed is used in case of ties in classification mode.
- **num.threads**: Number of threads. Default is number of CPUs available.
- **verbose**: Verbose output on or off.
- **inbag.counts**: Number of times the observations are in-bag in the trees.
- **...**: further arguments passed to or from other methods.

Details

For type = 'response' (the default), the predicted classes (classification), predicted numeric values (regression), predicted probabilities (probability estimation) or survival probabilities (survival) are returned. For type = 'se', the standard error of the predictions are returned (regression only). The jackknife-after-bootstrap or infinitesimal jackknife for bagging is used to estimate the standard errors based on out-of-bag predictions. See Wager et al. (2014) for details. For type = 'terminalNodes', the IDs of the terminal node in each tree for each observation in the given dataset are returned.

If type = 'se' is selected, the method to estimate the variances can be chosen with se.method. Set se.method = 'jack' for jackknife after bootstrap and se.method = 'infjack' for the infinitesimal jackknife for bagging.

For classification and predict.all = TRUE, a factor levels are returned as numerics. To retrieve the corresponding factor levels, use rf$forest$levels, if rf is the ranger object.

Value

Object of class ranger.prediction with elements
predictions.ranger

predictions uniqueNdeathNtimes chf survival numNtrees numNindependentNvariables treetype numNsamples

Predicted classes/values (only for classification and regression)
Unique death times (only for survival).
Estimated cumulative hazard function for each sample (only for survival).
Estimated survival function for each sample (only for survival).
Number of trees.
Number of independent variables.
Type of forest/tree. Classification, regression or survival.
Number of samples.

Author(s)

Marvin N. Wright

References


See Also

ranger

description

Extract training data predictions of Ranger object.

Usage

## S3 method for class 'ranger'
predictions(x, ...)

Arguments

x Ranger object.
...
Further arguments passed to or from other methods.

Value

Predictions: Classes for Classification forests, Numerical values for Regressions forests and the estimated survival functions for all individuals for Survival forests.
Author(s)
Marvin N. Wright

See Also
ranger
print.ranger

**Print Ranger**

**Description**
Print contents of Ranger object.

**Usage**
```r
## S3 method for class 'ranger'
print(x, \ldots)
```

**Arguments**
- `x`: Object of class `ranger`.
- `\ldots`: Further arguments passed to or from other methods.

**Author(s)**
Marvin N. Wright

**See Also**
- ranger

print.ranger.forest

**Print Ranger forest**

**Description**
Print contents of Ranger forest object.

**Usage**
```r
## S3 method for class 'ranger.forest'
print(x, \ldots)
```

**Arguments**
- `x`: Object of class `ranger.forest`.
- `\ldots`: Further arguments passed to or from other methods.

**Author(s)**
Marvin N. Wright
print.ranger.prediction

*Print Ranger prediction*

**Description**

Print contents of Ranger prediction object.

**Usage**

```r
## S3 method for class 'ranger.prediction'
print(x, ...)
```

**Arguments**

- `x`: Object of class `ranger.prediction`.
- `...`: further arguments passed to or from other methods.

**Author(s)**

Marvin N. Wright

---

*ranger*  

*Ranger*

**Description**

Ranger is a fast implementation of random forests (Breiman 2001) or recursive partitioning, particularly suited for high dimensional data. Classification, regression, and survival forests are supported. Classification and regression forests are implemented as in the original Random Forest (Breiman 2001), survival forests as in Random Survival Forests (Ishwaran et al. 2008). Includes implementations of extremely randomized trees (Geurts et al. 2006) and quantile regression forests (Meinshausen 2006).

**Usage**

```r
ranger(formula = NULL, data = NULL, num.trees = 500, mtry = NULL, importance = "none", write.forest = TRUE, probability = FALSE, min.node.size = NULL, max.depth = NULL, replace = TRUE, sample.fraction = ifelse(replace, 1, 0.632), case.weights = NULL, class.weights = NULL, splitrule = NULL, num.random.splits = 1, alpha = 0.5, minprop = 0.1, split.select.weights = NULL, always.split.variables = NULL, respect.unordered.factors = NULL, scale.permutation.importance = FALSE, keep.inbag = FALSE, inbag = NULL, holdout = FALSE, quantreg = FALSE,)
```
Arguments

**formula**
Object of class `formula` or character describing the model to fit. Interaction terms supported only for numerical variables.

**data**
Training data of class `data.frame`, `matrix`, `dgCMatrix` (Matrix) or `gwaa.data` (GenABEL).

**num.trees**
Number of trees.

**mtry**
Number of variables to possibly split at in each node. Default is the (rounded down) square root of the number variables.

**importance**
Variable importance mode, one of 'none', 'impurity', 'impurity_corrected', 'permutation'. The 'impurity' measure is the Gini index for classification, the variance of the responses for regression and the sum of test statistics (see `splitrule`) for survival.

**write.forest**
Save `ranger.forest` object, required for prediction. Set to `FALSE` to reduce memory usage if no prediction intended.

**probability**
Grow a probability forest as in Malley et al. (2012).

**min.node.size**
Minimal node size. Default 1 for classification, 5 for regression, 3 for survival, and 10 for probability.

**max.depth**
Maximal tree depth. A value of NULL or 0 (the default) corresponds to unlimited depth, 1 to tree stumps (1 split per tree).

**replace**
Sample with replacement.

**sample.fraction**
Fraction of observations to sample. Default is 1 for sampling with replacement and 0.632 for sampling without replacement. For classification, this can be a vector of class-specific values.

**case.weights**
Weights for sampling of training observations. Observations with larger weights will be selected with higher probability in the bootstrap (or subsampled) samples for the trees.

**class.weights**
Weights for the outcome classes (in order of the factor levels) in the splitting rule (cost sensitive learning). Classification and probability prediction only. For classification the weights are also applied in the majority vote in terminal nodes.

**splitrule**
Splitting rule. For classification and probability estimation "gini" or "extratrees" with default "gini". For regression "variance", "extratrees" or "maxstat" with default "variance". For survival "logrank", "extratrees", "C" or "maxstat" with default "logrank".

**num.random.splits**
For "extratrees" `splitrule`. Number of random splits to consider for each candidate splitting variable.

**alpha**
For "maxstat" `splitrule`: Significance threshold to allow splitting.

**minprop**
For "maxstat" `splitrule`: Lower quantile of covariate distribution to be considered for splitting.
split.select.weights
   Numeric vector with weights between 0 and 1, representing the probability to select variables for splitting. Alternatively, a list of size num.trees, containing split select weight vectors for each tree can be used.

always.split.variables
   Character vector with variable names to be always selected in addition to the mtry variables tried for splitting.

respect.unordered.factors
   Handling of unordered factor covariates. One of 'ignore', 'order' and 'partition'. For the "extratrees" splitrule the default is "partition" for all other splitrules 'ignore'. Alternatively TRUE (= 'order') or FALSE (= 'ignore') can be used. See below for details.

scale.permutation.importance
   Scale permutation importance by standard error as in (Breiman 2001). Only applicable if permutation variable importance mode selected.

keep.inbag
   Save how often observations are in-bag in each tree.

inbag
   Manually set observations per tree. List of size num.trees, containing inbag counts for each observation. Can be used for stratified sampling.

holdout
   Hold-out mode. Hold-out all samples with case weight 0 and use these for variable importance and prediction error.

quantreg
   Prepare quantile prediction as in quantile regression forests (Meinshausen 2006). Regression only. Set keep.inbag = TRUE to prepare out-of-bag quantile prediction.

oob.error
   Compute OOB prediction error. Set to FALSE to save computation time, e.g. for large survival forests.

num.threads
   Number of threads. Default is number of CPUs available.

save.memory
   Use memory saving (but slower) splitting mode. No effect for survival and GWAS data. Warning: This option slows down the tree growing, use only if you encounter memory problems.

verbose
   Show computation status and estimated runtime.

seed
   Random seed. Default is NULL, which generates the seed from R. Set to 0 to ignore the R seed.

dependent.variable.name
   Name of dependent variable, needed if no formula given. For survival forests this is the time variable.

status.variable.name
   Name of status variable, only applicable to survival data and needed if no formula given. Use 1 for event and 0 for censoring.

classification
   Only needed if data is a matrix. Set to TRUE to grow a classification forest.

Details

The tree type is determined by the type of the dependent variable. For factors classification trees are grown, for numeric values regression trees and for survival objects survival trees. The Gini index is used as default splitting rule for classification. For regression, the estimated response variances or
maximally selected rank statistics (Wright et al. 2016) can be used. For Survival the log-rank test, a C-index based splitting rule (Schmid et al. 2015) and maximally selected rank statistics (Wright et al. 2016) are available. For all tree types, forests of extremely randomized trees (Geurts et al. 2006) can be grown.

With the probability option and factor dependent variable a probability forest is grown. Here, the node impurity is used for splitting, as in classification forests. Predictions are class probabilities for each sample. In contrast to other implementations, each tree returns a probability estimate and these estimates are averaged for the forest probability estimate. For details see Malley et al. (2012).

Note that for classification and regression nodes with size smaller than \( \text{minNnodeNsize} \) can occur, as in original Random Forests. For survival all nodes contain at \( \text{minNnodeNsize} \) samples. Variables selected with \( \text{alwaysNsplitNvariables} \) are tried additionally to the \( \text{mtry} \) variables randomly selected. In \( \text{splitNselectNweights} \) variables weighted with 0 are never selected and variables with 1 are always selected. Weights do not need to sum up to 1, they will be normalized later. The weights are assigned to the variables in the order they appear in the formula or in the data if no formula is used. Names of the \( \text{splitNselectNweights} \) vector are ignored. The usage of \( \text{splitNselectNweights} \) can increase the computation times for large forests.

Unordered factor covariates can be handled in 3 different ways by using \( \text{respectNunorderedNfactors} \): For ‘ignore’ all factors are regarded ordered, for ‘partition’ all possible 2-partitions are considered for splitting. For ‘order’ and 2-class classification the factor levels are ordered by their proportion falling in the second class, for regression by their mean response, as described in Hastie et al. (2009), chapter 9.2.4. For multiclass classification the factor levels are ordered by the first principal component of the weighted covariance matrix of the contingency table (Coppersmith et al. 1999), for survival by the median survival (or the largest available quantile if the median is not available). The use of ‘order’ is recommended, as it computationally fast and can handle an unlimited number of factor levels. Note that the factors are only reordered once and not again in each split.

The ‘impurity_corrected’ importance measure is unbiased in terms of the number of categories and category frequencies and is almost as fast as the standard impurity importance. It is a modified version of the method by Sandri & Zuccolotto (2008), which is faster and more memory efficient. See Nembrini et al. (2018) for details. This importance measure can be combined with the methods to estimate p-values in \( \text{importanceNpvalues} \).

For a large number of variables and data frames as input data the formula interface can be slow or impossible to use. Alternatively dependent.\( \text{variableNname} \) (and status.\( \text{variableNname} \) for survival) can be used. Consider setting save.\( \text{memory} = \text{TRUE} \) if you encounter memory problems for very large datasets, but be aware that this option slows down the tree growing.

For GWAS data consider combining \text{ranger} with the GenABEL package. See the Examples section below for a demonstration using Plink data. All SNPs in the GenABEL object will be used for splitting. To use only the SNPs without sex or other covariates from the phenotype file, use 0 on the right hand side of the formula. Note that missing values are treated as an extra category while splitting.

See \url{https://github.com/imbs-hl/ranger} for the development version.

With recent R versions, multithreading on Windows platforms should just work. If you compile yourself, the new RT tools toolchain is required.

\textbf{Value}

Object of class \text{ranger} with elements
forest     Saved forest (If write.forest set to TRUE). Note that the variable IDs in the
split.varIDs object do not necessarily represent the column number in R.
predictions Predicted classes/values, based on out of bag samples (classification and regression only).
variable.importance Variable importance for each independent variable.
prediction.error Overall out of bag prediction error. For classification this is the fraction of misclassified samples, for probability estimation the Brier score, for regression the mean squared error and for survival one minus Harrell’s C-index.
r.squared R squared. Also called explained variance or coefficient of determination (regression only). Computed on out of bag data.
confusion.matrix Contingency table for classes and predictions based on out of bag samples (classification only).
unique.death.times Unique death times (survival only).
chf Estimated cumulative hazard function for each sample (survival only).
survival Estimated survival function for each sample (survival only).
call Function call.
num.trees Number of trees.
num.independent.variables Number of independent variables.
mtry Value of mtry used.
min.node.size Value of minimal node size used.
treetype Type of forest/tree. classification, regression or survival.
importance.mode Importance mode used.
num.samples Number of samples.
inbag.counts Number of times the observations are in-bag in the trees.

Author(s)
Marvin N. Wright

References


See Also
predict.ranger

Examples
require(ranger)

## Classification forest with default settings
ranger(Species ~ ., data = iris)

## Prediction
train.idx <- sample(nrow(iris), 2/3 * nrow(iris))
iris.train <- iris[train.idx, ]
iris.test <- iris[-train.idx, ]
rg.iris <- ranger(Species ~ ., data = iris.train)
pred.iris <- predict(rg.iris, data = iris.test)
table(iris.test$Species, pred.iris$predictions)

## Quantile regression forest
rf <- ranger(mpg ~ ., mtcars[1:26, ], quantreg = TRUE)
pred <- predict(rf, mtcars[27:32, ], type = "quantiles")
pred$predictions
### Variable importance

gg.iris <- ranger(Species ~ ., data = iris, importance = "impurity")
gg.iris$variable.importance

### Survival forest

require(survival)
grg.veteran <- ranger(Surv(time, status) ~ ., data = veteran)
plot(rgg.veteran$unique.death.times, gg.veteran$survival[1])

### Alternative interface

ranger(dependent.variable.name = "Species", data = iris)

---

**Description**

Extract unique death times of Ranger Survival forest

**Usage**

```r
## S3 method for class 'ranger'
timepoints(x, ...)
```

**Arguments**

- `x` Ranger Survival forest object.
- `...` Further arguments passed to or from other methods.

**Value**

Unique death times

**Author(s)**

Marvin N. Wright
timepoints.ranger.prediction

Ranger timepoints

Description

Extract unique death times of Ranger Survival prediction object.

Usage

```r
## S3 method for class 'ranger.prediction'
timepoints(x, ...)
```

Arguments

- `x` Ranger Survival prediction object.
- `...` Further arguments passed to or from other methods.

Value

Unique death times

Author(s)

Marvin N. Wright

See Also

ranger

treeInfo

Tree information in human readable format

Description

Extract tree information of a `ranger` object.

Usage

```r
treeInfo(object, tree = 1)
```
Arguments

object ranger object.
tree Number of the tree of interest.

Details

Node and variable ID’s are 0-indexed, i.e., node 0 is the root node. If the formula interface is used in the \texttt{ranger} call, the variable ID’s are usually different to the original data used to grow the tree. Refer to the variable name instead to be sure.

Splitting at unordered factors (nominal variables) depends on the option \texttt{respect.unordered.factors} in the \texttt{ranger} call. For the “ignore” and “order” approaches, all values smaller or equal the \texttt{splitval} value go to the left and all values larger go to the right, as usual. However, with “order” the values correspond to the order in \texttt{object$forest$covariate.levels} instead of the original order (usually alphabetical). In the “partition” mode, the \texttt{splitval} values for unordered factor are comma separated lists of values, representing the factor levels (in the original order) going to the right.

Value

A data.frame with the columns

- \texttt{nodeID} The nodeID, 0-indexed.
- \texttt{leftChild} ID of the left child node, 0-indexed.
- \texttt{rightChild} ID of the right child node, 0-indexed.
- \texttt{splitvarID} ID of the splitting variable, 0-indexed. Caution, the variable order changes if the formula interface is used.
- \texttt{splitvarName} Name of the splitting variable.
- \texttt{splitval} The splitting value. For numeric or ordinal variables, all values smaller or equal go to the left, larger values to the right.
- \texttt{terminal} Logical, TRUE for terminal nodes.
- \texttt{prediction} One column with the predicted class (factor) for classification and the predicted numerical value for regression.

Author(s)

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See Also

\texttt{ranger}

Examples

```r
require(ranger)
rf <- ranger(Species ~ ., data = iris)
treeInfo(rf, 1)
```
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