Package ‘mlr’

January 10, 2020

Title Machine Learning in R
Version 2.17.0
Description Interface to a large number of classification and regression techniques, including machine-readable parameter descriptions. There is also an experimental extension for survival analysis, clustering and general, example-specific cost-sensitive learning. Generic resampling, including cross-validation, bootstrapping and subsampling. Hyperparameter tuning with modern optimization techniques, for single- and multi-objective problems. Filter and wrapper methods for feature selection. Extension of basic learners with additional operations common in machine learning, also allowing for easy nested resampling. Most operations can be parallelized.

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BugReports https://github.com/mlr-org/mlr/issues

Depends ParamHelpers (>= 1.10), R (>= 3.0.2)

Imports backports (>= 1.1.0), BBmisc (>= 1.11), checkmate (>= 1.8.2), data.table (>= 1.12.4), ggplot2, methods, parallelMap (>= 1.3), stats, stringi, survival, utils, XML

Suggests ada, adabag, bartMachine, batchtools, bran, bst, C50, care, caret (>= 6.0-57), class, classiFunc, cluse, cluster, ClusterR, clusterSim (>= 0.44-5), cValid, cmaes, cowplot, CoxBoost, crs, Cubist, deepnet, DiceKriging, DiscrMiner, e1071, earth, elasticnet, emoa, evtree, extraTrees, fda.usc, FDboost, flare, FNN, forecast (>= 8.3), fpc, frbs, FSelector, FSelectorRcpp (>= 0.2.1), gbm, GenSA, ggrepur, glmnet, GPfit, h2o (>= 3.6.0.8), Hmisc, hrbrthemes, irace (>= 2.0), kernlab, kknn, klaR, knitr, laGP, LiblineaR, lintr (>= 1.0.0.9001), MASS, mboost, mco, mda, memoise, mlbench, mldr, mlrMBO, mmpf, modeltools, mRMRe, neuralnet, nnet, nodeHarvest (>= 0.7-3), numDeriv, pamr, pander, party, penalized (>= 0.9-47), pls, PMCMR (>= 4.1), praznik (>= 5.0.0), randomForest, randomForestSRC (>= 2.7.0),
ranger (>= 0.8.0), rappdirs, refund, rex, rFerns, rgenoud, rknn, rmarkdown, ROCR, rotationForest, rpart, RRF, rrlda, rsm, RSNNS, rucrdtw, RWeka, sda, sf, smoof, sparseLDA, stepPlr, survAUC, svglite, SwarmSVM, testthat, tgp, TH.data, tsfeatures, vdiffr, wavelets, xgboost (>= 0.7)

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Interface to a large number of classification and regression techniques, including machine-readable parameter descriptions. There is also an experimental extension for survival analysis, clustering and general, example-specific cost-sensitive learning. Generic resampling, including cross-validation, bootstrapping and subsampling. Hyperparameter tuning with modern optimization techniques, for single- and multi-objective problems. Filter and wrapper methods for feature selection. Extension of basic learners with additional operations common in machine learning, also allowing for easy nested resampling. Most operations can be parallelized.
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See Also

Useful links:
- https://mlr.mlr-org.com
- https://github.com/mlr-org/mlr
- Report bugs at https://github.com/mlr-org/mlr/issues
addRRMeasure  

Compute new measures for existing ResampleResult

Description

Adds new measures to an existing ResampleResult.

Usage

addRRMeasure(res, measures)

Arguments

res  
(ResampleResult)
The result of resample run with keep.pred = TRUE.

measures  
(Measure | list of Measure)
Performance measure(s) to evaluate. Default is the default measure for the task, see here getDefaultMeasure.

Value

(ResampleResult).

See Also

Other resample: ResamplePrediction, ResampleResult, getRRPredictionList(), getRRPredictions(), getRRTaskDescription(), getRRTaskDesc(), makeResampleDesc(), makeResampleInstance(), resample()  

Aggregation  

Aggregation object.

Description

An aggregation method reduces the performance values of the test (and possibly the training sets) to a single value. To see all possible implemented aggregations look at aggregations.

The aggregation can access all relevant information of the result after resampling and combine them into a single value. Though usually something very simple like taking the mean of the test set performances is done.

Object members:

id (character(1)) Name of the aggregation method.

name (character(1)) Long name of the aggregation method.

properties (character) Properties of the aggregation.

fun (‘function(task, perf.test, perf.train, measure, group, pred)’) Aggregation function.
aggregations

See Also

makeAggregation

aggregations

Aggregation methods.

Description

- **test.mean**
  Mean of performance values on test sets.

- **test.sd**
  Standard deviation of performance values on test sets.

- **test.median**
  Median of performance values on test sets.

- **test.min**
  Minimum of performance values on test sets.

- **test.max**
  Maximum of performance values on test sets.

- **test.sum**
  Sum of performance values on test sets.

- **train.mean**
  Mean of performance values on training sets.

- **train.sd**
  Standard deviation of performance values on training sets.

- **train.median**
  Median of performance values on training sets.

- **train.min**
  Minimum of performance values on training sets.

- **train.max**
  Maximum of performance values on training sets.

- **train.sum**
  Sum of performance values on training sets.

- **b632**
  Aggregation for B632 bootstrap.

- **b632plus**
  Aggregation for B632+ bootstrap.

- **testgroup.mean**
  Performance values on test sets are grouped according to resampling method. The mean for every group is calculated, then the mean of those means. Mainly used for repeated CV.

- **testgroup.sd**
  Similar to **testgroup.mean** - after the mean for every group is calculated, the standard deviation of those means is obtained. Mainly used for repeated CV.
• **test.join**
  Performance measure on joined test sets. This is especially useful for small sample sizes where unbalanced group sizes have a significant impact on the aggregation, especially for cross-validation. `test.join` might make sense now. For the repeated CV, the performance is calculated on each repetition and then aggregated with the arithmetic mean.

**Usage**

- `test.mean`
- `test.sd`
- `test.median`
- `test.min`
- `test.max`
- `test.sum`
- `test.range`
- `test.rmse`
- `train.mean`
- `train.sd`
- `train.median`
- `train.min`
- `train.max`
- `train.sum`
- `train.range`
- `train.rmse`
- `b632`
- `b632plus`
- `testgroup.mean`
- `testgroup.sd`
- `test.join`
**agri.task**

**Format**

None

**See Also**

Aggregation

---

**agri.task**  *European Union Agricultural Workforces clustering task.*

**Description**

Contains the task (agri.task).

**References**

See cluster::agriculture.

---

**analyzeFeatSelResult**  *Show and visualize the steps of feature selection.*

**Description**

This function prints the steps selectFeatures took to find its optimal set of features and the reason why it stopped. It can also print information about all calculations done in each intermediate step. Currently only implemented for sequential feature selection.

**Usage**

analyzeFeatSelResult(res, reduce = TRUE)

**Arguments**

- **res**  *(FeatSelResult)*
  The result of of selectFeatures.

- **reduce**  *(logical(1))*
  Per iteration: Print only the selected feature (or all features that were evaluated)? Default is TRUE.

**Value**

(invisible(NULL)).

**See Also**

Other featsel: FeatSelControl, getFeatSelResult(), makeFeatSelWrapper(), selectFeatures()
asROCRPrediction  Converts predictions to a format package ROCR can handle.

Description

Converts predictions to a format package ROCR can handle.

Usage

asROCRPrediction(pred)

Arguments

pred  (Prediction)
      Prediction object.

See Also

Other roc: calculateROCMeasures()
Other predict: getPredictionProbabilities(), getPredictionResponse(), getPredictionTaskDesc(), predict.WrappedModel(), setPredictThreshold(), setPredictType()

batchmark  Run machine learning benchmarks as distributed experiments.

Description

This function is a very parallel version of [benchmark] using batchtools. Experiments are created in the provided registry for each combination of learners, tasks and resamplings. The experiments are then stored in a registry and the runs can be started via [batchtools::submitJobs]. A job is one train/test split of the outer resampling. In case of nested resampling (e.g. with [makeTuneWrapper]), each job is a full run of inner resampling, which can be parallelized in a second step with ParallelMap. For details on the usage and support backends have a look at the batchtools tutorial page: <https://github.com/mlg/batchtools>.

The general workflow with ‘batchmark’ looks like this:

1. Create an ExperimentRegistry using [batchtools::makeExperimentRegistry].
2. Call ‘batchmark(...)’ which defines jobs for all learners and tasks in an [base::expand.grid] fashion.
3. Submit jobs using [batchtools::submitJobs].
4. Babysit the computation, wait for all jobs to finish using [batchtools::waitForJobs].
5. Call ‘reduceBatchmarkResult()’ to reduce results into a [BenchmarkResult].

If you want to use this with OpenML datasets you can generate tasks from a vector of dataset IDs easily with ‘tasks = lapply(data.ids, function(x) convertOMLDataSetToMlr(getOMLDataSet(x)))’.
**Usage**

```r
batchmark(
  learners,
  tasks,
  resamplings,
  measures,
  keep.pred = TRUE,
  keep.extract = FALSE,
  models = FALSE,
  reg = batchtools::getDefaultRegistry()
)
```

**Arguments**

- **learners** (list of Learner | character)
  Learning algorithms which should be compared, can also be a single learner. If you pass strings the learners will be created via `makeLearner`.

- **tasks** list of Task
  Tasks that learners should be run on.

- **resamplings** [(list of) [ResampleDesc]]
  Resampling strategy for each tasks. If only one is provided, it will be replicated to match the number of tasks. If missing, a 10-fold cross validation is used.

- **measures** (list of Measure)
  Performance measures for all tasks. If missing, the default measure of the first task is used.

- **keep.pred** (logical(1))
  Keep the prediction data in the `pred` slot of the result object. If you do many experiments (on larger data sets) these objects might unnecessarily increase object size / mem usage, if you do not really need them. The default is set to `TRUE`.

- **keep.extract** (logical(1))
  Keep the `extract` slot of the result object. When creating a lot of benchmark results with extensive tuning, the resulting R objects can become very large in size. That is why the tuning results stored in the `extract` slot are removed by default (`keep.extract = FALSE`). Note that when `keep.extract = FALSE` you will not be able to conduct analysis in the tuning results.

- **models** (logical(1))
  Should all fitted models be stored in the `ResampleResult`? Default is `FALSE`.

- **reg** ([batchtools::Registry])
  Registry, created by `batchtools::makeExperimentRegistry`. If not explicitly passed, uses the last created registry.

**Value**

([data.table]). Generated job ids are stored in the column “job.id”.
See Also

Other benchmark: BenchmarkResult, benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFatSelResults(), getBMRFilteredFeatures(), getBMRFeatSelResults(), getBMRFeatSelResults(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRLearners(), getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMBoxplots(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

bc.task  
Wisconsin Breast Cancer classification task.

Description

Contains the task (bc.task).

References

See mlbench::BreastCancer. The column "Id" and all incomplete cases have been removed from the task.

benchmark  
Benchmark experiment for multiple learners and tasks.

Description

Complete benchmark experiment to compare different learning algorithms across one or more tasks w.r.t. a given resampling strategy. Experiments are paired, meaning always the same training / test sets are used for the different learners. Furthermore, you can of course pass “enhanced” learners via wrappers, e.g., a learner can be automatically tuned using makeTuneWrapper.

Usage

```
benchmark(
  learners,
  tasks,
  resamplings,
  measures,
  keep.pred = TRUE,
  keep.extract = FALSE,
  models = FALSE,
  show.info = getMlrOption("show.info")
)
```
Arguments

learners (list of Learner | character)
Learning algorithms which should be compared, can also be a single learner. If you pass strings the learners will be created via makeLearner.

tasks list of Task
Tasks that learners should be run on.

resamplings (list of ResampleDesc | ResampleInstance)
Resampling strategy for each tasks. If only one is provided, it will be replicated to match the number of tasks. If missing, a 10-fold cross validation is used.

measures (list of Measure)
Performance measures for all tasks. If missing, the default measure of the first task is used.

keep.pred (logical(1))
Keep the prediction data in the pred slot of the result object. If you do many experiments (on larger data sets) these objects might unnecessarily increase object size / mem usage, if you do not really need them. The default is set to TRUE.

keep.extract (logical(1))
Keep the extract slot of the result object. When creating a lot of benchmark results with extensive tuning, the resulting R objects can become very large in size. That is why the tuning results stored in the extract slot are removed by default (keep.extract = FALSE). Note that when keep.extract = FALSE you will not be able to conduct analysis in the tuning results.

models (logical(1))
Should all fitted models be stored in the ResampleResult? Default is FALSE.

show.info (logical(1))
Print verbose output on console? Default is set via configureMlr.

Value

BenchmarkResult.

See Also

Other benchmark: BenchmarkResult, batchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFitSelResults(), getBMRFittedFeatures(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRLearners(), getBMRMergeIds(), getBMRMerge(), getBMRMerge(), getBMRMerge(), getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMRXboxplots(), plotBMXRanksAsBarChart(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

Examples

lrns = list(makeLearner("classif.lda"), makeLearner("classif.rpart"))
tasks = list(iris.task, sonar.task)
desc = makeResampleDesc("CV", iters = 2L)
meas = list(acc, ber)
bmr = benchmark(lrns, tasks, rdesc, measures = meas)
bh.task

BenchmarkResult

Description

Result of a benchmark experiment conducted by benchmark with the following members:

- **results** (list of ResampleResult): A nested list of resample results, first ordered by task id, then by learner id.
- **measures** (list of Measure): The performance measures used in the benchmark experiment.
- **learners** (list of Learner): The learning algorithms compared in the benchmark experiment.

The print method of this object shows aggregated performance values for all tasks and learners. It is recommended to retrieve required information via the getBMR* getter functions. You can also convert the object using as.data.frame.

See Also

Other benchmark: batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFixedFeatures(), getBMRFeatSelResults(), getBMRFeatSelResults(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRModels(), getBMRLearners(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRModels(), getBMRLearners(), getBMRTaskDescs(), getBMRTaskDescs(), getBMRTuneResults(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

bh.task

Boston Housing regression task.

Description

Contains the task (bh.task).

References

See mlbench::BostonHousing.
cache_helpers

---

**cache_helpers**  
*Get or delete mlr cache directory*

---

**Description**

Helper functions to deal with mlr caching.

**Usage**

- `getCacheDir()`
- `deleteCacheDir()`

**Details**

- `getCacheDir()` returns the default mlr cache directory
- `deleteCacheDir()` clears the default mlr cache directory. Custom cache directories must be deleted by hand.

---

**calculateConfusionMatrix**

*Confusion matrix.*

---

**Description**

Calculates the confusion matrix for a (possibly resampled) prediction. Rows indicate true classes, columns predicted classes. The marginal elements count the number of classification errors for the respective row or column, i.e., the number of errors when you condition on the corresponding true (rows) or predicted (columns) class. The last bottom right element displays the total amount of errors.

A list is returned that contains multiple matrices. If `relative = TRUE` we compute three matrices, one with absolute values and two with relative. The relative confusion matrices are normalized based on rows and columns respectively, if `FALSE` we only compute the absolute value matrix.

The `print` function returns the relative matrices in a compact way so that both row and column marginals can be seen in one matrix. For details see `ConfusionMatrix`.

Note that for resampling no further aggregation is currently performed. All predictions on all test sets are joined to a vector `yhat`, as are all labels joined to a vector `y`. Then `yhat` is simply tabulated vs. `y`, as if both were computed on a single test set. This probably mainly makes sense when cross-validation is used for resampling.

**Usage**

- `calculateConfusionMatrix(pred, relative = FALSE, sums = FALSE, set = "both")`

  ```r
  # S3 method for class 'ConfusionMatrix'
  print(x, both = TRUE, digits = 2, ...)
  ```
Arguments

pred (Prediction)
Prediction object.

relative (logical(1))
If TRUE two additional matrices are calculated. One is normalized by rows and one by columns.

sums (logical(1))
If TRUE add absolute number of observations in each group.

set (character(1))
Specifies which part(s) of the data are used for the calculation. If set equals train or test, the pred object must be the result of a resampling, otherwise an error is thrown. Defaults to “both”. Possible values are “train”, “test”, or “both”.

x (ConfusionMatrix)
Object to print.

both (logical(1))
If TRUE both the absolute and relative confusion matrices are printed.

digits (integer(1))
How many numbers after the decimal point should be printed, only relevant for relative confusion matrices.

... (any)
Currently not used.

Value
(ConfusionMatrix).

Methods (by generic)

• print:

See Also

Other performance: ConfusionMatrix, calculateROCMeasures(), estimateRelativeOverfitting(), makeCostMeasure(), makeCustomResampledMeasure(), makeMeasure(), measures, performance(), setAggregation(), setMeasurePars()

Examples

# get confusion matrix after simple manual prediction
allinds = 1:150
test = sample(allinds, 75)
mod = train("classif.lda", iris.task, subset = train)
pred = predict(mod, iris.task, subset = test)
print(calculateConfusionMatrix(pred))
print(calculateConfusionMatrix(pred, sums = TRUE))
print(calculateConfusionMatrix(pred, relative = TRUE))
# now after cross-validation
r = crossval("classif.lda", iris.task, iters = 2L)
print(calculateConfusionMatrix(r$pred))

---

calculateROCMeasures Calculate receiver operator measures.

Description

Calculate the absolute number of correct/incorrect classifications and the following evaluation measures:

- tpr True positive rate (Sensitivity, Recall)
- fpr False positive rate (Fall-out)
- fnr False negative rate (Miss rate)
- tnr True negative rate (Specificity)
- ppv Positive predictive value (Precision)
- for False omission rate
- 1rp Positive likelihood ratio (LR+)
- fdr False discovery rate
- npv Negative predictive value
- acc Accuracy
- 1rm Negative likelihood ratio (LR-)
- dor Diagnostic odds ratio

For details on the used measures see measures and also [https://en.wikipedia.org/wiki/Receiver_operating_characteristic](https://en.wikipedia.org/wiki/Receiver_operating_characteristic).

The element for the false omission rate in the resulting object is not called for but fomr since for should never be used as a variable name in an object.

Usage

calculateROCMeasures(pred)

## S3 method for class 'ROCMeasures'
print(x, abbreviations = TRUE, digits = 2, ...)
Arguments

pred (Prediction)
Prediction object.

x (ROCMeasures)
Created by calculateROCMeasures.

abbreviations (logical(1))
If TRUE a short paragraph with explanations of the used measures is printed additionally.

digits (integer(1))
Number of digits the measures are rounded to.

... (any)
Currently not used.

Value

(ROCMeasures). A list containing two elements confusion.matrix which is the 2 times 2 confusion matrix of absolute frequencies and measures, a list of the above mentioned measures.

Methods (by generic)

• print:

See Also

Other roc: asROCRPrediction()

Other performance: ConfusionMatrix, calculateConfusionMatrix(), estimateRelativeOverfitting(), makeCostMeasure(), makeCustomResampledMeasure(), makeMeasure(), measures, performance(), setAggregation(), setMeasurePars()

Examples

lrn = makeLearner("classif.rpart", predict.type = "prob")
fit = train(lrn, sonar.task)
pred = predict(fit, task = sonar.task)
calculateROCMeasures(pred)

capLargeValues Convert large/infinite numeric values in a data.frame or task.

Description

Convert numeric entries which large/infinite (absolute) values in a data.frame or task. Only numeric/integer columns are affected.
**Usage**

```r
capLargeValues(
  obj,
  target = character(0L),
  cols = NULL,
  threshold = Inf,
  impute = threshold,
  what = "abs"
)
```

**Arguments**

- `obj` *(data.frame | Task)*
  Input data.

- `target` *(character)*
  Name of the column(s) specifying the response. Target columns will not be capped. Default is `character(0)`.

- `cols` *(character)*
  Which columns to convert. Default is all numeric columns.

- `threshold` *(numeric(1))*
  Threshold for capping. Every entry whose absolute value is equal or larger is converted. Default is `Inf`.

- `impute` *(numeric(1))*
  Replacement value for large entries. Large negative entries are converted to `-impute`. Default is `threshold`.

- `what` *(character(1))*
  What kind of entries are affected? “abs” means `abs(x) > threshold`, “pos” means `abs(x) > threshold & x > 0`, “neg” means `abs(x) > threshold & x < 0`. Default is “abs”.

**Value**

*(data.frame)*

**See Also**

Other eda_and_preprocess: `createDummyFeatures()`, `dropFeatures()`, `mergeSmallFactorLevels()`, `normalizeFeatures()`, `removeConstantFeatures()`, `summarizeColumns()`, `summarizeLevels()`

**Examples**

```r
capLargeValues(iris, threshold = 5, impute = 5)
```
configureMlr  
**Configures the behavior of the package.**

### Description
Configuration is done by setting custom options.
If you do not set an option here, its current value will be kept.
If you call this function with an empty argument list, everything is set to its defaults.

### Usage
```
configureMlr(
  show.info,  
on.learner.error,  
on.learner.warning,  
on.par.without.desc,  
on.par.out.of.bounds,  
on.measure.not.applicable,  
show.learner.output,  
on.error.dump
)
```

### Arguments
- **show.info** (logical(1))
  Some methods of mlr support a `show.info` argument to enable verbose output on the console. This option sets the default value for these arguments. Setting the argument manually in one of these functions will overwrite the default value for that specific function call. Default is `TRUE`.

- **on.learner.error** (character(1))
  What should happen if an error in an underlying learning algorithm is caught:
  - “stop”: R exception is generated.
  - “warn”: A FailureModel will be created, which predicts only NAs and a warning will be generated.
  - “quiet”: Same as “warn” but without the warning.
  Default is “stop”.

- **on.learner.warning** (character(1))
  What should happen if a warning in an underlying learning algorithm is generated:
  - “warn”: The warning is generated as usual.
  - “quiet”: The warning is suppressed.
  Default is “warn”.

on.par.without.desc
(character(1))
What should happen if a parameter of a learner is set to a value, but no parameter
description object exists, indicating a possibly wrong name:
“stop”: R exception is generated.
“warn”: Warning, but parameter is still passed along to learner.
“quiet”: Same as “warn” but without the warning.
Default is “stop”.

on.par.out.of.bounds
(character(1))
What should happen if a parameter of a learner is set to an out of bounds value.
“stop”: R exception is generated.
“warn”: Warning, but parameter is still passed along to learner.
“quiet”: Same as “warn” but without the warning.
Default is “stop”.

on.measure.not.applicable
(logical(1))
What should happen if a measure is not applicable to a learner.
“stop”: R exception is generated.
“warn”: Warning, but value of the measure will be NA.
“quiet”: Same as “warn” but without the warning.
Default is “stop”.

show.learner.output
(logical(1))
Should the output of the learning algorithm during training and prediction be
shown or captured and suppressed? Default is TRUE.

on.error.dump
(logical(1))
Specify whether FailureModel models and failed predictions should contain an
error dump that can be used with debugger to inspect an error. This option is
only effective if on.learner.error is “warn” or “quiet”. If it is TRUE, the dump
can be accessed using getFailureModelDump on the FailureModel, getPredictionDump
on the failed prediction, and getRRDump on resample predictions. Default is FALSE.

Value

(invisible(NULL)).

See Also

Other configure: getMlrOptions()
Description

The result of `calculateConfusionMatrix`.

Object members:

- **result** *(matrix)*: Confusion matrix of absolute values and marginals. Can also contain row and column sums of observations.
- **task.desc** *(TaskDesc)*: Additional information about the task.
- **sums** *(logical(1))*: Flag if marginal sums of observations are calculated.
- **relative** *(logical(1))*: Flag if the relative confusion matrices are calculated.
- **relative.row** *(matrix)*: Confusion matrix of relative values and marginals normalized by row.
- **relative.col** *(matrix)*: Confusion matrix of relative values and marginals normalized by column.
- **relative.error** *(numeric(1))*: Relative error overall.

See Also

Other performance: `calculateConfusionMatrix()`, `calculateROCMeasures()`, `estimateRelativeOverfitting()`, `makeCostMeasure()`, `makeCustomResampledMeasure()`, `makeMeasure()`, `measures`, `performance()`, `setAggregation()`, `setMeasurePars()`

---

**convertBMRToRankMatrix**

Convert BenchmarkResult to a rank-matrix.

Description

Computes a matrix of all the ranks of different algorithms over different datasets (tasks). Ranks are computed from aggregated measures. Smaller ranks imply better methods, so for measures that are minimized, small ranks imply small scores. for measures that are maximized, small ranks imply large scores.

Usage

```r
convertBMRToRankMatrix(
  bmr,
  measure = NULL,
  ties.method = "average",
  aggregation = "default"
)
```
convertMLBenchObjToTask

Convert a machine learning benchmark / demo object from package mlbench to a task.

Description

We auto-set the target column, drop any column which is called “Id” and convert logicals to factors.

Usage

convertMLBenchObjToTask(x, n = 100L, ...)

Arguments

- **bmr** (BenchmarkResult): Benchmark result.
- **measure** (Measure): Performance measure. Default is the first measure used in the benchmark experiment.
- **ties.method** (character(1)): See base::rank for details.
- **aggregation** (character(1)): “mean” or “default”. See getBMRAggrPerformances for details on “default”.

Value

(matrix) with measure ranks as entries. The matrix has one row for each learner, and one column for each task.

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMR LearnerIds(), getBMR LearnerShortNames(), getBMR Learners(), getBMR MeasureIds(), getBMR Measures(), getBMR Models(), getBMR Performances(), getBMR Predictions(), getBMR TaskDescs(), getBMR TaskIds(), getBMR Tune Results(), plotBMR Boxplots(), plotBMR RanksAsBarChart(), plotBMR Summary(), plot Crit Differences(), reduceBatchmarkResults()
Arguments

- **x**  
  (character(1))  
  Name of an mlbench function or dataset.

- **n**  
  (integer(1))  
  Number of observations for data simul functions. Note that for a few mlbench function this setting is not exactly respected by mlbench. Default is 100.

- **...**  
  (any)  
  Passed on to data simul functions.

Examples

```r
print(convertMLBenchObjToTask("Ionosphere"))
print(convertMLBenchObjToTask("mlbench.spirals", n = 100, sd = 0.1))
```

costiris.task  
Iris cost-sensitive classification task.

Description

Contains the task (costiris.task).

References

See `datasets::iris`. The cost matrix was generated artificially following  

creatDummyFeatures  
Generate dummy variables for factor features.

Description

Replace all factor features with their dummy variables. Internally `model.matrix` is used. Non factor features will be left untouched and passed to the result.

Usage

```r
createDummyFeatures(
  obj,
  target = character(0L),
  method = "1-of-n",
  cols = NULL
)
```
createSpatialResamplingPlots

**Arguments**

- **obj** *(data.frame | Task)*
  Input data.

- **target** *(character(1) | character(2) | character(n.classes))*
  Name(s) of the target variable(s). Only used when `obj` is a data.frame, otherwise ignored. If survival analysis is applicable, these are the names of the survival time and event columns, so it has length 2. For multilabel classification these are the names of logical columns that indicate whether a class label is present and the number of target variables corresponds to the number of classes.

- **method** *(character(1))*
  Available are:
  "1-of-n": For n factor levels there will be n dummy variables.
  "reference": There will be n-1 dummy variables leaving out the first factor level of each variable.
  Default is “1-of-n”.

- **cols** *(character)*
  Columns to create dummy features for. Default is to use all columns.

**Value**

`data.frame | Task`. Same type as `obj`.

**See Also**

Other eda_and_preprocess: `capLargeValues()`, `dropFeatures()`, `mergeSmallFactorLevels()`, `normalizeFeatures()`, `removeConstantFeatures()`, `summarizeColumns()`, `summarizeLevels()`

---

**createSpatialResamplingPlots**

Create (spatial) resampling plot objects.

**Description**

Visualize partitioning of resample objects with spatial information.

**Usage**

```r
cREATE spatialResamplingPlots(
  task = NULL,
  resample = NULL,
  crs = NULL,
  datum = 4326,
  repetitions = 1,
  color.train = "#0072B5",
  color.test = "#E18727",
)```
point.size = 0.5,
axis.text.size = 14,
x.axis.breaks = waiver(),
y.axis.breaks = waiver()
)

Arguments

task Task
Task object.

resample ResampleResult or named list with (multiple) ResampleResult
As returned by resample.

crs integer
Coordinate reference system (EPSG code number) for the supplied coordinates
in the Task.

datum integer
Coordinate reference system which should be used in the resulting map.

repetitions integer
Number of repetitions.

color.train character
Color for train set.

color.test character
Color for test set.

point.size integer
Point size.

axis.text.size integer
Font size of axis labels.

x.axis.breaks numeric
Custom x axis breaks

y.axis.breaks numeric
Custom y axis breaks

Details

If a named list is given to resample, names will appear in the title of each fold. If multiple inputs
are given to resample, these must be named.

This function makes a hard cut at five columns of the resulting gridded plot. This means if the
resample object consists of folds > 5, these folds will be put into the new row.

For file saving, we recommend to use cowplot::save_plot.

When viewing the resulting plot in RStudio, margins may appear to be different than they really
are. Make sure to save the file to disk and inspect the image.

When modifying axis breaks, negative values need to be used if the area is located in either the
western or southern hemisphere. Use positive values for the northern and eastern hemisphere.
Value

(list of 2L containing (1) multiple ‘gg’ objects and (2) their corresponding labels.

CRS

The crs has to be suitable for the coordinates stored in the Task. For example, if the coordinates are UTM, crs should be set to a UTM projection. Due to a limited axis space in the resulting grid (especially on the x-axis), the data will by default projected into a lat/lon projection, specifically EPSG 4326. If other projections are desired for the resulting map, please set argument datum accordingly. This argument will be passed onto ggplot2::coord_sf.

Author(s)

Patrick Schratz

See Also

Other plot: plotBMRBoxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCalibration(), plotCritDifferences(), plotLearningCurve(), plotPartialDependence(), plotROCCurves(), plotResiduals(), plotThreshVsPerf()

Examples

```r
cr = makeResampleDesc("SpRepCV", folds = 5, reps = 4)
r = resample(makeLearner("classif.qda"), spatial.task, cr)

# single unnamed resample input with 5 folds and 2 repetitions
plots = createSpatialResamplingPlots(spatial.task, r, crs = 32717,
  repetitions = 2, x.axis.breaks = c(-79.065, -79.085),
  y.axis.breaks = c(-3.970, -4))
cowplot::plot_grid(plotlist = plots[["Plots"]], ncol = 5, nrow = 2,
  labels = plots[["Labels"]])
```

```
# single named resample input with 5 folds and 1 repetition and 32717 datum
plots = createSpatialResamplingPlots(spatial.task, list("Resamp" = r),
  crs = 32717, datum = 32717, repetitions = 1)
cowplot::plot_grid(plotlist = plots[["Plots"]], ncol = 5, nrow = 1,
  labels = plots[["Labels"]])
```

```
# multiple named resample inputs with 5 folds and 1 repetition
r = makeResampleDesc("SpRepCV", folds = 5, reps = 4)
```
r1 = resample(makeLearner("classif.qda"), spatial.task, rdesc1)
rdesc2 = makeResampleDesc("RepCV", folds = 5, reps = 4)
r2 = resample(makeLearner("classif.qda"), spatial.task, rdesc2)

plots = createSpatialResamplingPlots(spatial.task,
    list("SpRepCV" = r1, "RepCV" = r2), crs = 32717, repetitions = 1,
    x.axis.breaks = c(-79.055, -79.085), y.axis.breaks = c(-3.975, -4))
cowplot::plot_grid(plotlist = plots[["Plots"]], ncol = 5, nrow = 2,
    labels = plots[["Labels"]])

## Complex arrangements of multiple named resample inputs with 5 folds and 1 repetition
##-------------------------------------------------------------------------------------

p1 = plot_grid(plist[["Plots"]][[1]], plist[["Plots"]][[2]],
    plist[["Plots"]][[3]], ncol = 3, nrow = 1, labels = plist[["Labels"]][1:3],
    label_size = 18)
p12 = plot_grid(plist[["Plots"]][[4]], plist[["Plots"]][[5]], ncol = 2,
    nrow = 1, labels = plist[["Labels"]][4:5], label_size = 18)
p2 = plot_grid(plist[["Plots"]][[6]], plist[["Plots"]][[7]],
    plist[["Plots"]][[8]], ncol = 3, nrow = 1, labels = plist[["Labels"]][6:8],
    label_size = 18)
p22 = plot_grid(plist[["Plots"]][[9]], plist[["Plots"]][[10]], ncol = 2,
    nrow = 1, labels = plist[["Labels"]][9:10], label_size = 18)
cowplot::plot_grid(p1, p12, p2, p22, ncol = 1)

crossover

---

**Crossover.**

**Description**

Takes two bit strings and creates a new one of the same size by selecting the items from the first string or the second, based on a given rate (the probability of choosing an element from the first string).

**Arguments**

- **x** (logical)
  First parent string.

- **y** (logical)
  Second parent string.

- **rate** (numeric(1))
  A number representing the probability of selecting an element of the first string. Default is 0.5.
**downsample**

Downsample (subsample) a task or a data.frame.

**Description**

Decrease the observations in a task or a ResampleInstance to a given percentage of observations.

**Usage**

downsample(obj, perc = 1, stratify = FALSE)

**Arguments**

- **obj** *(Task | ResampleInstance)*
  Input data or a ResampleInstance.

- **perc** *(numeric(1))*
  Percentage from (0, 1). Default is 1.

- **stratify** *(logical(1))*
  Only for classification: Should the downsampled data be stratified according to the target classes? Default is FALSE.

**Value**

([data.frame | Task] | [ResampleInstance]). Same type as obj.'

**See Also**

- makeResampleInstance
- Other downsample: makeDownsampleWrapper()

---

**dropFeatures**

Drop some features of task.

**Description**

Drop some features of task.

**Usage**

dropFeatures(task, features)
**estimateRelativeOverfitting**

**Arguments**

- **task** *(Task)*
  The task.

- **features** *(character)*
  Features to drop.

**Value**

*Task.*

**See Also**

Other eda_and_preprocess: `capLargeValues()`, `createDummyFeatures()`, `mergeSmallFactorLevels()`, `normalizeFeatures()`, `removeConstantFeatures()`, `summarizeColumns()`, `summarizeLevels()`

---

**estimateRelativeOverfitting**

_Estimate relative overfitting._

**Description**

Estimates the relative overfitting of a model as the ratio of the difference in test and train performance to the difference of test performance in the no-information case and train performance. In the no-information case the features carry no information with respect to the prediction. This is simulated by permuting features and predictions.

**Usage**

```r
estimateRelativeOverfitting(
  predish,
  measures,
  task,
  learner = NULL,
  pred.train = NULL,
  iter = 1
)
```

**Arguments**

- **predish** *(ResampleDesc | ResamplePrediction | Prediction)*
  Resampling strategy or resampling prediction or test predictions.

- **measures** *(Measure | list of Measure)*
  Performance measure(s) to evaluate. Default is the default measure for the task, see here `getDefaultMeasure`.

- **task** *(Task)*
  The task.
learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.
pred.train (Prediction)
Training predictions. Only needed if test predictions are passed.
iter (integer)
Iteration number. Default 1, usually you don’t need to specify this. Only needed if test predictions are passed.

Details

Currently only support for classification and regression tasks is implemented.

Value

(data.frame). Relative overfitting estimate(s), named by measure(s), for each resampling iteration.

References


See Also

Other performance: ConfusionMatrix, calculateConfusionMatrix(), calculateROCMeasures(), makeCostMeasure(), makeCustomResampledMeasure(), makeMeasure(), measures, performance(), setAggregation(), setMeasurePars()

Examples

task = makeClassifTask(data = iris, target = "Species")
rdesc = makeResampleDesc("CV", iters = 2)
estimateRelativeOverfitting(rdesc, acc, task, makeLearner("classif.knn"))
estimateRelativeOverfitting(rdesc, acc, task, makeLearner("classif.lda"))
rpred = resample("classif.knn", task, rdesc)$pred
estimateRelativeOverfitting(rpred, acc, task)

estimateResidualVariance

Estimate the residual variance.

Description

Estimate the residual variance of a regression model on a given task. If a regression learner is provided instead of a model, the model is trained (see train) first.

Usage

estimateResidualVariance(x, task, data, target)
extractFDABsignal

Arguments

- **x** (Learner or WrappedModel)
  Learner or wrapped model.
- **task** (RegrTask)
  Regression task. If missing, data and target must be supplied.
- **data** (data.frame)
  A data frame containing the features and target variable. If missing, task must be supplied.
- **target** (character(1))
  Name of the target variable. If missing, task must be supplied.

**extractFDABsignal**  
*Bspline mlq features*

**Description**

The function extracts features from functional data based on the Bspline fit. For more details refer to `FDboost::bsignal()`.

**Usage**

```r
eextractFDABsignal(bsignal.knots = 10L, bsignal.df = 3)
```

**Arguments**

- **bsignal.knots** (integer(1))  
The number of knots for bspline.
- **bsignal.df** (numeric(1))  
The effective degree of freedom of penalized bspline.

**Value**

(data.frame).

**See Also**

Other fda_featextractor: `extractFDADTWKernel()`, `extractFDAFPCA()`, `extractFDAFourier()`, `extractFDAMultiResFeatures()`, `extractFDATsfeatures()`, `extractFDAWavelets()`
**extractFDADTWKernel**  
*DTW kernel features*

**Description**

The function extracts features from functional data based on the DTW distance with a reference dataframe.

**Usage**

```r
extractFDADTWKernel(
  ref.method = "random",
  n.refs = 0.05,
  refs = NULL,
  dtwwindow = 0.05
)
```

**Arguments**

- `ref.method` (character(1))
  How should the reference curves be obtained? Method random draws `n.refs` random reference curves, while all uses all curves as references. In order to use user-provided reference curves, this parameter is set to fixed.

- `n.refs` (numeric(1))
  Number of reference curves to be drawn (as a fraction of the number of observations in the training data).

- `refs` (matrix|integer(n))
  Integer vector of training set row indices or a matrix of reference curves with the same length as the functionals in the training data. Overwrites `ref.method` and `n.refs`.

- `dtwwindow` (numeric(1))
  Size of the warping window size (as a proportion of query length).

**Value**

(data.frame).

**See Also**

Other fda_featextractor: `extractFDABsignal()`, `extractFDAFPCA()`, `extractFDAFourier()`, `extractFDAMultiResFeatures()`, `extractFDATsFeatures()`, `extractFDAWavelets()`
extractFDAFeatures  
*Extract features from functional data.*

**Description**

Extract non-functional features from functional features using various methods.

The function `extractFDAFeatures` performs the extraction for all functional features via the methods specified in `feat.methods` and transforms all mentioned functional (matrix) features into regular data.frame columns. Additionally, a “extractFDAFeatDesc” object which contains learned coefficients and other helpful data for re-extraction during the predict-phase is returned. This can be used with `reextractFDAFeatures` in order to extract features during the prediction phase.

**Usage**

```r
extractFDAFeatures(obj, target = character(0L), feat.methods = list(), ...)
```

**Arguments**

- `obj` (Task | data.frame)
  Task or data.frame to extract functional features from. Must contain functional features as matrix columns.
- `target` (character(1))
  Task target column. Only neccessary for data.frames Default is character(0).
- `feat.methods` (named list)
  List of functional features along with the desired methods for each functional feature. “all” applies the `extractFDAFeatures` method to each functional feature. Names of `feat.methods` must match column names of functional features. Available feature extraction methods are available under family `fda_featextractor`. Specifying a functional feature multiple times with different extraction methods allows for the extraction of different features from the same functional. Default is `list()` which does nothing.
- `...` (any)
  Further hyperparameters passed on to the `feat.methods` specified above.

**Details**

The description object contains these slots:

- **target character**: See argument.
- **coln character**: Colum names of data.
- **fd.cols character**: Functional feature names.
- **extractFDAFeat list**: Contains `feature.methods` and relevant parameters for reextraction.
Value

(list)
data|task (data.frame | Task)
   Extracted features, same type as obj.
desc (extracFDAFeatDesc)
   Description object. See description for details.

See Also

Other fda: makeExtractFDAFeatMethod(), makeExtractFDAFeatsWrapper()

Examples

df = data.frame(x = matrix(rnorm(24), ncol = 8), y = factor(c("a", "a", "b")))
fdf = makeFunctionalData(df, fd.features = list(x1 = 1:4, x2 = 5:8), exclude.cols = "y")
task = makeClassifTask(data = fdf, target = "y")
extracted = extractFDAFeatures(task,
   feat.methods = list("x1" = extractFDAFourier(), "x2" = extractFDAWavelets(filter = "haar")))
print(extracted$task)
reextractFDAFeatures(task, extracted$desc)

extractFDAFourier

Fast Fourier transform features.

Description

The function extracts features from functional data based on the fast fourier transform. For more details refer to stats::fft.

Usage

extractFDAFourier(trafo.coeff = "phase")

Arguments

trafo.coeff (character(1))
   Specifies which transformation of the complex frequency domain representation should be calculated as a feature representation. Must be one of “amplitude” or “phase”. Default is “phase”. The phase shift is returned in Rad, i.e. values lie in [-180, 180].

Value

(data.frame).

See Also

Other fda_featextractor: extractFDABsignal(), extractFDADTWKernel(), extractFDAFPCA(), extractFDAMultiResFeatures(), extractFDATsfeatures(), extractFDAWavelets()
extractFDAFPCA  
*Extract functional principal component analysis features.*

**Description**

The function extracts the functional principal components from a data.frame containing functional features. Uses `stats::prcomp`.

**Usage**

```r
extractFDAFPCA(rank. = NULL, center = TRUE, scale. = FALSE)
```

**Arguments**

- `rank.`: (integer(1))
  Number of principal components to extract. Default is `NULL`.
- `center`: (logical(1))
  Should data be centered before applying PCA?
- `scale.`: (logical(1))
  Should data be scaled before applying PCA?

**Value**

`(data.frame)`.

**See Also**

Other `fda_featextractor`: `extractFDABsignal()`, `extractFDADTWKernel()`, `extractFDAFourier()`, `extractFDAMultiResFeatures()`, `extractFDATsfeatures()`, `extractFDAWavelets()`

extractFDAMultiResFeatures  
*Multiresolution feature extraction.*

**Description**

The function extracts currently the mean of multiple segments of each curve and stacks them as features. The segments length are set in a hierarchy way so the features cover different resolution levels.

**Usage**

```r
extractFDAMultiResFeatures(res.level = 3L, shift = 0.5, seg.lens = NULL)
```
extractFDATsfeatures

Arguments

- `res.level` (integer(1))
  The number of resolution hierarchy, each length is divided by a factor of 2.
- `shift` (numeric(1))
  The overlapping proportion when slide the window for one step.
- `seg.lens` (integer(1))
  Curve subsequence lengths. Needs to sum up to the length of the functional.

Value

(data.frame).

See Also

Other fda_features: `extractFDABsignal()`, `extractFDADTWKernel()`, `extractFDAOscillation()`, `extractFDATsfeatures()`, `extractFDAWavelets()`

---

extractFDATsfeatures  Time-Series Feature Heuristics

Description

The function extracts features from functional data based on known Heuristics. For more details refer to `tsfeatures::tsfeatures()`. Under the hood this function uses the package `tsfeatures::tsfeatures()`. For more information see Hyndman, Wang and Laptev, Large-Scale Unusual Time Series Detection, ICDM 2015.

Note: Currently computes the following features:

Usage

```r
extractFDATsfeatures(
  scale = TRUE,
  trim = FALSE,
  trim_amount = 0.1,
  parallel = FALSE,
  na.action = na.pass,
  feats = NULL,
  ...
)
```
extractFDAWavelets

Discrete Wavelet transform features.

Description

The function extracts discrete wavelet transform coefficients from the raw functional data. See wavelets::dwt for more information.

Arguments

- **scale** (logical(1))
  If TRUE, time series are scaled to mean 0 and sd 1 before features are computed.

- **trim** (logical(1))
  If TRUE, time series are trimmed by trim_amount before features are computed. Values larger than trim_amount in absolute value are set to NA.

- **trim_amount** (numeric(1))
  Default level of trimming if trim==TRUE.

- **parallel** (logical(1))
  If TRUE, multiple cores (or multiple sessions) will be used. This only speeds things up when there are a large number of time series.

- **na.action** (logical(1))
  A function to handle missing values. Use na.interp to estimate missing values.

- **feats** (character)
  A character vector of function names to apply to each time-series in order to extract features.
  Default:

- **...** (any)
  Further arguments passed on to the respective tsfeatures functions.

Value

(data.frame)

References


See Also

Other fda_featextractor: extractFDABsignal(), extractFDADTWKernel(), extractFDAFPCA(), extractFDAFourier(), extractFDAMultiResFeatures(), extractFDAWavelets()
Usage

extractFDAWavelets(filter = "la8", boundary = "periodic")

Arguments

filter (character(1))
Specifies which filter should be used. Must be one of dlla1|b1c followed by an even number for the level of the filter. The level of the filter needs to be smaller or equal then the time-series length. For more information and acceptable filters see help(wt.filter). Defaults to la8.

boundary (character(1))
Boundary to be used. “periodic” assumes circular time series, for “reflection” the series is extended to twice its length. Default is “periodic”.

Value

(data.frame).

See Also

Other fda_featextractor: extractFDABsignal(), extractFDADTWKernel(), extractFDAFPCA(), extractFDAFourier(), extractFDAMultiResFeatures(), extractFDATsfeatures()
Examples

```r
configureMlr(on.learner.error = "warn")
data = iris
data$newfeat = 1 # will make LDA crash
task = makeClassifTask(data = data, target = "Species")
m = train("classif.lda", task) # LDA crashed, but mlr catches this
print(m)
print(m$learner.model) # the error message
p = predict(m, task) # this will predict NAs
print(p)
print(performance(p))
configureMlr(on.learner.error = "stop")
```

FeatSelControl

Create control structures for feature selection.

Description

Feature selection method used by selectFeatures. The following optimization algorithms are available:

- **FeatSelControlExhaustive** Exhaustive search. All feature sets (up to a certain number of features max.features) are searched.

- **FeatSelControlRandom** Random search. Features vectors are randomly drawn, up to a certain number of features max.features. A feature is included in the current set with probability prob. So we are basically drawing (0,1)-membership-vectors, where each element is Bernoulli(prob) distributed.

- **FeatSelControlSequential** Deterministic forward or backward search. That means extending (forward) or shrinking (backward) a feature set. Depending on the given method different approaches are taken.
  - sfs Sequential Forward Search: Starting from an empty model, in each step the feature increasing the performance measure the most is added to the model.
  - sbs Sequential Backward Search: Starting from a model with all features, in each step the feature decreasing the performance measure the least is removed from the model.
  - sffs Sequential Floating Forward Search: Starting from an empty model, in each step the algorithm chooses the best model from all models with one additional feature and from all models with one feature less.
  - sfbs Sequential Floating Backward Search: Similar to sffs but starting with a full model.

- **FeatSelControlGA** Search via genetic algorithm. The GA is a simple (mu, lambda) or (mu + lambda) algorithm, depending on the comma setting. A comma strategy selects a new population of size mu out of the lambda > mu offspring. A plus strategy uses the joint pool of mu parents and lambda offspring for selecting mu new candidates. Out of those mu features, the new lambda features are generated by randomly choosing pairs of parents. These are crossed over and crossover_rate represents the probability of choosing a feature from the first parent instead of the second parent. The resulting offspring is mutated, i.e., its bits are flipped.
with probability `mutation.rate`. If `max.features` is set, offspring are repeatedly generated until the setting is satisfied.

Usage

```r
makeFeatSelControlExhaustive(
  same.resampling.instance = TRUE,
  maxit = NA_integer_,
  max.features = NA_integer_,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default"
)
```

```r
makeFeatSelControlGA(
  same.resampling.instance = TRUE,
  impute.val = NULL,
  maxit = NA_integer_,
  max.features = NA_integer_,
  comma = FALSE,
  mu = 10L,
  lambda,
  crossover.rate = 0.5,
  mutation.rate = 0.05,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default"
)
```

```r
makeFeatSelControlRandom(
  same.resampling.instance = TRUE,
  maxit = 100L,
  max.features = NA_integer_,
  prob = 0.5,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default"
)
```

```r
makeFeatSelControlSequential(
  same.resampling.instance = TRUE,
  impute.val = NULL,
  method,
  alpha = 0.01,
  beta = -0.001,
  maxit = NA_integer_,
  max.features = NA_integer_,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
```
log.fun = "default"
)

Arguments

same.resampling.instance
  (logical(1))
  Should the same resampling instance be used for all evaluations to reduce variance? Default is TRUE.

maxit
  (integer(1))
  Maximal number of iterations. Note, that this is usually not equal to the number of function evaluations.

max.features
  (integer(1))
  Maximal number of features.

tune.threshold
  (logical(1))
  Should the threshold be tuned for the measure at hand, after each feature set evaluation, via tuneThreshold? Only works for classification if the predict type is "prob". Default is FALSE.

tune.threshold.args
  (list)
  Further arguments for threshold tuning that are passed down to tuneThreshold. Default is none.

log.fun
  (function | character(1))
  Function used for logging. If set to "default" (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to "memory" the memory usage for each evaluation will also be displayed, with character(1) small increase in run time. Otherwise character(1) function with arguments learner, resampling, measures, par.set, control, opt_path, dob, x, y, remove.nas, stage and prev.stage is expected. The default displays the performance measures, the time needed for evaluating, the currently used memory and the max memory ever used before (the latter two both taken from go). See the implementation for details.

impute.val
  (numeric)
  If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or Inf instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.

comma
  (logical(1))
  Parameter of the GA feature selection, indicating whether to use a (mu, lambda) or (mu + lambda) GA. The default is FALSE.

mu
  (integer(1))
  Parameter of the GA feature selection. Size of the parent population.
lambda (integer(1))
Parameter of the GA feature selection. Size of the children population (should be smaller or equal to mu).

crossover.rate (numeric(1))
Parameter of the GA feature selection. Probability of choosing a bit from the first parent within the crossover mutation.

mutation.rate (numeric(1))
Parameter of the GA feature selection. Probability of flipping a feature bit, i.e. switch between selecting / deselecting a feature.

prob (numeric(1))

method (character(1))
Parameter of the sequential feature selection. A character representing the method. Possible values are sfs (forward search), sbs (backward search), sffs (floating forward search) and sfbs (floating backward search).

alpha (numeric(1))
Parameter of the sequential feature selection. Minimal required value of improvement difference for a forward / adding step. Default is 0.01.

beta (numeric(1))
Parameter of the sequential feature selection. Minimal required value of improvement difference for a backward / removing step. Negative values imply that you allow a slight decrease for the removal of a feature. Default is -0.001.

Value
(FeatSelControl). The specific subclass is one of FeatSelControlExhaustive, FeatSelControlRandom, FeatSelControlSequential, FeatSelControlGA.

References

See Also
Other featsel: analyzeFeatSelResult(), getFeatSelResult(), makeFeatSelWrapper(), selectFeatures()
Details

Object members:

- **learner** (**Learner**): Learner that was optimized.
- **control** (**FeatSelControl**): Control object from feature selection.
- **x** (**character**): Vector of feature names identified as optimal.
- **y** (**numeric**): Performance values for optimal \( x \).
- **threshold** (**numeric**): Vector of finally found and used thresholds if tune.threshold was enabled in **FeatSelControl**, otherwise not present and hence **NULL**.
- **opt.path** (**ParamHelpers::OptPath**): Optimization path which lead to \( x \).

---

**filterFeatures**  
*Filter features by thresholding filter values.*

Description

First, calls **generateFilterValuesData**. Features are then selected via **select** and **val**.

Usage

```r
filterFeatures(task,
  method = "randomForestSRC_importance",
  fval = NULL,
  perc = NULL,
  abs = NULL,
  threshold = NULL,
  fun = NULL,
  fun.args = NULL,
  mandatory.feat = NULL,
  select.method = NULL,
  base.methods = NULL,
  cache = FALSE,
  ...
)
```

Arguments

- **task** (**Task**): The task.
- **method** (**character(1)**): See **listFilterMethods**. Default is “randomForestSRC_importance”.
- **fval** (**FilterValues**): Result of **generateFilterValuesData**. If you pass this, the filter values in the object are used for feature filtering. **method** and ... are ignored then. Default is **NULL** and not used.
filterFeatures

perc (numeric(1))
If set, select \( \text{perc} \times 100 \) top scoring features. \( \text{perc} = 1 \) means to select all features. Mutually exclusive with arguments abs, threshold and fun.

abs (numeric(1))
If set, select abs top scoring features. Mutually exclusive with arguments perc, threshold and fun.

threshold (numeric(1))
If set, select features whose score exceeds threshold. Mutually exclusive with arguments perc, abs and fun.

fun (function)
If set, select features via a custom thresholding function, which must return the number of top scoring features to select. Mutually exclusive with arguments perc, abs and threshold.

fun.args (any)
Arguments passed to the custom thresholding function.

mandatory.feat (character)
Mandatory features which are always included regardless of their scores

select.method
If multiple methods are supplied in argument method, specify the method that is used for the final subsetting.

base.methods
If method is an ensemble filter, specify the base filter methods which the ensemble method will use.

cache (character(1) | logical)
Whether to use caching during filter value creation. See details.

... (any)
Passed down to selected filter method.

Value

Task.

Caching

If cache = TRUE, the default mlr cache directory is used to cache filter values. The directory is operating system dependent and can be checked with getCacheDir(). The default cache can be cleared with deleteCacheDir(). Alternatively, a custom directory can be passed to store the cache.

Note that caching is not thread safe. It will work for parallel computation on many systems, but there is no guarantee.

Simple and ensemble filters

Besides passing (multiple) simple filter methods you can also pass an ensemble filter method (in a list). The ensemble method will use the simple methods to calculate its ranking. See listFilterEnsembleMethods() for available ensemble methods.
friedmanPostHocTestBMR

Perform a posthoc Friedman-Nemenyi test.

Description

Performs a [PMCMR::posthoc.friedman.nemenyi.test] for a [BenchmarkResult] and a selected measure. This means *all pairwise comparisons* of ‘learners’ are performed. The null hypothesis of the post hoc test is that each pair of learners is equal. If the null hypothesis of the included ad hoc [stats::friedman.test] can be rejected an object of class ‘pairwise.htest’ is returned. If not, the function returns the corresponding friedman.test. Note that benchmark results for at least two learners on at least two tasks are required.

Usage

friedmanPostHocTestBMR(
  bmr,
  measure = NULL,
  p.value = 0.05,
  aggregation = "default"
)

Arguments

  bmr (BenchmarkResult)
  Benchmark result.

  measure (Measure)
  Performance measure. Default is the first measure used in the benchmark experiment.

  p.value (numeric(1))
  p-value for the tests. Default: 0.05

  aggregation (character(1))
  “mean” or “default”. See getBMRAggrPerformances for details on “default”.

Examples

# simple filter
filterFeatures(iris.task, method = "FSelectorRcpp_gain.ratio", abs = 2)

# ensemble filter
filterFeatures(iris.task, method = "E-min",
  base.methods = c("FSelectorRcpp_gain.ratio",
                    "FSelectorRcpp_information.gain"), abs = 2)
Value

([pairwise.htest]): See [PMCMR::posthoc.friedman.nemenyi.test] for details. Additionally two components are added to the list:

f.rejnull ('logical(1)') Whether the according friedman.test rejects the Null hypothesis at the selected p.value

crit.difference ('list(2)') Minimal difference the mean ranks of two learners need to have in order to be significantly different

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMR LearnerIds(), getBMR Learner ShortNames(), getBMR Learners(), getBMRMeasureIds(), getBMR Measures(), getBMR Models(), getBMR Performances(), getBMR Predictions(), getBMR TaskDescs(), getBMR TaskIds(), getBMR TuneResults(), plotBMR Boxplots(), plotBMR Ranks As Bar Chart(), plotBMR Summary(), plotCrit Differences(), reduceBatchmarkResults()

Examples

# see benchmark

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**friedmanTestBMR**

Perform overall Friedman test for a BenchmarkResult.

**Description**

Performs a stats::friedman.test for a selected measure. The null hypothesis is that apart from an effect of the different (Task), the location parameter (aggregated performance measure) is the same for each Learner. Note that benchmark results for at least two learners on at least two tasks are required.

**Usage**

friedmanTestBMR(bmr, measure = NULL, aggregation = "default")

**Arguments**

- **bmr** (BenchmarkResult)
  Benchmark result.
- **measure** (Measure)
  Performance measure. Default is the first measure used in the benchmark experiment.
- **aggregation** (character(1))
  “mean” or “default”. See getBMRAggrPerformances for details on “default”. 
generateCalibrationData

Value
(h.test): See stats::friedman.test for details.

See Also
Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(),
friedmanPostHocTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(),
generateCritDifferencesData(), getBMRFeatSelResults(), getBMRFILTERedFeatures(), getBMRFeatureIds(), getBMRFeatureShortNames(),
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generateCritDifferencesData(), getBMRFeatSelResults(), getBMRFeatureIds(), getBMRFeatureShortNames(),
gen
Usage

generateCalibrationData(obj, breaks = "Sturges", groups = NULL, task.id = NULL)

Arguments

obj (list of Prediction | list of ResampleResult | BenchmarkResult)
Single prediction object, list of them, single resample result, list of them, or a
benchmark result. In case of a list probably produced by different learners you
want to compare, then name the list with the names you want to see in the plots,
probably learner shortnames or ids.

breaks (character(1) | numeric)
If character(1), the algorithm to use in generating probability bins. See hist
for details. If numeric, the cut points for the bins. Default is “Sturges”.

groups (integer(1))
The number of bins to construct. If specified, breaks is ignored. Default is
NULL.

task.id (character(1))
Selected task in BenchmarkResult to do plots for, ignored otherwise. Default is
first task.

Value

CalibrationData. A list containing:

proportion data.frame with columns:
  • Learner Name of learner.
  • bin Bins calculated according to the breaks or groups argument.
  • Class Class labels (for binary classification only the positive class).
  • Proportion Proportion of observations from class Class among all obser-
vations with posterior probabilities of class Class within the interval given
in bin.

data data.frame with columns:
  • Learner Name of learner.
  • truth True class label.
  • Class Class labels (for binary classification only the positive class).
  • Probability Predicted posterior probability of Class.
  • bin Bin corresponding to Probability.

task (TaskDesc)
Task description.

References

generateCritDifferencesData

Generate data for critical-differences plot.

Description

Generates data that can be used to plot a critical differences plot. Computes the critical differences according to either the "Bonferroni-Dunn" test or the "Nemenyi" test. "Bonferroni-Dunn" usually yields higher power as it does not compare all algorithms to each other, but all algorithms to a baseline instead. Learners are drawn on the y-axis according to their average rank. For test = "nemenyi" a bar is drawn, connecting all groups of not significantly different learners. For test = "bd" an interval is drawn around the algorithm selected as a baseline. All learners within this interval are not significantly different from the baseline.

Calculation:

\[ CD = q_\alpha \sqrt{\frac{k(k+1)}{6N}} \]

Where \( q_\alpha \) is based on the studentized range statistic. See references for details.

Usage

```r
generateCritDifferencesData(
  bmr,
  measure = NULL,
  p.value = 0.05,
  baseline = NULL,
  test = "bd"
)
```

Arguments

- `bmr` *(BenchmarkResult)* Benchmark result.
- `measure` *(Measure)* Performance measure. Default is the first measure used in the benchmark experiment.
- `p.value` *(numeric(1))* P-value for the critical difference. Default: 0.05

See Also

Other generate_plot_data: `generateCritDifferencesData()`, `generateFeatureImportanceData()`, `generateFilterValuesData()`, `generateLearningCurveData()`, `generatePartialDependenceData()`, `generateThreshVsPerfData()`, `plotFilterValues()` Other calibration: `plotCalibration()`
baseline (character(1)): (learner.id)
Select a learner.id as baseline for the test = "bd" ("Bonferroni-Dunn") critical differences diagram. The critical difference interval will then be positioned around this learner. Defaults to best performing algorithm. For test = "nemenyi", no baseline is needed as it performs all pairwise comparisons.

test (character(1))
Test for which the critical differences are computed.
"bd" for the Bonferroni-Dunn Test, which is comparing all classifiers to a baseline, thus performing a comparison of one classifier to all others. Algorithms not connected by a single line are statistically different from the baseline.
"nemenyi" for the PMCMR::posthoc.friedman.nemenyi.test which is comparing all classifiers to each other. The null hypothesis that there is a difference between the classifiers can not be rejected for all classifiers that have a single grey bar connecting them.

Value
(critDifferencesData). List containing:

data (data.frame) containing the info for the descriptive part of the plot
friedman.nemenyi.test (list) of class pairwise.htest contains the calculated PMCMR::posthoc.friedman.nemenyi.test
cd.info (list) containing info on the critical difference and its positioning
baseline baseline chosen for plotting
p.value p.value used for the PMCMR::posthoc.friedman.nemenyi.test and for computation of the critical difference

See Also
Other generate_plot_data: generateCalibrationData(), generateFeatureImportanceData(), generateFilterValuesData(), generateLearningCurveData(), generatePartialDependenceData(), generateThreshVsPerfData()

Other benchmark: BenchmarkResult, benchmark(), benchmark().convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMRLearnedIds(), getBMRLearnedShortNames(), getBMRLearners(), getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMRXboxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()
generateFeatureImportanceData

Generate feature importance.

Description

Estimate how important individual features or groups of features are by contrasting prediction performances. For method “permutation.importance” compute the change in performance from permuting the values of a feature (or a group of features) and compare that to the predictions made on the unpermuted data.

Usage

```r
generateFeatureImportanceData(
  task,
  method = "permutation.importance",
  learner,
  features = getTaskFeatureNames(task),
  interaction = FALSE,
  measure,
  contrast = function(x, y) x - y,
  aggregation = mean,
  nmc = 50L,
  replace = TRUE,
  local = FALSE,
  show.info = FALSE
)
```

Arguments

- **task** (Task)
  The task.

- **method** (character(1))
  The method used to compute the feature importance. The only method available is “permutation.importance”. Default is “permutation.importance”.

- **learner** (Learner | character(1))
  The learner. If you pass a string the learner will be created via `makeLearner`.

- **features** (character)
  The features to compute the importance of. The default is all of the features contained in the `Task`.

- **interaction** (logical(1))
  Whether to compute the importance of the `features` argument jointly. For `method = "permutation.importance"` this entails permuting the values of all features together and then contrasting the performance with that of the performance without the features being permuted. The default is `FALSE`.
**generateFeatureImportanceData**

<table>
<thead>
<tr>
<th><strong>measure</strong></th>
<th>(Measure)</th>
<th>Performance measure. Default is the first measure used in the benchmark experiment.</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>contrast</strong></td>
<td>(function)</td>
<td>A difference function that takes a numeric vector and returns a numeric vector of the same length. The default is element-wise difference between the vectors.</td>
</tr>
<tr>
<td><strong>aggregation</strong></td>
<td>(function)</td>
<td>A function which aggregates the differences. This function must take a numeric vector and return a numeric vector of length 1. The default is mean.</td>
</tr>
<tr>
<td><strong>nmc</strong></td>
<td>(integer(1))</td>
<td>The number of Monte-Carlo iterations to use in computing the feature importance. If nmc == -1 and method = ”permutation.importance” then all permutations of the features are used. The default is 50.</td>
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<tr>
<td><strong>replace</strong></td>
<td>(logical(1))</td>
<td>Whether or not to sample the feature values with or without replacement. The default is TRUE.</td>
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<tr>
<td><strong>local</strong></td>
<td>(logical(1))</td>
<td>Whether to compute the per-observation importance. The default is FALSE.</td>
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<tr>
<td><strong>show.info</strong></td>
<td>(logical(1))</td>
<td>Whether progress output (feature name, time elapsed) should be displayed.</td>
</tr>
</tbody>
</table>

**Value**

(FeatureImportance). A named list which contains the computed feature importance and the input arguments.

Object members:

| **res** | (data.frame) | Has columns for each feature or combination of features (colon separated) for which the importance is computed. A row corresponds to importance of the feature specified in the column for the target. |
| **interaction** | (logical(1)) | Whether or not the importance of the features was computed jointly rather than individually. |
| **measure** | (Measure) | The measure used to compute performance. |

**contrast** (function) The function used to compare the performance of predictions.

**aggregation** (function) The function which is used to aggregate the contrast between the performance of predictions across Monte-Carlo iterations.

**replace** (logical(1)) Whether or not, when method = ”permutation.importance”, the feature values are sampled with replacement.
generateFilterValuesData

Calculates feature filter values.

Description

Calculates numerical filter values for features. For a list of features, use listFilterMethods.

Usage

generateFilterValuesData(
  task,  
  method = "randomForestSRC_importance",  
  nselect = getTaskNFeats(task),  
  ...,
  more.args = list()
)
Arguments

- task: (Task)
The task.
- method: (character | list)
  Filter method(s). In case of ensemble filters the list notation needs to be used. See the examples for more information. Default is "randomForestSRC_importance".
- nselect: (integer(1))
  Number of scores to request. Scores are getting calculated for all features per default.
- ...: (any)
  Passed down to selected method. Can only be use if method contains one element.
- more.args: (named list)
  Extra args passed down to filter methods. List elements are named with the filter method name the args should be passed down to. A more general and flexible option than .... Default is empty list.

Value

(FilterValues). A list containing:

- task.desc: [TaskDesc]
  Task description.
- data: (data.frame) with columns:
  - name: (character)
    Name of feature.
  - type: (character)
    Feature column type.
  - method: (numeric)
    One column for each method with the feature importance values.

Simple and ensemble filters

Besides passing (multiple) simple filter methods you can also pass an ensemble filter method (in a list). The ensemble method will use the simple methods to calculate its ranking. See listFilterEnsembleMethods() for available ensemble methods.

See Also

Other generate_plot_data: generateCalibrationData(), generateCritDifferencesData(), generateFeatureImportanceData(), generateLearningCurveData(), generatePartialDependenceData(), generateThreshVsPerfData(), plotFilterValues()

Other filter: filterFeatures(), getFilteredFeatures(), listFilterEnsembleMethods(), listFilterMethods(), makeFilterEnsemble(), makeFilterWrapper(), makeFilter(), plotFilterValues()
Examples

```r
# two simple filter methods
gval = generateFilterValuesData(iris.task, 
    method = c("FSelectorRcpp_gain.ratio", "FSelectorRcpp_information.gain"))
# using ensemble method "E-mean"
gval = generateFilterValuesData(iris.task, 
    method = list("E-mean", c("FSelectorRcpp_gain.ratio", 
        "FSelectorRcpp_information.gain")))
```

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**generateHyperParsEffectData**

*Generate hyperparameter effect data.*

**Description**

Generate cleaned hyperparameter effect data from a tuning result or from a nested cross-validation tuning result. The object returned can be used for custom visualization or passed downstream to an out of the box mlr method, `plotHyperParsEffect`.

**Usage**

```r
generateHyperParsEffectData(
    tune.result, 
    include.diagnostics = FALSE, 
    trafo = FALSE, 
    partial.dep = FALSE
)
```

**Arguments**

- `tune.result` *(TuneResult | ResampleResult)*
  
  Result of `tuneParams` (or `resample` ONLY when used for nested cross-validation). The tuning result (or results if the output is from nested cross-validation), also containing the optimizer results. If nested CV output is passed, each element in the list will be considered a separate run, and the data from each run will be included in the dataframe within the returned `HyperParsEffectData`.

- `include.diagnostics` *(logical(1))*
  
  Should diagnostic info (eol and error msg) be included? Default is `FALSE`.

- `trafo` *(logical(1))*
  
  Should the units of the hyperparameter path be converted to the transformed scale? This is only useful when `trafo` was used to create the path. Default is `FALSE`.

- `partial.dep` *(logical(1))*
  
  Should partial dependence be requested based on converting to reg task? This sets a flag so that we know to use partial dependence downstream. This should
generateLearningCurveData

Generates a learning curve.

Description

Observe how the performance changes with an increasing number of observations.
Usage

generateLearningCurveData(
    learners,
    task,
    resampling = NULL,
    percs = seq(0.1, 1, by = 0.1),
    measures,
    stratify = FALSE,
    show.info = getMlrOption("show.info")
)

Arguments

learners   [(list of) [Learner])
Learning algorithms which should be compared.

task       (Task)
The task.

resampling ([ResampleDesc] | [ResampleInstance])
Resampling strategy to evaluate the performance measure. If no strategy is given a
default "Holdout" will be performed.

percs      ([numeric])
Vector of percentages to be drawn from the training split. These values represent
the x-axis. Internally [makeDownsampleWrapper] is used in combination with
[benchmark]. Thus for each percentage a different set of observations is drawn
resulting in noisy performance measures as the quality of the sample can differ.

measures    [(list of) [Measure])
Performance measures to generate learning curves for, representing the y-axis.

stratify   (logical(1))
Only for classification: Should the downsampled data be stratified according to
the target classes?

show.info  (logical(1))
Print verbose output on console? Default is set via configureMlr.

Value

([LearningCurveData]). A ‘list’ containing:

task       ([Task])
The task.

measures    [(list of) [Measure])
Performance measures.

data        ([data.frame]) with columns:
  - ‘learner’ Names of learners.
  - ‘percentage’ Percentages drawn from the training split.
  - One column for each [Measure] passed to [generateLearningCurveData].
generatePartialDependenceData

See Also

Other generate_plot_data: generateCalibrationData(), generateCritDifferencesData(), generateFeatureImportanceData(), generateFilterValuesData(), generatePartialDependenceData(), generateThreshVsPerfData(), plotFilterValues()

Other learning_curve: plotLearningCurve()

Examples

r = generateLearningCurveData(list("classif.rpart", "classif.knn"),
  task = sonar.task, percs = seq(0.2, 1, by = 0.2),
  measures = list(tp, fp, tn, fn), resampling = makeResampleDesc(method = "Subsample", iters = 5),
  show.info = FALSE)
plotLearningCurve(r)

generatePartialDependenceData

Generate partial dependence.

Description

Estimate how the learned prediction function is affected by one or more features. For a learned
function f(x) where x is partitioned into x_s and x_c, the partial dependence of f on x_s can be
summarized by averaging over x_c and setting x_s to a range of values of interest, estimating
E_(x_c)(f(x_s, x_c)). The conditional expectation of f at observation i is estimated similarly. Addi-
tionally, partial derivatives of the marginalized function w.r.t. the features can be computed.

Usage

generatePartialDependenceData(
  obj,
  input,
  features = NULL,
  interaction = FALSE,
  derivative = FALSE,
  individual = FALSE,
  fun = mean,
  bounds = c(qnorm(0.025), qnorm(0.975)),
  uniform = TRUE,
  n = c(10, NA),
  ...
)

Arguments

obj (WrappedModel)
  Result of train.
generatePartialDependenceData

input (data.frame | Task)
Input data.

features character
A vector of feature names contained in the training data. If not specified all
features in the input will be used.

interaction (logical(1))
Whether the features should be interacted or not. If TRUE then the Cartesian
product of the prediction grid for each feature is taken, and the partial depen-
dence at each unique combination of values of the features is estimated. Note
that if the length of features is greater than two, plotPartialDependence cannot
be used. If FALSE each feature is considered separately. In this case features
can be much longer than two. Default is FALSE.

derivative (logical(1))
Whether or not the partial derivative of the learned function with respect to the
features should be estimated. If TRUE interaction must be FALSE. The partial
derivative of individual observations may be estimated. Note that computation
time increases as the learned prediction function is evaluated at gridsize points
* the number of points required to estimate the partial derivative. Additional
arguments may be passed to numDeriv::grad (for regression or survival tasks) or
numDeriv::jacobian (for classification tasks). Note that functions which are not
smooth may result in estimated derivatives of 0 (for points where the function
does not change within +/- epsilon) or estimates trending towards +/- infinity (at
discontinuities). Default is FALSE.

individual (logical(1))
Whether to plot the individual conditional expectation curves rather than the ag-
ggregated curve, i.e., rather than aggregating (using fun) the partial dependences
of features, plot the partial dependences of all observations in data across all
values of the features. The algorithm is developed in Goldstein, Kapelner,
Bleich, and Pitkin (2015). Default is FALSE.

fun function
A function which operates on the output on the predictions made on the input
data. For regression this means a numeric vector, and, e.g., for a multiclass
classification problem, this might instead be probabilities which are returned as a
numeric matrix. This argument can return vectors of arbitrary length, however,
if their length is greater than one, they must be named, e.g., fun = mean or fun =
function(x) c("mean" = mean(x),"variance" = var(x)). The default is the
mean, unless obj is classification with predict.type = "response" in which
case the default is the proportion of observations predicted to be in each class.

bounds (numeric(2))
The value (lower, upper) the estimated standard error is multiplied by to es-
timate the bound on a confidence region for a partial dependence. Ignored if
predict.type != "se" for the learner. Default is the 2.5 and 97.5 quantiles
(-1.96, 1.96) of the Gaussian distribution.

uniform (logical(1))
Whether or not the prediction grid for the features is a uniform grid of size
n[1] or sampled with replacement from the input. Default is TRUE.
\[ n \] (integer21)
The first element of \( n \) gives the size of the prediction grid created for each feature. The second element of \( n \) gives the size of the sample to be drawn without replacement from the input data. Setting \( n[2] \) less than the number of rows in the input will decrease computation time. The default for \( n[1] \) is 10, and the default for \( n[2] \) is the number of rows in the input.

\[
\text{...}\]
additional arguments to be passed to \texttt{mmpf::marginalPrediction}.

**Value**

\textbf{PartialDependenceData}. A named list, which contains the partial dependence, input data, target, features, task description, and other arguments controlling the type of partial dependences made.

Object members:

- **data** data.frame
  - Has columns for the prediction: one column for regression and survival analysis, and a column for class and the predicted probability for classification as well as a column for each element of features. If \texttt{individual = TRUE} then there is an additional column \texttt{idx} which gives the index of the data that each prediction corresponds to.

- **task.desc** TaskDesc
  - Task description.

- **target**
  - Target feature for regression, target feature levels for classification, survival and event indicator for survival.

- **features** character
  - Features argument input.

- **interaction** \( (\text{logical}(1)) \)
  - Whether or not the features were interacted (i.e. conditioning).

- **derivative** \( (\text{logical}(1)) \)
  - Whether or not the partial derivative was estimated.

- **individual** \( (\text{logical}(1)) \)
  - Whether the partial dependences were aggregated or the individual curves are retained.

**References**


**See Also**

Other \texttt{partial_dependence}: \texttt{plotPartialDependence()}

Other \texttt{generate_plot_data}: \texttt{generateCalibrationData()}, \texttt{generateCritDifferencesData()}, \texttt{generateFeatureImportanceData()}, \texttt{generateFilterValuesData()}, \texttt{generateLearningCurveData()}, \texttt{generateThreshVsPerfData()}, \texttt{plotFilterValues()}
generateThreshVsPerfData

Generate threshold vs. performance(s) for 2-class classification.

Usage

```r
generateThreshVsPerfData(
  obj, 
  measures, 
  gridsize = 100L, 
  aggregate = TRUE, 
  task.id = NULL 
)
```

Arguments

- **obj** (list of Prediction | list of ResampleResult | BenchmarkResult)
  - Single prediction object, list of them, single resample result, list of them, or a benchmark result. In case of a list probably produced by different learners you want to compare, then name the list with the names you want to see in the plots, probably learner shortnames or ids.

- **measures** (Measure | list of Measure)
  - Performance measure(s) to evaluate. Default is the default measure for the task, see here `getDefaultValue`.

- **gridsize** (integer(1))
  - Grid resolution for x-axis (threshold). Default is 100.

- **aggregate** (logical(1))
  - Whether to aggregate ResamplePredictions or to plot the performance of each iteration separately. Default is TRUE.

- **task.id** (character(1))
  - Selected task in BenchmarkResult to do plots for, ignored otherwise. Default is first task.
**getBMRAggrPerformances**

Extract the aggregated performance values from a benchmark result.

**Description**

Either a list of lists of “aggr” numeric vectors, as returned by resample, or these objects are rbind-ed with extra columns “task.id” and “learner.id”.

**Usage**

```r
getBMRAggrPerformances(
  bmr,
  task.ids = NULL,
  learner.ids = NULL,
  as.df = FALSE,
  drop = FALSE
)
```

**Arguments**

- `bmr` *(BenchmarkResult)*
  Benchmark result.
- `task.ids` *(character(1))*
  Restrict result to certain tasks. Default is all.
- `learner.ids` *(character(1))*
  Restrict result to certain learners. Default is all.
- `as.df` *(character(1))*
  Return one data.frame as result - or a list of lists of objects?. Default is FALSE.
- `drop` *(logical(1))*
  If drop is FALSE (the default), a nested list with the following structure is returned: `res[task.ids][learner.ids]`.

**Value**

`(ThresholdVsPerfData)`. A named list containing the measured performance across the threshold grid, the measures, and whether the performance estimates were aggregated (only applicable for (list of) ResampleResults).

**See Also**

Other generate_plot_data: generateCalibrationData(), generateCritDifferencesData(), generateFeatureImportanceData(), generateFilterValuesData(), generateLearningCurveData(), generatePartialDependenceData(), plotFilterValues()

Other thresh_vs_perf: plotROCCurves(), plotThreshVsPerf()
If drop is set to TRUE it is checked if the list structure can be simplified. If only one learner was passed, a list with entries for each task is returned. If only one task was passed, the entries are named after the corresponding learner. For an experiment with both one task and learner, the whole list structure is removed. Note that the name of the task/learner will be dropped from the return object.

Value

(list | data.frame). See above.

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMR LearnerIds(), getBMR Learner ShortNames(), getBMR Learners(), getBMR MeasureIds(), getBMR Measures(), getBMR Models(), getBMR Performances(), getBMR Predictions(), getBMR TaskDescs(), getBMR TaskIds(), getBMR Tune Results(), plotBM RBoxplots(), plotBM R RanksAsBarChart(), plotBMR Summary(), plotCritDifferences(), reduceBatchmarkResults()

getBMRFeatSelResults

Extract the feature selection results from a benchmark result.

Description

Returns a nested list of FeatSelResults. The first level of nesting is by data set, the second by learner, the third for the benchmark resampling iterations. If as.df is TRUE, a data frame with “task.id”, “learner.id”, the resample iteration and the selected features is returned. Note that if more than one feature is selected and a data frame is requested, there will be multiple rows for the same dataset-learner-iteration; one for each selected feature.

Usage

getBMRFeatSelResults(  
  bmr,  
  task.ids = NULL,  
  learner.ids = NULL,  
  as.df = FALSE,  
  drop = FALSE  
)

Arguments

bmr (BenchmarkResult) Benchmark result.
getBMRFilteredFeatures

Extract the feature selection results from a benchmark result.

Description

Returns a nested list of characters. The first level of nesting is by data set, the second by learner, the third for the benchmark resampling iterations. The list at the lowest level is the list of selected features. If `as.df` is `TRUE`, a data frame with “task.id”, “learner.id”, the resample iteration and the selected features is returned.

Note that if more than one feature is selected and a data frame is requested, there will be multiple rows for the same dataset-learner-iteration; one for each selected feature.
getBMRFilteredFeatures

Usage

getBMRFilteredFeatures(bmr, task.ids = NULL, learner.ids = NULL, as.df = FALSE, drop = FALSE)

Arguments

bmr (BenchmarkResult) Benchmark result.
task.ids (character(1)) Restrict result to certain tasks. Default is all.
learner.ids (character(1)) Restrict result to certain learners. Default is all.
as.df (character(1)) Return one data.frame as result - or a list of lists of objects? Default is FALSE.
drop (logical(1)) If drop is FALSE (the default), a nested list with the following structure is returned:
res[task.ids][learner.ids]. If drop is set to TRUE it is checked if the list structure can be simplified.
If only one learner was passed, a list with entries for each task is returned.
If only one task was passed, the entries are named after the corresponding learner.
For an experiment with both one task and learner, the whole list structure is removed.
Note that the name of the task/learner will be dropped from the return object.

Value

(list | data.frame). See above.

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRLearners(), getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMRRootBoxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()
getBMRLearnerIds

Return learner ids used in benchmark.

Description

Gets the IDs of the learners used in a benchmark experiment.

Usage

getBMRLearnerIds(bmr)

Arguments

bmr (BenchmarkResult)
Benchmark result.

Value

(character).

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMRLearnerShortNames(), getBMRLearners(), getBMRMetricIds(), getBMRMetrics(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMRBBoxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

getBMRLearners

Return learners used in benchmark.

Description

Gets the learners used in a benchmark experiment.

Usage

getBMRLearners(bmr)

Arguments

bmr (BenchmarkResult)
Benchmark result.
getBMRLearnerShortNames

Value

(list).

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(),
friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(),
getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMRLearnerIds(), getBMRLearnerShortNames(),
getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(),
getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMRBoxplots(), plotBMRanksAsBarChart(),
plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

getBMRLearnerShortNames

Return learner short.names used in benchmark.

Description

Gets the learner short.names of the learners used in a benchmark experiment.

Usage

getBMRLearnerShortNames(bmr)

Arguments

bmr

(BenchmarkResult)

Benchmark result.

Value

(character).

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(),
friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(),
getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMRLearnerIds(), getBMRLearnerShortNames(),
getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(),
getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMRBoxplots(), plotBMRanksAsBarChart(),
plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()
getBMRMeasureIds

Return measures IDs used in benchmark.

Description

Gets the IDs of the measures used in a benchmark experiment.

Usage

getBMRMeasureIds(bmr)

Arguments

bmr (BenchmarkResult)
Benchmark result.

Value

(list). See above.

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(),
friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(),
getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMR LearnerIds(), getBMR Learner Short Names(),
getBMR Learners(), getBMR Measures(), getBMR Models(), getBMR Performances(), getBMR Predictions(),
getBMR Task Descs(), getBMR Task Ids(), getBMR Tune Results(), plotBMR Boxplots(), plotBMR Ranks As Bar Chart(),
plotBMR Summary(), plot Crit Differences(), reduceBatchmarkResults()

getBMRMeasures

Return measures used in benchmark.

Description

Gets the measures used in a benchmark experiment.

Usage

getBMRMeasures(bmr)

Arguments

bmr (BenchmarkResult)
Benchmark result.
Value

(list). See above.

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMR LearnerIds(), getBMR Learner ShortNames(), getBMR Learners(), getBMRAggregateIds(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMR TuneResults(), plotBMRRboxplots(), plotBMRRanksAsBarChart(), plotBMRS Summary(), plotCritDifferences(), reduceBatchmarkResults()

getBMRModels Extract all models from benchmark result.

Description

A list of lists containing all WrappedModels trained in the benchmark experiment.

If models is FALSE in the call to benchmark, the function will return NULL.

Usage

getBMRModels(bmr, task.ids = NULL, learner.ids = NULL, drop = FALSE)

Arguments

  bmr  (BenchmarkResult)
  Benchmark result.

  task.ids  (character(1))
  Restrict result to certain tasks. Default is all.

  learner.ids  (character(1))
  Restrict result to certain learners. Default is all.

  drop  (logical(1))
  If drop is FALSE (the default), a nested list with the following structure is returned:
  res[task.ids][learner.ids].
  If drop is set to TRUE it is checked if the list structure can be simplified.
  If only one learner was passed, a list with entries for each task is returned.
  If only one task was passed, the entries are named after the corresponding learner.
  For an experiment with both one task and learner, the whole list structure is removed.
  Note that the name of the task/learner will be dropped from the return object.

Value

(list).
getBMRPerformances

Extract the test performance values from a benchmark result.

Description

Either a list of lists of “measure.test” data.frames, as returned by resample, or these objects are rbind-ed with extra columns “task.id” and “learner.id”.

Usage

getBMRPerformances(
  bmr,
  task.ids = NULL,
  learner.ids = NULL,
  as.df = FALSE,
  drop = FALSE
)

Arguments

bmr  (BenchmarkResult)
  Benchmark result.

task.ids  (character(1))
  Restrict result to certain tasks. Default is all.

learner.ids  (character(1))
  Restrict result to certain learners. Default is all.

as.df  (character(1))
  Return one data.frame as result - or a list of lists of objects?. Default is FALSE.

drop  (logical(1))
  If drop is FALSE (the default), a nested list with the following structure is returned:
  res[task.ids][learner.ids].
  If drop is set to TRUE it is checked if the list structure can be simplified.
  If only one learner was passed, a list with entries for each task is returned.
  If only one task was passed, the entries are named after the corresponding learner.
  For an experiment with both one task and learner, the whole list structure is removed.
  Note that the name of the task/learner will be dropped from the return object.
getBMRPredictions

Extract the predictions from a benchmark result.

Description

Either a list of lists of ResamplePrediction objects, as returned by resample, or these objects are rbind-ed with extra columns “task.id” and “learner.id”.

If predict.type is “prob”, the probabilities for each class are returned in addition to the response.

If keep.pred is FALSE in the call to benchmark, the function will return NULL.

Usage

getBMRPredictions(
  bmr,
  task.ids = NULL,
  learner.ids = NULL,
  as.df = FALSE,
  drop = FALSE
)

Arguments

bmr (BenchmarkResult)
Benchmark result.

task.ids (character(1))
Restrict result to certain tasks. Default is all.

learner.ids (character(1))
Restrict result to certain learners. Default is all.

as.df (character(1))
Return one data.frame as result - or a list of lists of objects? . Default is FALSE.
getBMRTaskDescriptions

Extract all task descriptions from benchmark result (DEPRECATED).

Description

A list containing all TaskDescs for each task contained in the benchmark experiment.

Usage

getBMRTaskDescriptions(bmr)

Arguments

bmr  (BenchmarkResult)

Benchmark result.

Value

(list)
getBMRTaskDescs

Extract all task descriptions from benchmark result.

Description

A list containing all TaskDescs for each task contained in the benchmark experiment.

Usage

getBMRTaskDescs(bmr)

Arguments

bmr (BenchmarkResult)

Benchmark result.

Value

(list).

See Also

Other benchmark: BenchmarkResult, benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFixedFeatures(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRLearners(), getBMRFixedFeatures(), getBMRFixedMeasureIds(), getBMRFixedMeasures(), getBMRFixedModels(), getBMRFixedPerformances(), getBMRFixedPredictions(), getBMRTaskIds(), getBMRTuneResults(), plotBMRSBoxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

getBMRTaskIds

Return task ids used in benchmark.

Description

Gets the task IDs used in a benchmark experiment.

Usage

getBMRTaskIds(bmr)

Arguments

bmr (BenchmarkResult)

Benchmark result.
**getBMRTuneResults**

Extract the tuning results from a benchmark result.

**Value**

(character).

**See Also**

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMR LearnerIds(), getBMR LearnerShortNames(), getBMR Learners(), getBMRMeasureIds(), getBMR Measures(), getBMR Models(), getBMR Performances(), getBMRPredictions(), getBMR TaskDescs(), getBMRTuneResults(), plotBMRBoxplots(), plotBMR RanksAsBarChart(), plotBMR Summary(), plotCritDifferences(), reduceBatchmarkResults()

**Description**

Returns a nested list of TuneResults. The first level of nesting is by data set, the second by learner, the third for the benchmark resampling iterations. If as.df is TRUE, a data frame with the “task.id”, “learner.id”, the resample iteration, the parameter values and the performances is returned.

**Usage**

getBMRTuneResults(
  bmr,  
  task.ids = NULL,
  learner.ids = NULL,
  as.df = FALSE,
  drop = FALSE
)

**Arguments**

- **bmr** (BenchmarkResult)
  Benchmark result.
- **task.ids** (character(1))
  Restrict result to certain tasks. Default is all.
- **learner.ids** (character(1))
  Restrict result to certain learners. Default is all.
- **as.df** (character(1))
  Return one data.frame as result - or a list of lists of objects?. Default is FALSE.
- **drop** (logical(1))
  If drop is FALSE (the default), a nested list with the following structure is returned: res[task.ids][learner.ids]. If drop is set to TRUE it is checked if the list structure can be simplified.
getCaretParamSet

If only one learner was passed, a list with entries for each task is returned.
If only one task was passed, the entries are named after the corresponding
learner.
For an experiment with both one task and learner, the whole list structure is re-
moved.
Note that the name of the task/learner will be dropped from the return object.

Value

(list | data.frame). See above.

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(),
friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(),
getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMRLearnerIds(), getBMRLearnerShortNames(),
getBMRLearners(), getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(),
getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), plotBMROboxplots(), plotBMRRanksAsBarChart(),
plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

getCaretParamSet  Get tuning parameters from a learner of the caret R-package.

Description

Constructs a grid of tuning parameters from a learner of the caret R-package. These values are then
converted into a list of non-tunable parameters (par.vals) and a tunable ParamHelpers::ParamSet
(par.set), which can be used by tuneParams for tuning the learner. Numerical parameters will
either be specified by their lower and upper bounds or they will be discretized into specific values.

Usage

getcaretParamSet(learner, length = 3L, task, discretize = TRUE)

Arguments

  learner  (character(1))
The name of the learner from caret (cf. https://topepo.github.io/caret/
available-models.html). Note that the names in caret often differ from the
ones in mlr.

  length  (integer(1))
A length / precision parameter which is used by caret for generating the grid of
tuning parameters. caret generates either as many values per tuning parameter
/ dimension as defined by length or only a single value (in case of non-tunable
par.vals).

  task  (Task)
Learning task, which might be requested for creating the tuning grid.
Should the numerical parameters be discretized? Alternatively, they will be defined by their lower and upper bounds. The default is TRUE.

Value

(list(2)). A list of parameters:

- `par.vals` contains a list of all constant tuning parameters
- `par.set` is a `ParamHelpers::ParamSet`, containing all the configurable tuning parameters

Examples

```r
if (requireNamespace("caret") && requireNamespace("mlbench")) {
  library(caret)
  classifTask = makeClassifTask(data = iris, target = "Species")

  # (1) classification (random forest) with discretized parameters
  getCaretParamSet("rf", length = 9L, task = classifTask, discretize = TRUE)

  # (2) regression (gradient boosting machine) without discretized parameters
  library(mlbench)
  data(BostonHousing)
  regrTask = makeRegrTask(data = BostonHousing, target = "medv")
  getCaretParamSet("gbm", length = 9L, task = regrTask, discretize = FALSE)
}
```

`getClassWeightParam` Get the class weight parameter of a learner.

Description

Gets the class weight parameter of a learner.

Usage

`getClassWeightParam(learner, lrn.id = NULL)`

Arguments

- `learner` (Learner | character(1))
  The learner. If you pass a string the learner will be created via `makeLearner`.
- `lrn.id` (character)
  Only used for BaseEnsembles. It is possible that multiple learners in a base ensemble have a class weight param. Specify the learner from which the class weight should be extracted.
**Value**

numeric LearnerParam: A numeric parameter object, containing the class weight parameter of the given learner.

**See Also**

Other learner: LearnerProperties, getHyperPars(), getLearnerId(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(), setPredictThreshold(), setPredictType()
getFailureModelDump

Return the error dump of FailureModel.

Description

Returns the error dump that can be used with debugger() to evaluate errors. If configureMlr configuration on.error.dump is FALSE, this returns NULL.

Usage

getFailureModelDump(model)

Arguments

model (WrappedModel)
The model.

defaultMeasure

Get default measure.

Description

Get the default measure for a task type, task, task description or a learner. Currently these are:

classif mmce
regr mse
cluster db
surv cindex
costsens mcp
multilabel multilabel.hamloss

Usage

defaultMeasure(x)

Arguments

x (character(1) | Task | TaskDesc | Learner)
Task type, task, task description, learner name, a learner, or a type of learner (e.g. "classif").

Value

(Measure).
getFeatSelResult

Description

Returns the selected feature set and optimization path after training.

Usage

getFeatSelResult(object)

Arguments

object (WrappedModel)
Trained Model created with makeFeatSelWrapper.

Value

(FeatSelResult).

getFailureModelMsg

Return error message of FailureModel.

Description

Such a model is created when one sets the corresponding option in configureMlr. If no failure occurred, NA is returned.

For complex wrappers this getter returns the first error message encountered in ANY model that failed.

Usage

getFailureModelMsg(model)

Arguments

model (WrappedModel)
The model.

Value

(character(1)).

getFeatSelResult

Returns the selected feature set and optimization path after training.

Description

Returns the selected feature set and optimization path after training.

Usage

getFeatSelResult(object)

Arguments

object (WrappedModel)
Trained Model created with makeFeatSelWrapper.

Value

(FeatSelResult).
getFeatureImportance

See Also

Other featsel: FeatSelControl, analyzeFeatSelResult(), makeFeatSelWrapper(), selectFeatures()

getFeatureImportance Calculates feature importance values for trained models.

Description

For some learners it is possible to calculate a feature importance measure. 'getFeatureImportance' extracts those values from trained models. See below for a list of supported learners.

- boosting
  Measure which accounts the gain of Gini index given by a feature in a tree and the weight of that tree.
- cforest
  Permutation principle of the 'mean decrease in accuracy' principle in randomForest. If 'auc=TRUE' (only for binary classification), area under the curve is used as measure. The algorithm used for the survival learner is 'extremely slow and experimental; use at your own risk'. See varimp for details and further parameters.
- gbm
  Estimation of relative influence for each feature. See relative.influence for details and further parameters.
- h2o
  Relative feature importances as returned by varimp.
- randomForest
  For 'type = 2' (the default) the 'MeanDecreaseGini' is measured, which is based on the Gini impurity index used for the calculation of the nodes. Alternatively, you can set 'type' to 1, then the measure is the mean decrease in accuracy calculated on OOB data. Note, that in this case the learner’s parameter ‘importance’ needs to be set to be able to compute feature importance values. See importance for details.
- RRF
  This is identical to randomForest.
- randomForestSRC
  This method can calculate feature importance for various measures. By default the Breiman-Cutler permutation method is used. See vimp for details.
- ranger
  Supports both measures mentioned above for the randomForest learner. Note, that you need to specifically set the learners parameter ‘importance‘, to be able to compute feature importance measures. See importance and ranger for details.
- rpart
  Sum of decrease in impurity for each of the surrogate variables at each node.
- xgboost
  The value implies the relative contribution of the corresponding feature to the model calculated by taking each feature’s contribution for each tree in the model. The exact computation of the importance in xgboost is undocumented.
getFeatureImportance

Usage

getFeatureImportance(object, ...)

Arguments

object ((WrappedModel))
Wrapped model, result of \[train\].

... (any)
Additional parameters, which are passed to the underlying importance value generating function.

Value

((FeatureImportance)) An object containing a ‘data.frame‘ of the variable importances and further information.

getFilteredFeatures

Returns the filtered features.

Description

Returns the filtered features.

Usage

getFilteredFeatures(model)

Arguments

model ((WrappedModel))
Trained Model created with \makeFilterWrapper\.

Value

(character).

See Also

Other filter: \filterFeatures()\, \generateFilterValuesData()\, \listFilterEnsembleMethods()\, \listFilterMethods()\, \makeFilterEnsemble()\, \makeFilterWrapper()\, \makeFilter()\, \plotFilterValues()
**getHomogeneousEnsembleModels**

*Deprecated, use getLearnerModel instead.*

**Description**

Deprecated, use getLearnerModel instead.

**Usage**

```r
getHomogeneousEnsembleModels(model, learner.models = FALSE)
```

**Arguments**

- `model` (Deprecated).
- `learner.models` (Deprecated).

---

**getHyperPars**

*Get current parameter settings for a learner.*

**Description**

Retrieves the current hyperparameter settings of a learner.

**Usage**

```r
getHyperPars(learner, for.fun = c("train", "predict", "both"))
```

**Arguments**

- `learner` *(Learner)*
  The learner.
- `for.fun` *(character(1))*
  Restrict the returned settings to hyperparameters corresponding to when they are used (see `ParamHelpers::LearnerParam`). Must be a subset of: “train”, “predict” or “both”. Default is `c("train", "predict", "both")`.

**Details**

This function only shows hyperparameters that differ from the learner default (because mlr changed the default) or if the user set hyperparameters manually during learner creation. If you want to have an overview of all available hyperparameters use `getParamSet()`.

**Value**

*(list). A named list of values.*
getLearnerId

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(), setPredictThreshold(), setPredictType()

Examples

getHyperPars(makeLearner("classif.ranger"))

## set learner hyperparameter 'mtry' manually
getHyperPars(makeLearner("classif.ranger", mtry = 100))

generateTable

getLearnerId

Get the ID of the learner.

Description

Get the ID of the learner.

Usage

generateTable

getLearnerId(learner)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

Value

(character(1)).

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(), setPredictThreshold(), setPredictType())
**getLearnerModel**  
Get underlying R model of learner integrated into mlr.

**Description**  
Get underlying R model of learner integrated into mlr.

**Usage**  
getLearnerModel(model, more.unwrap = FALSE)

**Arguments**
- **model** *(WrappedModel)*  
The model, returned by e.g., train.
- **more.unwrap** *(logical(1))*  
Some learners are not basic learners from R, but implemented in mlr as meta-techniques. Examples are everything that inherits from `HomogeneousEnsemble`. In these cases, the `learner.model` is often a list of mlr `WrappedModels`. This option allows to strip them further to basic R models. The option is simply ignored for basic learner models. Default is FALSE.

**Value**  
(any). A fitted model, depending the learner / wrapped package. E.g., a model of class `rpart::rpart` for learner “classif.rpart”.

**getLearnerNote**  
Get the note for the learner.

**Description**  
Get the note for the learner.

**Usage**  
getLearnerNote(learner)

**Arguments**
- **learner** *(Learner | character(1))*  
The learner. If you pass a string the learner will be created via `makeLearner`.

**Value**  
(character).
getLearnerParamSet

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(),
getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()

getLearnerPackages  
Get the required R packages of the learner.

Description
Get the R packages the learner requires.

Usage
getLearnerPackages(learner)

Arguments

learner  
(Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

Value
(character).

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(),
getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()

getLearnerParamSet  
Get the parameter set of the learner.

Description
Alias for getParamSet.

Usage
getLearnerParamSet(learner)
getLearnerParVals

Arguments

learner

(Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

Value

ParamSet.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerPredictType(),
getLearnerShortName(), getLearnerType(),getParamSet(), helpLearnerParam(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()

getLearnerParVals Get the parameter values of the learner.

Description

Alias for getHyperPars.

Usage

getLearnerParVals(learner, for.fun = c("train", "predict", "both"))

Arguments

learner

(Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

for.fun

(character(1))
Restrict the returned settings to hyperparameters corresponding to when the are
used (see ParamHelpers::LearnerParam). Must be a subset of: "train", "predict"
or "both". Default is c("train", "predict", "both").

Value

(list). A named list of values.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParamSet(), getLearnerPredictType(),
getLearnerShortName(), getLearnerType(),getParamSet(), helpLearnerParam(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()
getLearnerPredictType  
*Get the predict type of the learner.*

**Description**

Get the predict type of the learner.

**Usage**

```r
getLearnerPredictType(learner)
```

**Arguments**

- `learner` *(Learner | character(1))*
  
  The learner. If you pass a string the learner will be created via `makeLearner`.

**Value**

(character(1)).

**See Also**

Other learner: `LearnerProperties`, `getClassWeightParam()`, `getHyperPars()`, `getLearnerId()`, `getLearnerNote()`, `getLearnerPackages()`, `getLearnerParVals()`, `getLearnerParamSet()`, `getLearnerShortName()`, `getLearnerType()`, `getParamSet()`, `helpLearnerParam()`, `helpLearner()`, `makeLearners()`, `makeLearner()`, `removeHyperPars()`, `setHyperPars()`, `setId()`, `setLearnerId()`, `setPredictThreshold()`, `setPredictType()`

getLearnerShortName  
*Get the short name of the learner.*

**Description**

For an ordinary learner simply its short name is returned. For wrapped learners, the wrapper id is successively attached to the short name of the base learner. E.g: “rf.bagged.imputed”

**Usage**

```r
getLearnerShortName(learner)
```

**Arguments**

- `learner` *(Learner | character(1))*
  
  The learner. If you pass a string the learner will be created via `makeLearner`.
getLearnerType

Value

(character(1)).

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(),
getLearnerPredictType(), getLearnerType(), getParamSet().helpLearnerParam(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()

getLearnerType get the type of the learner.

Description

Get the type of the learner.

Usage

getLearnerType(learner)

Arguments

learner (Learner | character(1))

The learner. If you pass a string the learner will be created via makeLearner.

Value

(character(1)).

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(),
getLearnerPredictType(), getLearnerType(), getParamSet().helpLearnerParam(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(),
setLearnerId().setPredictThreshold(), setPredictType()
getMlrOptions

_returns a list of mlr's options._

Description

Gets the options for mlr.

Usage

getMlrOptions()

Value

(list).

See Also

Other configure: configureMlr()

getMultilabelBinaryPerformances

_retrieve binary classification measures for multilabel classification predictions._

Description

Measures the quality of each binary label prediction w.r.t. some binary classification performance measure.

Usage

getCodeMultilabelBinaryPerformances(pred, measures)

Arguments

pred (Prediction) Multilabel Prediction object.

measures (Measure | list of Measure) Performance measure(s) to evaluate, must be applicable to binary classification performance. Default is mmce.

Value

(named matrix). Performance value(s), column names are measure(s), row names are labels.
getNestedTuneResultsOptPathDf

See Also

Other multilabel: makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper()

Examples

# see makeMultilabelBinaryRelevanceWrapper

generateTuneResultsOptPathDf

Get the opt paths from each tuning step from the outer resampling.

Description

After you resampled a tuning wrapper (see makeTuneWrapper) with resample(..., extract = getTuneResult) this helper returns a data.frame with all opt.paths combined by rbind. An additional column iter indicates to what resampling iteration the row belongs.

Usage

generateTuneResultsOptPathDf(r, trafo = FALSE)

Arguments

r (ResampleResult) The result of resampling of a tuning wrapper.

trafo (logical(1)) Should the units of the hyperparameter path be converted to the transformed scale? This is only necessary when trafo was used to create the opt.paths. Note that opt.paths are always stored on the untransformed scale. Default is FALSE.

Value

(data.frame). See above.

See Also

Other tune: TuneControl, generateTuneResultsX(), getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(), makeTuneControlCMAES(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(), makeTuneControlIrace(), makeTuneControlMBO(), makeTuneControlRandom(), makeTuneWrapper(), tuneParams(), tuneThreshold()

Examples

# see example of makeTuneWrapper
getNestedTuneResultsX  
*Get the tuned hyperparameter settings from a nested tuning.*

**Description**

After you resampled a tuning wrapper (see `makeTuneWrapper`) with `resample(..., extract = getTuneResult) this helper returns a `data.frame` with the best found hyperparameter settings for each resampling iteration.

**Usage**

```r
getNestedTuneResultsX(r)
```

**Arguments**

- `r` **(ResampleResult)**
  The result of resampling of a tuning wrapper.

**Value**

- `data.frame`. One column for each tuned hyperparameter and one row for each outer resampling iteration.

**See Also**

Other `tune`: `TuneControl`, `getNestedTuneResultsOptPathDf()`, `getResamplingIndices()`, `getTuneResult()`, `makeModelMultiplexerParamSet()`, `makeModelMultiplexer()`, `makeTuneControlCMAES()`, `makeTuneControlDesign()`, `makeTuneControlGenSA()`, `makeTuneControlGrid()`, `makeTuneControlIrace()`, `makeTuneControlMBO()`, `makeTuneControlRandom()`, `makeTuneWrapper()`, `tuneParams()`, `tuneThreshold()`

**Examples**

```r
# see example of makeTuneWrapper
```

---

getOOBPreds  
*Extracts out-of-bag predictions from trained models.*

**Description**

Learners like `randomForest` produce out-of-bag predictions. `getOOBPreds` extracts this information from trained models and builds a prediction object as provided by `predict` (with prediction time set to NA). In the classification case: What is stored exactly in the `Prediction` object depends on the `predict.type` setting of the `Learner`.

You can call `listLearners(properties = "oobpreds")` to get a list of learners which provide this.
getParamSet

Usage

getOOBPreds(model, task)

Arguments

model (WrappedModel)
The model.

task (Task)
The task.

Value

(Prediction).

Examples

training.set = sample(1:150, 50)
lrn = makeLearner("classif.ranger", predict.type = "prob", predict.threshold = 0.6)
mod = train(lrn, sonar.task, subset = training.set)
oob = getOOBPreds(mod, sonar.task)
oob
performance(oob, measures = list(auc, mmce))

getParamSet

Get a description of all possible parameter settings for a learner.

Description

Returns the ParamHelpers::ParamSet from a Learner.

Value

ParamSet.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(),
getLearnerPredictType(), getLearnerShortName(), getLearnerType(), helpLearnerParam(),
helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(),
setLearnerId(), setPredictThreshold(), setPredictType()
**getPredictionDump**  
*Return the error dump of a failed Prediction.*

**Description**

Returns the error dump that can be used with debugger() to evaluate errors. If `configureMlr` configuration `on.error.dump` is `FALSE` or if the prediction did not fail, this returns NULL.

**Usage**

```r
getPredictionDump(pred)
```

**Arguments**

- `pred` *(Prediction)*
  - Prediction object.

**Value**

(last.dump).

**See Also**

Other debug: `FailureModel`, `ResampleResult`, `getRRDump()`

---

**getPredictionProbabilities**  
*Get probabilities for some classes.*

**Description**

Get probabilities for some classes.

**Usage**

```r
getPredictionProbabilities(pred, cl)
```

**Arguments**

- `pred` *(Prediction)*
  - Prediction object.
- `cl` *(character)*
  - Names of classes. Default is either all classes for multi-class / multilabel problems or the positive class for binary classification.
**getPredictionResponse**

**Value**

A `data.frame` with numerical columns or a numerical vector if length of `cl` is 1. Order of columns is defined by `cl`.

**See Also**

Other predict: `asROCRPrediction()`, `getPredictionResponse()`, `getPredictionTaskDesc()`, `predict.WrappedModel()`, `setPredictThreshold()`, `setPredictType()`

**Examples**

```r
task = makeClassifTask(data = iris, target = "Species")
lrn = makeLearner("classif.lda", predict.type = "prob")
mod = train(lrn, task)
# predict probabilities
pred = predict(mod, newdata = iris)

# Get probabilities for all classes
head(getPredictionProbabilities(pred))

# Get probabilities for a subset of classes
head(getPredictionProbabilities(pred, c("setosa", "virginica")))
```

---

**getDescription**  Get response / truth from prediction object.

**Description**

The following types are returned, depending on task type:

- **classif**: factor
- **regr**: numeric
- **se**: numeric
- **cluster**: integer
- **surv**: numeric
- **multilabel**: logical matrix, columns named with labels

**Usage**

```
getPredictionResponse(pred)
getPredictionSE(pred)
getPredictionTruth(pred)
```
getPredictionTaskDesc

Arguments

pred (Prediction)
Prediction object.

Value

See above.

See Also

Other predict: asROCRPrediction().getPredictionProbabilities().getPredictionTaskDesc().predict.WrappedModel().setPredictThreshold().setPredictType()
getProbabilities

Deprecated, use getPredictionProbabilities instead.

Description

Deprecated, use getPredictionProbabilities instead.

Usage

getProbabilities(pred, cl)

Arguments

pred	Deprecated.
cl	Deprecated.

getResamplingIndices

Get the resampling indices from a tuning or feature selection wrapper.

Description

After you resampled a tuning or feature selection wrapper (see makeTuneWrapper) with resample(..., extract = getTuneResult) or resample(..., extract = getFeatSelResult) this helper returns a list with the resampling indices used for the respective method.

Usage

getResamplingIndices(object, inner = FALSE)

Arguments

object	(ResampleResult)
The result of resampling of a tuning or feature selection wrapper.
inner	(logical)
If TRUE, returns the inner indices of a nested resampling setting.

Value

(list). One list for each outer resampling fold.

See Also

Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(),
getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(), makeTuneControlCMAES(),
makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(), makeTuneControlIrace(),
makeTuneControlMBO(), makeTuneControlRandom(), makeTuneWrapper(), tuneParams(), tuneThreshold()
Examples

```r
task = makeClassifTask(data = iris, target = "Species")
lnr = makeLearner("classif.rpart")
# stupid mini grid
ps = makeParamSet(
  makeDiscreteParam("cp", values = c(0.05, 0.1)),
  makeDiscreteParam("minsplit", values = c(10, 20))
)
ctrl = makeTuneControlGrid()
inner = makeResampleDesc("Holdout")
outer = makeResampleDesc("CV", iters = 2)
lnr = makeTuneWrapper(lnr, resampling = inner, par.set = ps, control = ctrl)
# nested resampling for evaluation
# we also extract tuned hyper pars in each iteration and by that the resampling indices
r = resample(lnr, task, outer, extract = getTuneResult)
# get tuning indices
getResamplingIndices(r, inner = TRUE)
```

getRRDump

Return the error dump of ResampleResult.

Description

Returns the error dumps generated during resampling, which can be used with debugger() to debug errors. These dumps are saved if configureMlr configuration on.error.dump, or the corresponding learner config, is TRUE.

The returned object is a list with as many entries as the resampling being used has folds. Each of these entries can have a subset of the following slots, depending on which step in the resampling iteration failed: “train” (error during training step), “predict.train” (prediction on training subset), “predict.test” (prediction on test subset).

Usage

```r
getRRDump(res)
```

Arguments

- `res` (ResampleResult)
  The result of resample.

Value

- `list`

See Also

Other debug: FailureModel, ResampleResult, getPredictionDump()
getRRPredictionList

Get list of predictions for train and test set of each single resample iteration.

Description

This function creates a list with two slots train and test where each slot is again a list of Prediction objects for each single resample iteration. In case that predict = "train" was used for the resample description (see makeResampleDesc), the slot test will be NULL and in case that predict = "test" was used, the slot train will be NULL.

Usage

classicalRss(res, ...) 

Arguments

res (ResampleResult) The result of resample run with keep.pred = TRUE.

Value

list.

See Also

Other resample: ResamplePrediction, ResampleResult, addRRMeasure(), getRRPredictions(), getRRTaskDescription(), getRRTaskDesc(), makeResampleDesc(), makeResampleInstance(), resample()

getRRPredictions

Get predictions from resample results.

Description

Very simple getter.

Usage

classicalRss(res)

Arguments

res (ResampleResult) The result of resample run with keep.pred = TRUE.
getRRTaskDesc

Value

(ResamplePrediction).

See Also

Other resample: ResamplePrediction, ResampleResult, addRRMeasure(), getRRPredictionList(), getRRTaskDescription(), getRRTaskDesc(), makeResampleDesc(), makeResampleInstance(), resample()

getRRTaskDesc Get task description from resample results (DEPRECATED).

Description

Get a summarizing task description.

Usage

getRRTaskDesc(res)

Arguments

res (ResampleResult) The result of resample.

Value

(TaskDesc).

See Also

Other resample: ResamplePrediction, ResampleResult, addRRMeasure(), getRRPredictionList(), getRRPredictions(), getRRTaskDescription(), makeResampleDesc(), makeResampleInstance(), resample()
getRRTaskDescription  
Get task description from resample results (DEPRECATED).

Description
Get a summarizing task description.

Usage
getRRTaskDescription(res)

Arguments
res  (ResampleResult)
The result of resample.

Value
(TaskDesc).

See Also
Other resample: ResamplePrediction, ResampleResult, addRRMeasure(), getRRPredictionList(), getRRPredictions(), getRRTaskDesc(), makeResampleDesc(), makeResampleInstance(), resample()

getStackedBaseLearnerPredictions
Returns the predictions for each base learner.

Description
Returns the predictions for each base learner.

Usage
getStackedBaseLearnerPredictions(model, newdata = NULL)

Arguments
model  (WrappedModel)
Wrapped model, result of train.
newdata  (data.frame)
New observations, for which the predictions using the specified base learners should be returned. Default is NULL and extracts the base learner predictions that were made during the training.
getTaskClassLevels

Get the class levels for classification and multilabel tasks.

Description

NB: For multilabel, getTaskTargetNames and getTaskClassLevels actually return the same thing.

Usage

getTaskClassLevels(x)

Arguments

x

(Task | TaskDesc)
Task or its description object.

Value

(character).

See Also

Other task: getTaskCosts(), getTaskData(), getTaskDesc(), getTaskFeatureNames(), getTaskFormula(), getTaskId(), getTaskNFeats(), getTaskSize(), getTaskTargetNames(), getTaskTargets(), getTaskType(), subsetTask()

getTaskCosts

Extract costs in task.

Description

Returns “NULL” if the task is not of type “costsens”.

Usage

getTaskCosts(task, subset = NULL)

Arguments

task

(CostSensTask)
The task.

subset

(integer | logical | NULL)
Selected cases. Either a logical or an index vector. By default NULL if all observations are used.
getTaskData

Extract data in task.

Description

Useful in trainLearner when you add a learning machine to the package.

Usage

getTaskData(
  task,
  subset = NULL,
  features,
  target.extra = FALSE,
  recode.target = "no",
  functionals.as = "dfcols"
)

Arguments

task  (Task)
The task.

subset  (integer | logical | NULL)
Selected cases. Either a logical or an index vector. By default NULL if all observations are used.

features  (character | integer | logical)
Vector of selected inputs. You can either pass a character vector with the feature names, a vector of indices, or a logical vector.
In case of an index vector each element denotes the position of the feature name returned by getTaskFeatureNames.
Note that the target feature is always included in the resulting task, you should not pass it here. Default is to use all features.

target.extra  (logical(1))
Should target vector be returned separately? If not, a single data.frame including the target columns is returned, otherwise a list with the input data.frame and an extra vector or data.frame for the targets. Default is FALSE.
recode.target (character(1))
Should target classes be recoded? Supported are binary and multilabel classification and survival. Possible values for binary classification are “01”, “-1+1” and “drop.levels”. In the two latter cases the target vector is converted into a numeric vector. The positive class is coded as “+1” and the negative class either as “0” or “-1”. “drop.levels” will remove empty factor levels in the target column. In the multilabel case the logical targets can be converted to factors with “multilabel.factor”. For survival, you may choose to recode the survival times to “left”, “right” or “interval2” censored times using “lcens”, “rcens” or “icens”, respectively. See survival::Surv for the format specification. Default for both binary classification and survival is “no” (do nothing).

functionals.as (character(1))
How to represents functional features? Option “matrix”: Keep them as matrix columns in the data.frame. Option “dfcols”: Convert them to individual numeric data.frame columns. Default is “dfcols”.

Value
Either a data.frame or a list with data.frame data and vector target.

See Also
Other task: getTaskClassLevels(), getTaskCosts(), getTaskDesc(), getTaskFeatureNames(), getTaskFormula(), getTaskId(), getTaskNFeats(), getTaskSize(), getTaskTargetNames(), getTaskTargets(), getTaskType(), subsetTask()
getTaskDescription

Arguments

x (Task | TaskDesc)
Task or its description object.

Value

ret_taskdesc

See Also

Other task: getTaskClassLevels(), getTaskCosts(), getTaskData(), getTaskFeatureNames(),
getTaskFormula(), getTaskId(), getTaskNFeats(), getTaskSize(), getTaskTargetNames(),
getTaskTargets(), getTaskType(), subsetTask()

getTaskDescription  Deprecated, use getTaskDesc instead.

Description

Deprecated, use getTaskDesc instead.

Usage

getTaskDescription(x)

Arguments

x (Task | TaskDesc)
Task or its description object.

getTaskFeatureNames  Get feature names of task.

Description

Target column name is not included.

Usage

getTaskFeatureNames(task)

Arguments

task (Task)
The task.
getTaskFormula

Value

(character).

See Also

Other task: getTaskClassLevels(), getTaskCosts(), getTaskData(), getTaskDesc(), getTaskFormula(), getTaskId(), getTaskNFeats(), getTaskSize(), getTaskTargetNames(), getTaskTargets(), getTaskType(), subsetTask()

getTaskFormula

Get formula of a task.

Description

This is usually simply <target> ~ . For multilabel it is <target_1> + ... + <target_k> ~.

Usage

getTaskFormula(
  x,
  target = getTaskTargetNames(x),
  explicit.features = FALSE,
  env = parent.frame()
)

Arguments

x (Task | TaskDesc)
Task or its description object.

target (character(1))
Left hand side of the formula. Default is defined by task x.

explicit.features (logical(1))
Should the features (right hand side of the formula) be explicitly listed? Default is FALSE, i.e., they will be represented as ".".

env (environment)
Environment of the formula. Default is parent.frame().

Value

(formula).

See Also

Other task: getTaskClassLevels(), getTaskCosts(), getTaskData(), getTaskDesc(), getTaskFeatureNames(), getTaskId(), getTaskNFeats(), getTaskSize(), getTaskTargetNames(), getTaskTargets(), getTaskType(), subsetTask()
getTaskId

Get the id of the task.

Description
See title.

Usage
getTaskId(x)

Arguments
x (Task | TaskDesc)
Task or its description object.

Value
(character(1)).

See Also
Other task: getTaskClassLevels(), getTaskCosts(), getTaskData(), getTaskDesc(), getTaskFeatureNames(), getTaskFormula(), getTaskNFeats(), getTaskSize(), getTaskTargetNames(), getTaskTargets(), getTaskType(), subsetTask()

getTaskNFeats
Get number of features in task.

Description
See title.

Usage
getTaskNFeats(x)

Arguments
x (Task | TaskDesc)
Task or its description object.

Value
(integer(1)).
getTaskSize

Get number of observations in task.

Description
See title.

Usage
getTaskSize(x)

Arguments
x (Task | TaskDesc)
Task or its description object.

Value
(integer(1)).

See Also
Other task: getTaskClassLevels(), getTaskCosts(), getTaskData(), getTaskDesc(), getTaskFeatureNames(), getTaskFormula(), getTaskId(), getTaskSize(), getTaskTargetNames(), getTaskTargets(), getTaskType(), subsetTask()

getTaskTargetNames
Get the name(s) of the target column(s).

Description
NB: For multilabel, getTaskTargetNames and getTaskClassLevels actually return the same thing.

Usage
getTaskTargetNames(x)

Arguments
x (Task | TaskDesc)
Task or its description object.
getTaskTargets

Value

(character).

See Also

Other task: getTaskClassLevels(), getTaskCosts(), getTaskData(), getTaskDesc(), getTaskFeatureNames(), getTaskFormula(), getTaskId(), getTaskNFeats(), getTaskSize(), getTaskTargets(), getTaskType(), subsetTask()

getTaskTargets  Get target data of task.

Description

Get target data of task.

Usage

getTaskTargets(task, recode.target = "no")

Arguments

  task  (Task)
  The task.

  recode.target  (character(1))
  Should target classes be recoded? Supported are binary and multilabel classification and survival. Possible values for binary classification are "01", "-1+1" and "drop.levels". In the two latter cases the target vector is converted into a numeric vector. The positive class is coded as "+1" and the negative class either as "0" or "-1". "drop.levels" will remove empty factor levels in the target column. In the multilabel case the logical targets can be converted to factors with "multilabel.factor". For survival, you may choose to recode the survival times to "left", "right" or "interval2" censored times using "lcens", "rcens" or "icens", respectively. See survival::Surv for the format specification. Default for both binary classification and survival is "no" (do nothing).

Value

A factor for classification or a numeric for regression, a data.frame of logical columns for multilabel.

See Also

Other task: getTaskClassLevels(), getTaskCosts(), getTaskData(), getTaskDesc(), getTaskFeatureNames(), getTaskFormula(), getTaskId(), getTaskNFeats(), getTaskSize(), getTaskTargetNames(), getTaskType(), subsetTask()
getTuneResult

Returns the optimal hyperparameters and optimization path after training.

Description
Returns the optimal hyperparameters and optimization path after training.

Usage
getTuneResult(object)

Arguments
object (WrappedModel)
Trained Model created with makeTuneWrapper.
**getTuneResultOptPath**

*Get the optimization path of a tuning result.*

**Value**

*(TuneResult).*

**See Also**

Other tune: `TuneControl`, `getNestedTuneResultsOptPathDf()`, `getNestedTuneResultsX()`, `getResamplingIndices()`, `makeModelMultiplexerParamSet()`, `makeModelMultiplexer()`, `makeTuneControlCMAES()`, `makeTuneControlDesign()`, `makeTuneControlGenSA()`, `makeTuneControlGrid()`, `makeTuneControlIrace()`, `makeTuneControlMBO()`, `makeTuneControlRandom()`, `makeTuneWrapper()`, `tuneParams()`, `tuneThreshold()`

---

**Description**

Returns the opt.path from a *(TuneResult)* object.

**Usage**

`getTuneResultOptPath(tune.result, as.df = TRUE)`

**Arguments**

- `tune.result` *(TuneResult)*
  
  A tuning result of the *(tuneParams)* function.

- `as.df` *(logical(1))*
  
  Should the optimization path be returned as a data frame? Default is *TRUE*.

**Value**

*(ParamHelpers::OptPath)* or *(data.frame)*.

---

**gunpoint.task**

*Gunpoint functional data classification task.*

**Description**

Contains the task *(gunpoint.task)*. You have to classify whether a person raises up a gun or just an empty hand.

**References**

hasFunctionalFeatures  
*Check whether the object contains functional features.*

**Description**

See title.

**Usage**

```r
hasFunctionalFeatures(obj)
```

**Arguments**

- `obj` *(Task | TaskDesc | data.frame)*
  
  Object to check.

**Value**

```r
logical(1)
```

---

**hasProperties**  
*Deprecated, use hasLearnerProperties instead.*

**Description**

Deprecated, use hasLearnerProperties instead.

**Usage**

```r
hasProperties(learner, props)
```

**Arguments**

- `learner`  
  
  Deprecated.

- `props`  
  
  Deprecated.
helpLearner

Access help page of learner functions.

Description

Interactive function that gives the user quick access to the help pages associated with various functions involved in the given learner.

Usage

helpLearner(learner)

Arguments

learner (Learner | character(1))

The learner. If you pass a string the learner will be created via makeLearner.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(),
getLearnerPredictType(), getLearnerShortName(), getLearnerType(),getParamSet(), helpLearnerParam(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()

Other help: helpLearnerParam()

helpLearnerParam

Get specific help for a learner’s parameters.

Description

Print the description of parameters of a given learner. The description is automatically extracted from the help pages of the learner, so it may be incomplete.

Usage

helpLearnerParam(learner, param = NULL)

Arguments

learner (Learner | character(1))

The learner. If you pass a string the learner will be created via makeLearner.

param (character | NULL)

Parameter(s) to describe. Defaults to NULL, which prints information on the documentation status of all parameters.
See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(),
getLearnerPredictType(), getLearnerShortName(), getLearnerType(),getParamSet(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()
Other help: helpLearner()

---

**imputations**

Built-in imputation methods.

**Description**

The built-ins are:

- `imputeConstant(const)` for imputation using a constant value,
- `imputeMedian()` for imputation using the median,
- `imputeMode()` for imputation using the mode,
- `imputeMin(multiplier)` for imputing constant values shifted below the minimum using
  \[ \text{min}(x) - \text{multiplier} \times \text{diff(range}(x)) \],
- `imputeMax(multiplier)` for imputing constant values shifted above the maximum using
  \[ \text{max}(x) + \text{multiplier} \times \text{diff(range}(x)) \],
- `imputeNormal(mean, sd)` for imputation using normally distributed random values. Mean
  and standard deviation will be calculated from the data if not provided.
- `imputeHist(breaks, use.mids)` for imputation using random values with probabilities calculated
  using `table` or `hist`.
- `imputeLearner(learner, features = NULL)` for imputations using the response of a classification
  or regression learner.

**Usage**

```r
imputeConstant(const)
imputeMedian()
imputeMean()
imputeMode()
imputeMin(multiplier = 1)
imputeMax(multiplier = 1)
imputeUniform(min = NA_real_, max = NA_real_)
```
imputeNormal(mu = NA_real_, sd = NA_real_)
imputeHist(breaks, use.mids = TRUE)
imputeLearner(learner, features = NULL)

Arguments

const (any)
Constant valued use for imputation.

multiplier (numeric(1))
Value that stored minimum or maximum is multiplied with when imputation is done.

min (numeric(1))
Lower bound for uniform distribution. If NA (default), it will be estimated from the data.

max (numeric(1))
Upper bound for uniform distribution. If NA (default), it will be estimated from the data.

mu (numeric(1))
Mean of normal distribution. If missing it will be estimated from the data.

sd (numeric(1))
Standard deviation of normal distribution. If missing it will be estimated from the data.

breaks (numeric(1))
Number of breaks to use in graphics::hist. If missing, defaults to auto-detection via “Sturges”.

use.mids (logical(1))
If x is numeric and a histogram is used, impute with bin mids (default) or instead draw uniformly distributed samples within bin range.

learner (Learner | character(1))
Supervised learner. Its predictions will be used for imputations. If you pass a string the learner will be created via makeLearner. Note that the target column is not available for this operation.

features (character)
Features to use in learner for prediction. Default is NULL which uses all available features except the target column of the original task.

See Also

Other impute: impute(), makeImputeMethod(), makeImputeWrapper(), reimpute()
impute

Impute and re-impute data

Description

Allows imputation of missing feature values through various techniques. Note that you have the possibility to re-impute a data set in the same way as the imputation was performed during training. This especially comes in handy during resampling when one wants to perform the same imputation on the test set as on the training set.

The function `impute` performs the imputation on a data set and returns, alongside with the imputed data set, an “ImputationDesc” object which can contain “learned” coefficients and helpful data. It can then be passed together with a new data set to `reimpute`.

The imputation techniques can be specified for certain features or for feature classes, see function arguments.

You can either provide an arbitrary object, use a built-in imputation method listed under `imputations` or create one yourself using `makeImputeMethod`.

Usage

```r
impute(
  obj,
  target = character(0L),
  classes = list(),
  cols = list(),
  dummy.classes = character(0L),
  dummy.cols = character(0L),
  dummy.type = "factor",
  force.dummies = FALSE,
  impute.new.levels = TRUE,
  recode.factor.levels = TRUE
)
```

Arguments

- `obj` (data.frame | Task): Input data.
- `target` (character): Name of the column(s) specifying the response. Default is character(0).
- `classes` (named list): Named list containing imputation techniques for classes of columns. E.g. `list(numeric = imputeMedian())`.
- `cols` (named list): Named list containing names of imputation methods to impute missing values in the data column referenced by the list element’s name. Overrules imputation set via `classes`.
dummy.classes (character)  
Classes of columns to create dummy columns for. Default is character(0).

dummy.cols (character)  
Column names to create dummy columns (containing binary missing indicator) for. Default is character(0).

dummy.type (character(1))  
How dummy columns are encoded. Either as 0/1 with type “numeric” or as “factor”. Default is “factor”.

force.dummies (logical(1))  
Force dummy creation even if the respective data column does not contain any NAs. Note that (a) most learners will complain about constant columns created this way but (b) your feature set might be stochastic if you turn this off. Default is FALSE.

impute.new.levels  
(logical(1))  
If new, unencountered factor level occur during reimputation, should these be handled as NAs and then be imputed the same way? Default is TRUE.

recode.factor.levels  
(logical(1))  
Recode factor levels after reimputation, so they match the respective element of lvls (in the description object) and therefore match the levels of the feature factor in the training data after imputation?. Default is TRUE.

Details

The description object contains these slots

target (character)  See argument.
features (character)  Feature names (column names of data).
classes (character)  Feature classes (storage type of data).
lvls (named list)  Mapping of column names of factor features to their levels, including newly created ones during imputation.
impute (named list)  Mapping of column names to imputation functions.
dummies (named list)  Mapping of column names to imputation functions.
impute.new.levels (logical(1))  See argument.
recode.factor.levels (logical(1))  See argument.

Value

(list)

data (data.frame)  
Imputed data.
desc (ImputationDesc)  
Description object.
### `isFailureModel`

#### Description

Such a model is created when one sets the corresponding option in `configureMlr`. For complex wrappers this getter returns `TRUE` if ANY model contained in it failed.

#### Usage

```r
isFailureModel(model)
```

#### Arguments

- `model` *(WrappedModel)*
  
  The model.

#### Value

`logical(1)`.
**joinClassLevels**

Join some class existing levels to new, larger class levels for classification problems.

**Description**

Join some class existing levels to new, larger class levels for classification problems.

**Usage**

```r
joinClassLevels(task, new.levels)
```

**Arguments**

- `task` *(Task)*
  - The task.

- `new.levels` *(list of character)*
  - Element names specify the new class levels to create, while the corresponding element character vector specifies the existing class levels which will be joined to the new one.

**Value**

- `Task`.

**Examples**

```r
joinClassLevels(iris.task, new.levels = list(foo = c("setosa", "virginica")))
```

**learnerArgsToControl**

Convert arguments to control structure.

**Description**

Find all elements in ... which are not missing and call `control` on them.

**Usage**

```r
learnerArgsToControl(control, ...)
```

**Arguments**

- `control` *(function)*
  - Function that creates control structure.

- `...` *(any)*
  - Arguments for control structure function.
Value

Control structure for learner.

Description

Properties can be accessed with `getLearnerProperties(learner)`, which returns a character vector.

The learner properties are defined as follows:

- **numerics, factors, ordered** Can numeric, factor or ordered factor features be handled?
- **functionals** Can an arbitrary number of functional features be handled?
- **single.functional** Can exactly one functional feature be handled?
- **missings** Can missing values in features be handled?
- **weights** Can observations be weighted during fitting?
- **oneclas, twoclass, multiclass** Only for classif: Can one-class, two-class or multi-class classification problems be handled?
- **class.weights** Only for classif: Can class weights be handled?
- **rcens, lcens, icens** Only for surv: Can right, left, or interval censored data be handled?
- **prob** For classif, cluster, multilabel, surv: Can probabilites be predicted?
- **se** Only for regr: Can standard errors be predicted?
- **oobpreds** Only for classif, regr and surv: Can out of bag predictions be extracted from the trained model?
- **featimp** For classif, regr, surv: Does the model support extracting information on feature importance?

Usage

- `getLearnerProperties(learner)`
- `hasLearnerProperties(learner, props)`

Arguments

- **learner** *(Learner | character(1))*
  The learner. If you pass a string the learner will be created via `makeLearner`.
- **props** *(character)*
  Vector of properties to query.
Value
getLearnerProperties returns a character vector with learner properties. hasLearnerProperties
returns a logical vector of the same length as props.

See Also
Other learner: getClassWeightParam(), getHyperPars(), getLearnerId(), getLearnerNote(),
getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(),
getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(),
makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(),
setPredictThreshold(), setPredictType()

learners
List of supported learning algorithms.

Description
All supported learners can be found by listLearners or as a table in the tutorial appendix: https://mlr.mlr-org.com/articles/tutorial/integrated_learners.html.

listFilterEnsembleMethods
List ensemble filter methods.

Description
Returns a subset-able dataframe with filter information.

Usage
listFilterEnsembleMethods(desc = TRUE)

Arguments
  desc (logical(1))
  Provide more detailed information about filters. Default is TRUE.

Value
(data.frame).

See Also
Other filter: filterFeatures(), generateFilterValuesData(), getFilteredFeatures(), listFilterMethods(),
makeFilterEnsemble(), makeFilterWrapper(), makeFilter(), plotFilterValues()
listFilterMethods  List filter methods.

Description
Returns a subset-able dataframe with filter information.

Usage
listFilterMethods(
  desc = TRUE,
  tasks = FALSE,
  features = FALSE,
  include.deprecated = FALSE
)

Arguments
desc (logical(1))
  Provide more detailed information about filters. Default is TRUE.

tasks (logical(1))
  Provide information on supported tasks. Default is FALSE.

features (logical(1))
  Provide information on supported features. Default is FALSE.

include.deprecated (logical(1))
  Should deprecated filter methods be included in the list. Default is FALSE.

Value
(data.frame).

See Also
Other filter: filterFeatures(), generateFilterValuesData(), getFilteredFeatures(), listFilterEnsembleMethods(), makeFilterEnsemble(), makeFilterWrapper(), makeFilter(), plotFilterValues()
listLearnerProperties  
List the supported learner properties

Description
This is useful for determining which learner properties are available.

Usage
listLearnerProperties(type = "any")

Arguments

  type (character(1))
  Only return properties for a specified task type. Default is “any”.

Value
(character).

listLearners  
Find matching learning algorithms.

Description
Returns learning algorithms which have specific characteristics, e.g. whether they support missing values, case weights, etc.

Note that the packages of all learners are loaded during the search if you create them. This can be a lot. If you do not create them we only inspect properties of the S3 classes. This will be a lot faster.

Note that for general cost-sensitive learning, mlr currently supports mainly “wrapper” approaches like CostSensWeightedPairsWrapper, which are not listed, as they are not basic R learning algorithms. The same applies for many multilabel methods, see, e.g., makeMultilabelBinaryRelevanceWrapper.

Usage
listLearners(
  obj = NA_character_,
  properties = character(0L),
  quiet = TRUE,
  warn.missing.packages = TRUE,
  check.packages = FALSE,
  create = FALSE
)
## Default S3 method:

```r
listLearners(
  obj = NA_character_,
  properties = character(0L),
  quiet = TRUE,
  warn.missing.packages = TRUE,
  check.packages = FALSE,
  create = FALSE
)
```

## S3 method for class 'character'

```r
listLearners(
  obj = NA_character_,
  properties = character(0L),
  quiet = TRUE,
  warn.missing.packages = TRUE,
  check.packages = FALSE,
  create = FALSE
)
```

## S3 method for class 'Task'

```r
listLearners(
  obj = NA_character_,
  properties = character(0L),
  quiet = TRUE,
  warn.missing.packages = TRUE,
  check.packages = TRUE,
  create = FALSE
)
```

### Arguments

#### obj
(character(1) | Task)
Either character(1) task or the type of the task, in the latter case one of: “classif” “regr” “surv” “costsens” “cluster” “multilabel”. Default is NA matching all types.

#### properties
(character)
Set of required properties to filter for. Default is character(0).

#### quiet
(logical(1))
Construct learners quietly to check their properties, shows no package startup messages. Turn off if you suspect errors. Default is TRUE.

#### warn.missing.packages
(logical(1))
If some learner cannot be constructed because its package is missing, should a warning be shown? Default is TRUE.

#### check.packages
(logical(1))
Check if required packages are installed. Calls find.package(). If create is TRUE, this is done implicitly and the value of this parameter is ignored. If
create is FALSE and check.packages is TRUE the returned table only contains learners whose dependencies are installed. If check.packages set to FALSE, learners that cannot actually be constructed because of missing packages may be returned. Default is FALSE.

create

(logical(1))

Instantiate objects (or return info table)? Packages are loaded if and only if this option is TRUE. Default is FALSE.

Value

(data.frame | list' of Learner). Either a descriptive data.frame that allows access to all properties of the learners or a list of created learner objects (named by ids of listed learners).

Examples

## Not run:
listLearners("classif", properties = c("multiclass", "prob"))
data = iris
task = makeClassifTask(data = data, target = "Species")
listLearners(task)

## End(Not run)
listMeasures  

*Find matching measures.*

**Description**

Returns the matching measures which have specific characteristics, e.g. whether they supports classification or regression.

**Usage**

```
listMeasures(obj, properties = character(0L), create = FALSE)
```

```{}
## Default S3 method:
listMeasures(obj, properties = character(0L), create = FALSE)

## S3 method for class 'character'
listMeasures(obj, properties = character(0L), create = FALSE)

## S3 method for class 'Task'
listMeasures(obj, properties = character(0L), create = FALSE)
```

**Arguments**

- **obj** (character(1) | Task)
  Either character(1) task or the type of the task, in the latter case one of: “classif” “regr” “surv” “costsens” “cluster” “multilabel”. Default is NA matching all types.

- **properties** (character)
  Set of required properties to filter for. See Measure for some standardized properties. Default is character(0).

- **create** (logical(1))
  Instantiate objects (or return strings)? Default is FALSE.

**Value**

( character | list’ of Measure). Class names of matching measures or instantiated objects.

---

listTaskTypes  

*List the supported task types in mlr*

**Description**

Returns a character vector with each of the supported task types in mlr.
Usage

listTaskTypes()

Value

(character).

lung.task

NCCTG Lung Cancer survival task.

Description

Contains the task (lung.task).

References

See survival::lung. Incomplete cases have been removed from the task.

makeAggregation

Specify your own aggregation of measures.

Description

This is an advanced feature of mlr. It gives access to some inner workings so the result might not be compatible with everything!

Usage

makeAggregation(id, name = id, properties, fun)

Arguments

id (character(1))
Name of the aggregation method (preferably the same name as the generated function).

name (character(1))
Long name of the aggregation method. Default is id.

properties (character)
Set of aggregation properties.

req.train Are prediction or train sets required to calculate the aggregation?

req.test Are prediction or test sets required to calculate the aggregation?

fun (function(task, perf.test, perf.train, measure, group, pred))
Calculates the aggregated performance. In most cases you will only need the performances perf.test and optionally perf.train on the test and training data sets.
task (Task) The task.
perf.test (numeric) performance results on the test data sets.
perf.train (numeric) performance results on the training data sets.
measure (Measure) Performance measure.
group (factor) Grouping of resampling iterations. This encodes whether specific iterations 'belong together' (e.g. repeated CV).
pred (Prediction) Prediction object.

Value

(Aggregation).

See Also

aggregations, setAggregation

Examples

# computes the interquartile range on all performance values
test.iqr = makeAggregation(
  id = "test.iqr", name = "Test set interquartile range",
  properties = "req.test",
  fun = function(task, perf.test, perf.train, measure, group, pred) IQR(perf.test)
)

makeBaggingWrapper

Fuse learner with the bagging technique.

Description

Fuses a learner with the bagging method (i.e., similar to what a randomForest does). Creates a learner object, which can be used like any other learner object. Models can easily be accessed via getLearnerModel.

Bagging is implemented as follows: For each iteration a random data subset is sampled (with or without replacement) and potentially the number of features is also restricted to a random subset. Note that this is usually handled in a slightly different way in the random forest where features are sampled at each tree split.

Prediction works as follows: For classification we do majority voting to create a discrete label and probabilities are predicted by considering the proportions of all predicted labels. For regression the mean value and the standard deviations across predictions is computed. Note that the passed base learner must always have predict.type = 'response', while the BaggingWrapper can estimate probabilities and standard errors, so it can be set, e.g., to predict.type = 'prob'. For this reason, when you call setPredictType, the type is only set for the BaggingWrapper, not passed down to the inner learner.
Usage

makeBaggingWrapper(
  learner,
  bw.iters = 10L,
  bw.replace = TRUE,
  bw.size,
  bw.feats = 1
)

Arguments

learner  (Learner | character(1))
  The learner. If you pass a string the learner will be created via makeLearner.

bw.iters  (integer(1))
  Iterations = number of fitted models in bagging. Default is 10.

bw.replace  (logical(1))
  Sample bags with replacement (bootstrapping)? Default is TRUE.

bw.size  (numeric(1))
  Percentage size of sampled bags. Default is 1 for bootstrapping and 0.632 for subsampling.

bw.feats  (numeric(1))
  Percentage size of randomly selected features in bags. Default is 1. At least one feature will always be selected.

Value

Learner.

See Also

Other wrapper: makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(),
makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(),
makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(),
makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(),
makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(),
makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(),
makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(),
makeUndersampleWrapper(), makeWeightedClassesWrapper()
Description
Builds regression models that predict for the positive class whether a particular example belongs to it (1) or not (-1).
Probabilities are generated by transforming the predictions with a softmax.
Inspired by WEKA's ClassificationViaRegression (http://weka.sourceforge.net/doc.dev/weka/classifiers/meta/ClassificationViaRegression.html).

Usage
makeClassificationViaRegressionWrapper(learner, predict.type = "response")

Arguments
learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.
predict.type (character(1))
"response" (= labels) or "prob" (= probabilities and labels by selecting the one with maximal probability).

Value
Learner.

See Also
Other wrapper: makeBaggingWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()

Examples
lrn = makeLearner("regr.rpart")
lrn = makeClassificationViaRegressionWrapper(lrn)
mod = train(lrn, sonar.task, subset = 1:140)
predictions = predict(mod, newdata = getTaskData(sonar.task)[141:208, 1:60])
makeClassifTask

Usage

makeClassifTask(
  id = deparse(substitute(data)),
  data,
  target,
  weights = NULL,
  blocking = NULL,
  coordinates = NULL,
  positive = NA_character_,
  fixup.data = "warn",
  check.data = TRUE
)

Arguments

id (character(1)) Id string for object. Default is the name of the R variable passed to data.
data (data.frame) A data frame containing the features and target variable(s).
target (character(1) | character(2) | character(n.classes)) Name(s) of the target variable(s). For survival analysis these are the names of the survival time and event columns, so it has length 2. For multilabel classification it contains the names of the logical columns that encode whether a label is present or not and its length corresponds to the number of classes.
weights (numeric) Optional, non-negative case weight vector to be used during fitting. Cannot be set for cost-sensitive learning. Default is NULL which means no (= equal) weights.
blocking (factor) An optional factor of the same length as the number of observations. Observations with the same blocking level “belong together”. Specifically, they are either put all in the training or the test set during a resampling iteration. Default is NULL which means no blocking.
coordinates (data.frame) Coordinates of a spatial data set that will be used for spatial partitioning of the data in a spatial cross-validation resampling setting. Coordinates have to be numeric values. Provided data.frame needs to have the same number of rows as data and consist of at least two dimensions.
positive (character(1)) Positive class for binary classification (otherwise ignored and set to NA). Default is the first factor level of the target attribute.
fixup.data (character(1)) Should some basic cleaning up of data be performed? Currently this means removing empty factor levels for the columns. Possible choices are: “no” = Don’t do it. “warn” = Do it but warn about it. “quiet” = Do it but keep silent. Default is “warn”.
check.data (logical(1))
Should sanity of data be checked initially at task creation? You should have
good reasons to turn this off (one might be speed). Default is TRUE.

See Also
Task CostSensTask ClusterTask MultilabelTask RegrTask SurvTask

makeClusterTask Create a cluster task.

Description
Create a cluster task.

Usage
makeClusterTask(
  id = deparse(substitute(data)),
  data,
  weights = NULL,
  blocking = NULL,
  coordinates = NULL,
  fixup.data = "warn",
  check.data = TRUE
)

Arguments
id (character(1))
Id string for object. Default is the name of the R variable passed to data.
data (data.frame)
A data frame containing the features and target variable(s).
weights (numeric)
Optional, non-negative case weight vector to be used during fitting. Cannot
be set for cost-sensitive learning. Default is NULL which means no (= equal)
weights.
blocking (factor)
An optional factor of the same length as the number of observations. Observ-
ations with the same blocking level “belong together”. Specifically, they are
either put all in the training or the test set during a resampling iteration. Default
is NULL which means no blocking.
coordinates (data.frame)
Coordinates of a spatial data set that will be used for spatial partitioning of the
data in a spatial cross-validation resampling setting. Coordinates have to be
numeric values. Provided data.frame needs to have the same number of rows as
data and consist of at least two dimensions.
Should some basic cleaning up of data be performed? Currently this means removing empty factor levels for the columns. Possible choices are: “no” = Don’t do it. “warn” = Do it but warn about it. “quiet” = Do it but keep silent. Default is “warn”.

Should sanity of data be checked initially at task creation? You should have good reasons to turn this off (one might be speed). Default is TRUE.

See Also

Task ClassifTask CostSensTask MultilabelTask RegrTask SurvTask

makeConstantClassWrapper

Wraps a classification learner to support problems where the class label is (almost) constant.

Description

If the training data contains only a single class (or almost only a single class), this wrapper creates a model that always predicts the constant class in the training data. In all other cases, the underlying learner is trained and the resulting model used for predictions.

Probabilities can be predicted and will be 1 or 0 depending on whether the label matches the majority class or not.

Usage

makeConstantClassWrapper(learner, frac = 0)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

frac numeric(1)
The fraction of labels in \([0, 1)\) that can be different from the majority label. Default is 0, which means that constant labels are only predicted if there is exactly one label in the data.

Value

Learner.
makeCostMeasure

Creates a measure for non-standard misclassification costs.

Description

Creates a cost measure for non-standard classification error costs.

Usage

makeCostMeasure(
  id = "costs",
  minimize = TRUE,
  costs,
  combine = mean,
  best = NULL,
  worst = NULL,
  name = id,
  note = ""
)

Arguments

id (character(1))
Name of measure. Default is “costs”.

minimize (logical(1))
Should the measure be minimized? Otherwise you are effectively specifying a benefits matrix. Default is TRUE.

costs (matrix)
Matrix of misclassification costs. Rows and columns have to be named with class labels, order does not matter. Rows indicate true classes, columns predicted classes.

combine (function)
How to combine costs over all cases for a SINGLE test set? Note this is not the same as the aggregate argument in makeMeasure You can set this as well via setAggregation, as for any measure. Default is mean.
**makeCostSensClassifWrapper**

Wraps a classification learner for use in cost-sensitive learning.

**Description**

Creates a wrapper, which can be used like any other learner object. The classification model can easily be accessed via `getLearnerModel`.

This is a very naive learner, where the costs are transformed into classification labels - the label for each case is the name of class with minimal costs. (If ties occur, the label which is better on average w.r.t. costs over all training data is preferred.) Then the classifier is fitted to that data and subsequently used for prediction.

**Usage**

```r
makeCostSensClassifWrapper(learner)
```

**Arguments**

- **learner**
  
  (Learner | character(1))
  
  The classification learner. If you pass a string the learner will be created via `makeLearner`.

**Value**

- **Learner**
makeCostSensRegrWrapper

Wraps a regression learner for use in cost-sensitive learning.

Description

Creates a wrapper, which can be used like any other learner object. Models can easily be accessed via getLearnerModel.

For each class in the task, an individual regression model is fitted for the costs of that class. During prediction, the class with the lowest predicted costs is selected.

Usage

makeCostSensRegrWrapper(learner)

Arguments

learner (Learner | character(1))

The regression learner. If you pass a string the learner will be created via makeLearner.

Value

Learner.

See Also

Other costsens: makeCostSensClassifWrapper(), makeCostSensTask(), makeCostSensWeightedPairsWrapper()

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()
makeCostSensTask

Create a cost-sensitive classification task.

Description

Create a cost-sensitive classification task.

Usage

makeCostSensTask(
  id = deparse(substitute(data)),
  data,
  costs,
  blocking = NULL,
  coordinates = NULL,
  fixup.data = "warn",
  check.data = TRUE
)

Arguments

id (character(1))
Id string for object. Default is the name of the R variable passed to data.

data (data.frame)
A data frame containing the features and target variable(s).

costs (data.frame)
A numeric matrix or data frame containing the costs of misclassification. We assume the general case of observation specific costs. This means we have n rows, corresponding to the observations, in the same order as data. The columns correspond to classes and their names are the class labels (if unnamed we use y1 to yk as labels). Each entry (i,j) of the matrix specifies the cost of predicting class j for observation i.

blocking (factor)
An optional factor of the same length as the number of observations. Observations with the same blocking level “belong together”. Specifically, they are either put all in the training or the test set during a resampling iteration. Default is NULL which means no blocking.

coordinates (data.frame)
Coordinates of a spatial data set that will be used for spatial partitioning of the data in a spatial cross-validation resampling setting. Coordinates have to be numeric values. Provided data.frame needs to have the same number of rows as data and consist of at least two dimensions.

fixup.data (character(1))
Should some basic cleaning up of data be performed? Currently this means removing empty factor levels for the columns. Possible choices are: “no” =
Don’t do it. “warn” = Do it but warn about it. “quiet” = Do it but keep silent. Default is “warn”.

check.data (logical(1))
Should sanity of data be checked initially at task creation? You should have good reasons to turn this off (one might be speed). Default is TRUE.

See Also
Task ClassifTask ClusterTask MultilabelTask RegrTask SurvTask
Other costsens: makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeCostSensWeightedPairsWrapper()

makeCostSensWeightedPairsWrapper
Wraps a classifier for cost-sensitive learning to produce a weighted pairs model.

Description
Creates a wrapper, which can be used like any other learner object. Models can easily be accessed via getLearnerModel.

For each pair of labels, we fit a binary classifier. For each observation we define the label to be the element of the pair with minimal costs. During fitting, we also weight the observation with the absolute difference in costs. Prediction is performed by simple voting.

This approach is sometimes called cost-sensitive one-vs-one (CS-OVO), because it is obviously very similar to the one-vs-one approach where one reduces a normal multi-class problem to multiple binary ones and aggregates by voting.

Usage
makeCostSensWeightedPairsWrapper(learner)

Arguments
learner (Learner | character(1))
The classification learner. If you pass a string the learner will be created via makeLearner.

Value
(Learner).

References
**makeCustomResampledMeasure**

Construct your own resampled performance measure.

**Description**

Construct your own performance measure, used after resampling. Note that individual training / test set performance values will be set to ‘NA’, you only calculate an aggregated value. If you can define a function that makes sense for every single training / test set, implement your own [Measure].

**Usage**

```r
makeCustomResampledMeasure(
  measure.id,
  aggregation.id,
  minimize = TRUE,
  properties = character(0L),
  fun,
  extra.args = list(),
  best = NULL,
  worst = NULL,
  measure.name = measure.id,
  aggregation.name = aggregation.id,
  note = ""
)
```

**Arguments**

- `measure.id` (`'character(1)'`)
  Short name of measure.

- `aggregation.id` (`'character(1)'`)
  Short name of aggregation.

- `minimize` (`'logical(1)'`)
  Should the measure be minimized? Default is ‘TRUE’.

- `properties` (`[character]`)
  Set of measure properties. For a list of values see [Measure]. Default is ‘character(0)’.

- `fun` (`'function(task, group, pred, extra.args)'`)
  Calculates performance value from [ResamplePrediction] object. For rare cases you can also use the task, the grouping or the extra arguments ‘extra.args’.

  - `task` ([Task ]): The task.
makeDownsampleWrapper

‘group’ ([factor]) Grouping of resampling iterations. This encodes whether specific iterations ‘belong together’ (e.g. repeated CV).

‘pred’ ([Prediction]) Prediction object.
‘extra.args’ ([list]) See below.

extra.args ([list])
List of extra arguments which will always be passed to ‘fun’. Default is empty list.

best (‘numeric(1)’) Best obtainable value for measure. Default is ‘-Inf’ or ‘Inf’, depending on ‘minimize’.

worst (‘numeric(1)’) Worst obtainable value for measure. Default is ‘Inf’ or ‘-Inf’, depending on ‘minimize’.

measure.name (‘character(1)’) Long name of measure. Default is ‘measure.id’.

aggregation.name (‘character(1)’) Long name of the aggregation. Default is ‘aggregation.id’.

note ([character]) Description and additional notes for the measure. Default is “”.

Value

Measure.

See Also

Other performance: ConfusionMatrix, calculateConfusionMatrix(), calculateROCMeasures(), estimateRelativeOverfitting(), makeCostMeasure(), makeMeasure(), measures, performance(), setAggregation(), setMeasurePars()

makeDownsampleWrapper  Fuse learner with simple downsampling (subsampling).

Description

Creates a learner object, which can be used like any other learner object. It will only be trained on a subset of the original data to save computational time.

Usage

makeDownsampleWrapper(learner, dw.perc = 1, dw.stratify = FALSE)
**makeDummyFeaturesWrapper**

_Fuse learner with dummy feature creator._

**Description**

Fuses a base learner with the dummy feature creator (see `createDummyFeatures`). Returns a learner which can be used like any other learner.

**Usage**

```r
makeDummyFeaturesWrapper(learner, method = "1-of-n", cols = NULL)
```

**Arguments**

- **learner** (`Learner | character(1)`)
  The learner. If you pass a string the learner will be created via `makeLearner`.

- **method** (`character(1)`)
  Available are:
  
  - "1-of-n": For n factor levels there will be n dummy variables.
  - "reference": There will be n-1 dummy variables leaving out the first factor level of each variable.

- **cols** (`numeric(1)`)
  See `downsample`. Default is 1.
makeExtractFDAFeatMethod

Constructor for FDA feature extraction methods.

Description

This can be used to implement custom FDA feature extraction. Takes a learn and a reextract function along with some optional parameters to those as argument.

Usage

makeExtractFDAFeatMethod(learn, reextract, args = list(), par.set = NULL)

Arguments

learn

(fixed) (function(data, target, col, ...))

Function to learn and extract information on functional column col. Arguments are:

• data data.frame

Data.frame containing matrices with one row per observation of a single functional or time series and one column per measurement time point. All entries need to be numeric.

• target (character(1))

Name of the target variable. Default: “NULL”. The variable is only set to be consistent with the API.

• col (character(1) | numeric(1))

column names or indices, the extraction should be performed on. The function has to return a named list of values.

Value

Learner.

See Also

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTETRansformer(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()
**makeExtractFDAFeatsWrapper**

Fuse learner with an extractFDAFeatures method.

### Description

Fuses a base learner with an extractFDAFeatures method. Creates a learner object, which can be used like any other learner object. Internally uses `extractFDAFeatures` before training the learner and `reextractFDAFeatures` before predicting.

### Usage

```r
makeExtractFDAFeatsWrapper(learner, feat.methods = list())
```

### Arguments

- **learner** *(Learner | character(1))*
  
The learner. If you pass a string the learner will be created via `makeLearner`.

- **feat.methods** *(named list)*
  
  List of functional features along with the desired methods for each functional feature. “all” applies the `extractFDAFeatures` method to each functional feature. Names of `feat.methods` must match column names of functional features. Available feature extraction methods are available under family `fda_featextractor`. Specifying a functional feature multiple times with different extraction methods allows for the extraction of different features from the same functional. Default is `list()` which does nothing.

### Value

Learner.
Description

Fuses a base learner with a search strategy to select variables. Creates a learner object, which
can be used like any other learner object, but which internally uses \selectFeatures. If the train
function is called on it, the search strategy and resampling are invoked to select an optimal set of
variables. Finally, a model is fitted on the complete training data with these variables and returned.
See \selectFeatures for more details.

After training, the optimal features (and other related information) can be retrieved with \getFeat-
SelResult.

Usage

makeFeatSelWrapper(
  learner,
  resampling,
  measures,
  bit.names,
  bits.to.features,
  control,
  show.info = getMlrOption("show.info")
)

Arguments

learner  (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

resampling  ([ResampleInstance] | [ResampleDesc])
Resampling strategy for feature selection. If you pass a description, it is instan-
tiated once at the beginning by default, so all points are evaluated on the same
training/test sets. If you want to change that behaviour, look at [FeatSelControl].
makeFeatSelWrapper

measures
(list of Measure | Measure)
Performance measures to evaluate. The first measure, aggregated by the first
aggregation function is optimized, others are simply evaluated. Default is the
default measure for the task, see here getDefaultMeasure.

bit.names
[character]
Names of bits encoding the solutions. Also defines the total number of bits in
the encoding. Per default these are the feature names of the task. Has to be used
together with 'bits.to.features'.

bits.to.features
[function(x, task)]
Function which transforms an integer-0-1 vector into a character vector of se-
lected features. Per default a value of 1 in the ith bit selects the ith feature to be
in the candidate solution. The vector ‘x’ will correspond to the ‘bit.names’ and
has to be of the same length.

control
[see [FeatSelControl]) Control object for search method. Also selects the opti-
mization algorithm for feature selection.

show.info
(logical(1))
Print verbose output on console? Default is set via configureMlr.

Value
Learner.

See Also
Other featsel: FeatSelControl, analyzeFeatSelResult(), getFeatSelResult(), selectFeatures()
Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(),
makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(),
makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFilterWrapper(), makeImputeWrapper(),
makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(),
makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(),
makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(),
makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()

Examples
# nested resampling with feature selection (with a pretty stupid algorithm for selection)
outer = makeResampleDesc("CV", iters = 2L)
inner = makeResampleDesc("Holdout")
ctrl = makeFeatSelControlRandom(maxit = 1)
lrn = makeFeatSelWrapper("classif.ksvm", resampling = inner, control = ctrl)
# we also extract the selected features for all iteration here
r = resample(lrn, iris.task, outer, extract = getFeatSelResult)
makeFilter

Create a feature filter.

Description

Creates and registers custom feature filters. Implemented filters can be listed with listFilterMethods. Additional documentation for the fun parameter specific to each filter can be found in the description.

Usage

makeFilter(name, desc, pkg, supported.tasks, supported.features, fun)

Arguments

name (character(1)) Identifier for the filter.
desc (character(1)) Short description of the filter.
pkg (character(1)) Source package where the filter is implemented.
supported.tasks (character) Task types supported.
supported.features (character) Feature types supported.
fun (function(task, nselect, ...)) Function which takes a task and returns a named numeric vector of scores, one score for each feature of task. Higher scores mean higher importance of the feature. At least nselect features must be calculated, the remaining may be set to NA or omitted, and thus will not be selected. The original order will be restored if necessary.

Value

Object of class “Filter”.

References


makeFilterEnsemble  

Create an ensemble feature filter.

Description

Creates and registers custom ensemble feature filters. Implemented ensemble filters can be listed with listFilterEnsembleMethods. Additional documentation for the fun parameter specific to each filter can be found in the description.

Usage

makeFilterEnsemble(name, base.methods, desc, fun)

Arguments

name (character(1))  
Identifier for the filter.

base.methods  
the base filter methods which the ensemble method will use.

desc (character(1))  
Short description of the filter.

fun (function(task, nselect, ...))  
Function which takes a task and returns a named numeric vector of scores, one score for each feature of task. Higher scores mean higher importance of the feature. At least nselect features must be calculated, the remaining may be set to NA or omitted, and thus will not be selected. the original order will be restored if necessary.

Value

Object of class “FilterEnsemble”.

See Also

Other filter: filterFeatures(), generateFilterValuesData(), getFilteredFeatures(), listFilterEnsembleMethods, listFilterMethods(), makeFilterEnsemble(), makeFilterWrapper(), plotFilterValues()
makeFilterWrapper

Fuse learner with a feature filter method.

Description

Fuses a base learner with a filter method. Creates a learner object, which can be used like any other learner object. Internally uses filterFeatures before every model fit.

Usage

makeFilterWrapper(
  learner,
  fw.method = "randomForestSRC_importance",
  fw.base.methods = NULL,
  fw.perc = NULL,
  fw.abs = NULL,
  fw.threshold = NULL,
  fw.fun = NULL,
  fw.fun.args = NULL,
  fw.mandatory.feat = NULL,
  cache = FALSE,
  ...
)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

fw.method (character(1))
Filter method. See listFilterMethods. Default is “randomForestSRC_importance”.

fw.base.methods (character(1))
Simple Filter methods for ensemble filters. See listFilterMethods. Can only be used in combination with ensemble filters. See listFilterEnsembleMethods.

fw.perc (numeric(1))
If set, select \( \text{fw.perc} \times 100 \) top scoring features. Mutually exclusive with arguments \( \text{fw.abs} \), \( \text{fw.threshold} \) and ‘fw.fun’.

fw.abs (numeric(1))
If set, select \( \text{fw.abs} \) top scoring features. Mutually exclusive with arguments \( \text{fw.perc} \), \( \text{fw.threshold} \) and \( \text{fw.fun} \).

fw.threshold (numeric(1))
If set, select features whose score exceeds \( \text{fw.threshold} \). Mutually exclusive with arguments \( \text{fw.perc} \), \( \text{fw.abs} \) and \( \text{fw.fun} \).

fw.fun (function)
If set, select features via a custom thresholding function, which must return the
number of top scoring features to select. Mutually exclusive with arguments fw.perc, fw.abs and fw.threshold.

fw.fun.args (any)
Arguments passed to the custom thresholding function

fw.mandatory.feat (character)
Mandatory features which are always included regardless of their scores

cache (character(1) | logical)
Whether to use caching during filter value creation. See details.

... (any)
Additional parameters passed down to the filter. If you are using more than one filter method, you need to pass the arguments in a named list via more.args. For example, more.args = list("FSelectorRcpp_information.gain" = list(equal = TRUE)).

Details
If ensemble = TRUE, ensemble feature selection using all methods specified in fw.method is performed. At least two methods need to be selected.

After training, the selected features can be retrieved with getFilteredFeatures.

Note that observation weights do not influence the filtering and are simply passed down to the next learner.

Value
Learner.

Caching
If cache = TRUE, the default mlr cache directory is used to cache filter values. The directory is operating system dependent and can be checked with getCacheDir(). Alternatively a custom directory can be passed to store the cache. The cache can be cleared with deleteCacheDir(). Caching is disabled by default. Care should be taken when operating on large clusters due to possible write conflicts to disk if multiple workers try to write the same cache at the same time.

See Also
Other filter: filterFeatures(), generateFilterValuesData(), getFilteredFeatures(), listFilterEnsembleMethods(), makeFilterEnsemble(), makeFilter(), plotFilterValues()

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()
Examples

```r
makeFixedHoldoutInstance
task = makeClassifTask(data = iris, target = "Species")
lnr = makeLearner("classif.lda")
inner = makeResampleDesc("Holdout")
outer = makeResampleDesc("CV", iters = 2)
lnr = makeFilterWrapper(lnr, fw.perc = 0.5)
mod = train(lnr, task)
print(getFilteredFeatures(mod))

# now nested resampling, where we extract the features that the filter method selected
r = resample(lnr, task, outer, extract = function(model) {
  getFilteredFeatures(model)
})
print(r$extract)

# usage of an ensemble filter
lnr = makeLearner("classif.lda")
lnr = makeFilterWrapper(lnr, fw.method = "E-Borda",
  fw.base.methods = c("FSelectorRcpp_gain.ratio", "FSelectorRcpp_information.gain"),
  fw.perc = 0.5)

r = resample(lnr, task, outer, extract = function(model) {
  getFilteredFeatures(model)
})
print(r$extract)

# usage of a custom thresholding function
biggest_gap = function(values, diff) {
  gap_size = 0
  gap_location = 0
  for (i in (diff + 1):length(values)) {
    gap = values[[i - diff]] - values[[i]]
    if (gap > gap_size) {
      gap_size = gap
      gap_location = i - 1
    }
  }
  return(gap_location)
}

lnr = makeLearner("classif.lda")
lnr = makeFilterWrapper(lnr, fw.method = "randomForestSRC_importance",
  fw.fun = biggest_gap, fw.fun.args = list("diff" = 1))

r = resample(lnr, task, outer, extract = function(model) {
  getFilteredFeatures(model)
})
print(r$extract)
```

makeFixedHoldoutInstance

*Generate a fixed holdout instance for resampling.*
**makeFunctionalData**

**Description**
Generate a fixed holdout instance for resampling.

**Usage**

```r
makeFixedHoldoutInstance(train.inds, test.inds, size)
```

**Arguments**

- `train.inds` (integer)
  Indices for training set.
- `test.inds` (integer)
  Indices for test set.
- `size` (integer(1))
  Size of the data set to resample. The function needs to know the largest possible index of the whole data set.

**Value**

(ResampleInstance).

**makeFunctionalData**

Create a data.frame containing functional features from a normal data.frame.

**Description**
To work with functional features, those features need to be stored as a matrix column in the data.frame, so mlr can automatically recognize them as functional features. This function allows for an easy conversion from a data.frame with numeric columns to the required format. If the data already contains matrix columns, they are left as-is if not specified otherwise in fd.features. See Examples for the structure of the generated output.

**Usage**

```r
makeFunctionalData(data, fd.features = NULL, exclude.cols = NULL)
```

**Arguments**

- `data` (data.frame)
  A data.frame that contains the functional features as numeric columns.
- `fd.features` (list)
  Named list containing integer column indices or character column names. Each element defines a functional feature, in the given order of the indices or column names. The name of the list element defines the name of the functional feature. All selected columns have to correspond to numeric data.frame entries. The default is NULL, which means all numeric features are considered to be a single functional “fd1”.


makeImputeMethod

Description
This is a constructor to create your own imputation methods.

Usage
makeImputeMethod(learn, impute, args = list())

Arguments

  learn (function(data, target, col, ...))
  Function to learn and extract information on column col out of data frame data.
  Argument target specifies the target column of the learning task. The function
  has to return a named list of values.

  impute (function(data, target, col, ...))
  Function to impute missing values in col using information returned by learn on
  the same column. All list elements of the return values of learn are passed to
  this function into ....

  args (list)
  Named list of arguments to pass to learn via ....

See Also
Other impute: imputations.impute(), makeImputeWrapper(), reimpute()
makeImputeWrapper

Fuse learner with an imputation method.

Description
Fuses a base learner with an imputation method. Creates a learner object, which can be used like any other learner object. Internally uses impute before training the learner and reimpute before predicting.

Usage
makeImputeWrapper(
  learner,
  classes = list(),
  cols = list(),
  dummy.classes = character(0L),
  dummy.cols = character(0L),
  dummy.type = "factor",
  force.dummies = FALSE,
  impute.new.levels = TRUE,
  recode.factor.levels = TRUE
)

Arguments

learner  (Learner | character(1))
   The learner. If you pass a string the learner will be created via makeLearner.

classes  (named list)
   Named list containing imputation techniques for classes of columns. E.g. list(numeric = imputeMedian()).

cols  (named list)
   Named list containing names of imputation methods to impute missing values in the data column referenced by the list element’s name. Overrules imputation set via classes.

dummy.classes  (character)
   Classes of columns to create dummy columns for. Default is character(0).

dummy.cols  (character)
   Column names to create dummy columns (containing binary missing indicator) for. Default is character(0).

dummy.type  (character(1))
   How dummy columns are encoded. Either as 0/1 with type “numeric” or as “factor”. Default is “factor”.

force.dummies  (logical(1))
   Force dummy creation even if the respective data column does not contain any NAs. Note that (a) most learners will complain about constant columns created
makeLearner

Create learner object.

Description

For a classification learner the predict.type can be set to "prob" to predict probabilities and the maximum value selects the label. The threshold used to assign the label can later be changed using the setThreshold function.

Usage

makeLearner(
  cl,
  id = cl,
  predict.type = "response",
  predict.threshold = NULL,
  fix.factors.prediction = FALSE,
  ...
)
makeLearner

par.vals = list(),
config = list()
)

Arguments

cl (character(1))
Class of learner. By convention, all classification learners start with “classif.” all regression learners with “regr.” all survival learners start with “surv.” all clustering learners with “cluster.” and all multilabel classification learners start with “multilabel.”. A list of all integrated learners is available on the learners help page.

id (character(1))
Id string for object. Used to display object. Default is cl.

predict.type (character(1))
Classification: “response” (= labels) or “prob” (= probabilities and labels by selecting the ones with maximal probability). Regression: “response” (= mean response) or “se” (= standard errors and mean response). Survival: “response” (= some sort of orderable risk) or “prob” (= time dependent probabilities). Clustering: “response” (= cluster IDS) or “prob” (= fuzzy cluster membership probabilities), Multilabel: “response” (= logical matrix indicating the predicted class labels) or “prob” (= probabilities and corresponding logical matrix indicating class labels). Default is “response”.

predict.threshold (numeric)
Threshold to produce class labels. Has to be a named vector, where names correspond to class labels. Only for binary classification it can be a single numerical threshold for the positive class. See setThreshold for details on how it is applied. Default is NULL which means 0.5 / an equal threshold for each class.

fix.factors.prediction (logical(1))
In some cases, problems occur in underlying learners for factor features during prediction. If the new features have LESS factor levels than during training (a strict subset), the learner might produce an error like “type of predictors in new data do not match that of the training data”. In this case one can repair this problem by setting this option to TRUE. We will simply add the missing factor levels missing from the test feature (but present in training) to that feature. Default is FALSE.

... (any)
Optional named (hyper)parameters. If you want to set specific hyperparameters for a learner during model creation, these should go here. You can get a list of available hyperparameters using getParamSet(<learner>). Alternatively hyperparameters can be given using the par.vals argument but ... should be preferred!

par.vals (list)
Optional list of named (hyper)parameters. The arguments in ... take precedence over values in this list. We strongly encourage you to use ... for passing hyperparameters.
config (named list)
Named list of config option to overwrite global settings set via configureMlr for this specific learner.

Value
(Learner).

par.vals vs ... 
The former aims at specifying default hyperparameter settings from mlr which differ from the actual defaults in the underlying learner. For example, respect.unordered.factors is set to order in mlr while the default in ranger::ranger depends on the argument splitrule. getHyperPars(<learner>) can be used to query hyperparameter defaults that differ from the underlying learner. This function also shows all hyperparameters set by the user during learner creation (if these differ from the learner defaults).

regr.randomForest
For this learner we added additional uncertainty estimation functionality (predict.type = "se") for the randomForest, which is not provided by the underlying package.

Currently implemented methods are:

- If se.method = "jackknife" the standard error of a prediction is estimated by computing the jackknife-after-bootstrap, the mean-squared difference between the prediction made by only using trees which did not contain said observation and the ensemble prediction.

- If se.method = "bootstrap" the standard error of a prediction is estimated by bootstrapping the random forest, where the number of bootstrap replicates and the number of trees in the ensemble are controlled by se.boot and se.ntree respectively, and then taking the standard deviation of the bootstrap predictions. The "brute force" bootstrap is executed when ntree = se.ntree, the latter of which controls the number of trees in the individual random forests which are bootstrapped. The "noisy bootstrap" is executed when se.ntree < ntree which is less computationally expensive. A Monte-Carlo bias correction may make the latter option preferable in many cases. Defaults are se.boot = 50 and se.ntree = 100.

- If se.method = "sd", the default, the standard deviation of the predictions across trees is returned as the variance estimate. This can be computed quickly but is also a very naive estimator.

For both “jackknife” and “bootstrap”, a Monte-Carlo bias correction is applied and, in the case that this results in a negative variance estimate, the values are truncated at 0.

Note that when using the “jackknife” procedure for se estimation, using a small number of trees can lead to training data observations that are never out-of-bag. The current implementation ignores these observations, but in the original definition, the resulting se estimation would be undefined.

Please note that all of the mentioned se.method variants do not affect the computation of the posterior mean "response" value. This is always the same as from the underlying randomForest.
**regr.featureless**

A very basic baseline method which is useful for model comparisons (if you don’t beat this, you very likely have a problem). Does not consider any features of the task and only uses the target feature of the training data to make predictions. Using observation weights is currently not supported.

Methods “mean” and “median” always predict a constant value for each new observation which corresponds to the observed mean or median of the target feature in training data, respectively.

The default method is “mean” which corresponds to the ZeroR algorithm from WEKA, see [https://weka.wikispaces.com/ZeroR](https://weka.wikispaces.com/ZeroR).

**classif.featureless**

Method “majority” predicts always the majority class for each new observation. In the case of ties, one randomly sampled, constant class is predicted for all observations in the test set. This method is used as the default. It is very similar to the ZeroR classifier from WEKA (see [https://weka.wikispaces.com/ZeroR](https://weka.wikispaces.com/ZeroR)). The only difference is that ZeroR always predicts the first class of the tied class values instead of sampling them randomly.

Method “sample-prior” always samples a random class for each individual test observation according to the prior probabilities observed in the training data.

If you opt to predict probabilities, the class probabilities always correspond to the prior probabilities observed in the training data.

**See Also**

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(), setPredictThreshold(), setPredictType()

**Examples**

```r
makeLearner("classif.rpart")
makeLearner("classif.lda", predict.type = "prob")
lrn = makeLearner("classif.lda", method = "t", nu = 10)
getHyperPars(lrn)
```

**Description**

Small helper function that can save some typing when creating multiple learner objects. Calls `makeLearner` multiple times internally.

**Usage**

```r
makeLearners(cls, ids = NULL, type = NULL, ...)
```
Argument

cls (character)
Classes of learners.
ids (character)
Id strings. Must be unique. Default is cls.
type (character(1))
Shortcut to prepend type string to cls so one can set cls = "rpart". Default is NULL, i.e., this is not used.
...
(any)
Optional named (hyper)parameters. If you want to set specific hyperparameters for a learner during model creation, these should go here. You can get a list of available hyperparameters using getParamSet(<learner>). Alternatively hyperparameters can be given using the par.vals argument but ... should be preferred!

Value

(named list of Learner). Named by ids.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(), setPredictThreshold(), setPredictType()

Examples

makeLearners(c("rpart", "lda"), type = "classif", predict.type = "prob")

Description

A measure object encapsulates a function to evaluate the performance of a prediction. Information about already implemented measures can be obtained here: [measures].
A learner is trained on a training set d1, results in a model m and predicts another set d2 (which may be a different one or the training set) resulting in the prediction. The performance measure can now be defined using all of the information of the original task, the fitted model and the prediction. Object slots:

id (character(1)) See argument.
minimize (logical(1)) See argument.
properties (character)  See argument.
fun (function)  See argument.
extra.args (list)  See argument.
aggr (Aggregation)  See argument.
best (numeric(1))  See argument.
worst (numeric(1))  See argument.
name (character(1))  See argument.
ote (character(1))  See argument.

Usage
makeMeasure(
  id,
  minimize,
  properties = character(0L),
  fun,
  extra.args = list(),
  aggr = test.mean,
  best = NULL,
  worst = NULL,
  name = id,
  note = ""
)

Arguments
id (‘character(1)’)
Name of measure.
minimize (‘logical(1)’)
Should the measure be minimized? Default is ‘TRUE’.
properties (character)
Set of measure properties. Some standard property names include:
classif  Is the measure applicable for classification?
classif.multi  Is the measure applicable for multi-class classification?
multilabel  Is the measure applicable for multilabel classification?
regr  Is the measure applicable for regression?
surv  Is the measure applicable for survival?
cluster  Is the measure applicable for cluster?
costsens  Is the measure applicable for cost-sensitive learning?
req.pred  Is prediction object required in calculation? Usually the case.
req.truth  Is truth column required in calculation? Usually the case.
req.task  Is task object required in calculation? Usually not the case.
req.model  Is model object required in calculation? Usually not the case.
req.feats  Are feature values required in calculation? Usually not the case.
**makeMeasure**

- **req.prob**: Are predicted probabilities required in calculation? Usually not the case, example would be AUC.
  
  Default is `character(0)`.

- **fun**: (`function(task, model, pred, feats, extra.args)`)  
  Calculates the performance value. Usually you will only need the prediction object `pred`.

  - `task` ([Task]) The task.
  - `model` ([WrappedModel]) The fitted model.
  - `pred` ([Prediction]) Prediction object.
  - `feats` ([data.frame]) The features.
  - `extra.args` ([list]) See below.

- **extra.args**: ([list])  
  List of extra arguments which will always be passed to `fun`. Can be changed after construction via `[setMeasurePars]<3>`.

- **aggr** ([Aggregation])  
  Aggregation function, which is used to aggregate the values measured on test / training sets of the measure to a single value. Default is `[test.mean]`.

- **best**: (`numeric(1)`)  
  Best obtainable value for measure. Default is `-Inf` or `Inf`, depending on `minimize`.

- **worst**: (`numeric(1)`)  
  Worst obtainable value for measure. Default is `Inf` or `-Inf`, depending on `minimize`.

- **name** ([character])  
  Name of the measure. Default is `id`.

- **note** ([character])  
  Description and additional notes for the measure. Default is `""`.

**Value**

- **Measure**.

**See Also**

Other performance: `ConfusionMatrix, calculateConfusionMatrix()`, `calculateROCMeasures()`, `estimateRelativeOverfitting()`, `makeCostMeasure()`, `makeCustomResampledMeasure()`, `measures`, `performance()`, `setAggregation()`, `setMeasurePars()`

**Examples**

```r
f = function(task, model, pred, extra.args) {  
  sum((pred$data$response - pred$data$truth)^2)
}
makeMeasure(id = "my.sse", minimize = TRUE, properties = c("regr", "response"), fun = f)
```
**makeModelMultiplexer**

Create model multiplexer for model selection to tune over multiple possible models.

**Description**

Combines multiple base learners by dispatching on the hyperparameter “selected.learner” to a specific model class. This allows to tune not only the model class (SVM, random forest, etc) but also their hyperparameters in one go. Combine this with [tuneParams] and [makeTuneControlIrace] for a very powerful approach, see example below.

The parameter set is the union of all (unique) base learners. In order to avoid name clashes all parameter names are prefixed with the base learner id, i.e. “[learner.id].[parameter.name]”.

The predict.type of the Multiplexer is inherited from the predict.type of the base learners.

The getter [getLearnerProperties] returns the properties of the selected base learner.

**Usage**

```r
makeModelMultiplexer(base.learners)
```

**Arguments**

- `base.learners` ([list' of [Learner])
  - List of Learners with unique IDs.

**Value**

- ([ModelMultiplexer]). A [Learner] specialized as ‘ModelMultiplexer’.

**Note**

Note that logging output during tuning is somewhat shortened to make it more readable. I.e., the artificial prefix before parameter names is suppressed.

**See Also**

Other multiplexer: `makeModelMultiplexerParamSet()`

Other tune: `TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(), getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeTuneControlCMAES(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(), makeTuneControlIrace(), makeTuneControlMBO(), makeTuneControlRandom(), makeTuneWrapper(), tuneParams(), tuneThreshold()`
Examples

```r
set.seed(123)

library(BBmisc)
bls = list(
  makeLearner("classif.ksvm"),
  makeLearner("classif.randomForest")
)

lrn = makeModelMultiplexer(bls)
# simple way to construct param set for tuning
# parameter names are prefixed automatically and the 'requires'
# element is set, too, to make all parameters subordinate to 'selected.learner'
ps = makeModelMultiplexerParamSet(lrn,
  makeNumericParam("sigma", lower = -10, upper = 10, trafo = function(x) 2^(x)),
  makeIntegerParam("ntree", lower = 1L, upper = 500L)
)

print(ps)

rdesc = makeResampleDesc("CV", iters = 2L)
# to save some time we use random search. but you probably want something like this:
# ctrl = makeTuneControlIrace(maxExperiments = 500L)
ctrl = makeTuneControlRandom(maxit = 10L)

res = tuneParams(lrn, iris.task, rdesc, par.set = ps, control = ctrl)
print(res)

df = as.data.frame(res$opt.path)
print(head(df[, -ncol(df)]))

# more unique and reliable way to construct the param set
ps = makeModelMultiplexerParamSet(lrn,
  classif.ksvm = makeParamSet(
    makeNumericParam("sigma", lower = -10, upper = 10, trafo = function(x) 2^(x))
  ),
  classif.randomForest = makeParamSet(
    makeIntegerParam("ntree", lower = 1L, upper = 500L)
  )
)

# this is how you would construct the param set manually, works too
ps = makeParamSet(
  makeDiscreteParam("selected.learner", values = extractSubList(bls, "id")),
  makeNumericParam("classif.ksvm.sigma", lower = -10, upper = 10, trafo = function(x) 2^(x),
    requires = quote(selected.learner == "classif.ksvm")),
  makeIntegerParam("classif.randomForest.ntree", lower = 1L, upper = 500L,
    requires = quote(selected.learner == "classif.randomForest"))
)

# all three ps-objects are exactly the same internally.
```
makeModelMultiplexerParamSet

Create a parameter set for model multiplexer tuning.

Description

Handy way to create the param set with less typing.
The following is done automatically:

• The selected.learner param is created
• Parameter names are prefixed.
• The requires field of each param is set. This makes all parameters subordinate to selected.learner

Usage

makeModelMultiplexerParamSet(multiplexer, ..., .check = TRUE)

Arguments

multiplexer  
(ModelMultiplexer)
The multiplexer learner.

...  
(ParamHelpers::ParamSet | ParamHelpers::Param)
(a) First option: Named param sets. Names must correspond to base learners. You only need to enter the parameters you want to tune without reference to the selected.learner field in any way.
(b) Second option. Just the params you would enter in the param sets. Even shorter to create. Only works when it can be uniquely identified to which learner each of your passed parameters belongs.

.check  
(logical)
Check that for each param in ... one param in found in the base learners. Default is TRUE

Value

ParamSet.

See Also

Other multiplexer: makeModelMultiplexer()
Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(), getResamplingIndices(), getTuneResult(), makeModelMultiplexer(), makeTuneControlCMAES(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(), makeTuneControlIrace(), makeTuneControlMBO(), makeTuneControlRandom(), makeWrapper(), tuneParams(), tuneThreshold()

Examples

# See makeModelMultiplexer
makeMulticlassWrapper

Fuse learner with multiclass method.

Description

Fuses a base learner with a multi-class method. Creates a learner object, which can be used like any other learner object. This way learners which can only handle binary classification will be able to handle multi-class problems, too.

We use a multiclass-to-binary reduction principle, where multiple binary problems are created from the multiclass task. How these binary problems are generated is defined by an error-correcting-output-code (ECOC) code book. This also allows the simple and well-known one-vs-one and one-vs-rest approaches. Decoding is currently done via Hamming decoding, see e.g. here http://jmlr.org/papers/volume11/escalera10a/escalera10a.pdf.

Currently, the approach always operates on the discrete predicted labels of the binary base models (instead of their probabilities) and the created wrapper cannot predict posterior probabilities.

Usage

makeMulticlassWrapper(learner, mcw.method = "onevsrest")

Arguments

learner (Learner | character(1))
   The learner. If you pass a string the learner will be created via makeLearner.

cmw.method (character(1) | function)
   "onevsone" or "onevsrest". You can also pass a function, with signature function(task) and which returns a ECOC codematrix with entries +1,-1,0. Columns define new binary problems, rows correspond to classes (rows must be named). 0 means class is not included in binary problem. Default is "onevsrest".

Value

Learner.

See Also

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()
makeMultilabelBinaryRelevanceWrapper

Description

Every learner which is implemented in mlr and which supports binary classification can be converted to a wrapped binary relevance multilabel learner. The multilabel classification problem is converted into simple binary classifications for each label/target on which the binary learner is applied.

Models can easily be accessed via `getLearnerModel`.

Note that it does not make sense to set a threshold in the used base learner when you predict probabilities. On the other hand, it can make a lot of sense, to call `setThreshold` on the `MultilabelBinaryRelevanceWrapper` for each label individually; Or to tune these thresholds with `tuneThreshold`; especially when you face very unbalanced class distributions for each binary label.

Usage

```r
makeMultilabelBinaryRelevanceWrapper(learner)
```

Arguments

- `learner` *(Learner | character(1))*
  The learner. If you pass a string the learner will be created via `makeLearner`.

Value

Learner.

References


See Also

Other wrapper: `makeBaggingWrapper()`, `makeClassificationViaRegressionWrapper()`, `makeConstantClassWrapper()`, `makeCostSensClassifWrapper()`, `makeCostSensRegrWrapper()`, `makeDownsampleWrapper()`, `makeDummyFeaturesWrapper()`, `makeExtractFDAFeatsWrapper()`, `makeFeatSelWrapper()`, `makeFilterWrapper()`, `makeImputeWrapper()`, `makeMulticlassWrapper()`, `makeMultilabelClassifierChainsWrapper()`, `makeMultilabelDBRWrapper()`, `makeMultilabelNestedStackingWrapper()`, `makeMultilabelStackingWrapper()`, `makeOverBaggingWrapper()`, `makePreprocWrapperCaret()`, `makePreprocWrapper()`, `makeRemoveConstantFeaturesWrapper()`, `makeSMOTETwrapper()`, `makeTuneWrapper()`, `makeUndersampleWrapper()`, `makeWeightedClassesWrapper()`

Other multilabel: `getMultilabelBinaryPerformances()`, `makeMultilabelClassifierChainsWrapper()`, `makeMultilabelDBRWrapper()`, `makeMultilabelNestedStackingWrapper()`, `makeMultilabelStackingWrapper()`
Examples

d = getTaskData(yeast.task)
# drop some labels so example runs faster
d = d[seq(1, nrow(d), by = 20), c(1:2, 15:17)]
task = makeMultilabelTask(data = d, target = c("label1", "label2"))
lrn = makeLearner("classif.rpart")
lrn = makeMultilabelBinaryRelevanceWrapper(lrn)
lrn = setPredictType(lrn, "prob")
# train, predict and evaluate
mod = train(lrn, task)
pred = predict(mod, task)
performance(pred, measure = list(multilabel.hamloss, multilabel.subset01, multilabel.f1))
# the next call basically has the same structure for any multilabel meta wrapper
getMultilabelBinaryPerformances(pred, measures = list(mmce, auc))
# above works also with predictions from resample!

makeMultilabelClassifierChainsWrapper

Use classifier chains method (CC) to create a multilabel learner.

Description

Every learner which is implemented in mlr and which supports binary classification can be con-
verted to a wrapped classifier chains multilabel learner. CC trains a binary classifier for each label
following a given order. In training phase, the feature space of each classifier is extended with true
label information of all previous labels in the chain. During the prediction phase, when true labels
are not available, they are replaced by predicted labels.

Models can easily be accessed via getLearnerModel.

Usage

makeMultilabelClassifierChainsWrapper(learner, order = NULL)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

order (character)
Specifies the chain order using the names of the target labels. E.g. for m target
labels, this must be a character vector of length m that contains a permutation
of the target label names. Default is NULL which uses a random ordering of the
target label names.

Value

Learner.
makeMultilabelDBRWrapper

References

Montanes, E. et al. (2013) Dependent binary relevance models for multi-label classification Artificial Intelligence Center, University of Oviedo at Gijon, Spain.

See Also

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()

Other multilabel: getMultilabelBinaryPerformances(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper()

Examples

d = getTaskData(yeast.task)
# drop some labels so example runs faster
d = d[seq(1, nrow(d), by = 20), c(1:2, 15:17)]
task = makeMultilabelTask(data = d, target = c("label1", "label2"))
lrn = makeLearner("classif.rpart")
lrn = makeMultilabelBinaryRelevanceWrapper(lrn)
clr = setPredictType(lrn, "prob")
# train, predict and evaluate
mod = train(lrn, task)
pred = predict(mod, task)
performance(pred, measure = list(multilabel.hamloss, multilabel.subset01, multilabel.f1))
# the next call basically has the same structure for any multilabel meta wrapper
getMultilabelBinaryPerformances(pred, measures = list(mmce, auc))
# above works also with predictions from resample!

makeMultilabelDBRWrapper

Use dependent binary relevance method (DBR) to create a multilabel learner.

Description

Every learner which is implemented in mlr and which supports binary classification can be converted to a wrapped DBR multilabel learner. The multilabel classification problem is converted into simple binary classifications for each label/target on which the binary learner is applied. For each target, actual information of all binary labels (except the target variable) is used as additional features. During prediction these labels need are obtained by the binary relevance method using the same binary learner.

Models can easily be accessed via getLearnerModel.
makeMultilabelDBRWrapper

Usage

makeMultilabelDBRWrapper(learner)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

Value

Learner.

References

Montanes, E. et al. (2013) *Dependent binary relevance models for multi-label classification* Artificial Intelligence Center, University of Oviedo at Gijon, Spain.

See Also

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEBinaryWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()

Other multilabel: getMultilabelBinaryPerformances(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper()

Examples

d = getTaskData(yeast.task)
# drop some labels so example runs faster
d = d[seq(1, nrow(d), by = 20), c(1:2, 15:17)]
task = makeMultilabelTask(data = d, target = c("label1", "label2"))
lrn = makeLearner("classif.rpart")
lrn = makeMultilabelBinaryRelevanceWrapper(lrn)
lrn = setPredictType(lrn, "prob")
# train, predict and evaluate
mod = train(lrn, task)
pred = predict(mod, task)
performance(pred, measure = list(multilabel.hamloss, multilabel.subset01, multilabel.f1))
# the next call basically has the same structure for any multilabel meta wrapper
getMultilabelBinaryPerformances(pred, measures = list(mmce, auc))
# above works also with predictions from resample!
makeMultilabelNestedStackingWrapper

Use nested stacking method to create a multilabel learner.

Description

Every learner which is implemented in mlr and which supports binary classification can be converted to a wrapped nested stacking multilabel learner. Nested stacking trains a binary classifier for each label following a given order. In training phase, the feature space of each classifier is extended with predicted label information (by cross validation) of all previous labels in the chain. During the prediction phase, predicted labels are obtained by the classifiers, which have been learned on all training data.

Models can easily be accessed via `getLearnerModel`.

Usage

```r
makeMultilabelNestedStackingWrapper(learner, order = NULL, cv.folds = 2)
```

Arguments

- **learner** *(Learner | character(1))*
  The learner. If you pass a string the learner will be created via `makeLearner`.

- **order** *(character)*
  Specifies the chain order using the names of the target labels. E.g. for `m` target labels, this must be a character vector of length `m` that contains a permutation of the target label names. Default is `NULL` which uses a random ordering of the target label names.

- **cv.folds** *(integer(1))*
  The number of folds for the inner cross validation method to predict labels for the augmented feature space. Default is 2.

Value

`Learner`.

References

Montanes, E. et al. (2013), Dependent binary relevance models for multi-label classification Artificial Intelligence Center, University of Oviedo at Gijon, Spain.

See Also

makeMultilabelStackingWrapper

Use stacking method (stacked generalization) to create a multilabel learner.

Description

Every learner which is implemented in mlr and which supports binary classification can be converted to a wrapped stacking multilabel learner. Stacking trains a binary classifier for each label using predicted label information of all labels (including the target label) as additional features (by cross validation). During prediction these labels need are obtained by the binary relevance method using the same binary learner.

Models can easily be accessed via getLearnerModel.

Usage

makeMultilabelStackingWrapper(learner, cv.folds = 2)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

cv.folds (integer(1))
The number of folds for the inner cross validation method to predict labels for the augmented feature space. Default is 2.
**makeMultilabelTask**

Create a multilabel task.

**Value**

Learner.

**References**

Montanes, E. et al. (2013) Dependent binary relevance models for multi-label classification Artificial Intelligence Center, University of Oviedo at Gijon, Spain.

**See Also**

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()

Other multilabel: getMultilabelBinaryPerformances(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper()

**Examples**

d = getTaskData(yeast.task)
# drop some labels so example runs faster
d = d[seq(1, nrow(d), by = 20), c(1:2, 15:17)]
task = makeMultilabelTask(data = d, target = c("label1", "label2"))
lnr = makeLearner("classif.rpart")
lnr = makeMultilabelBinaryRelevanceWrapper(lnr)
lnr = setPredictType(lnr, "prob")
# train, predict and evaluate
mod = train(lnr, task)
pred = predict(mod, task)
performance(pred, measure = list(multilabel.hamloss, multilabel.subset01, multilabel.f1))
# the next call basically has the same structure for any multilabel meta wrapper
getMultilabelBinaryPerformances(pred, measures = list(mmce, auc))
# above works also with predictions from resample!
Usage

makeMultilabelTask(
  id = deparse(substitute(data)),
  data,
  target,
  weights = NULL,
  blocking = NULL,
  coordinates = NULL,
  fixup.data = "warn",
  check.data = TRUE
)

Arguments

id (character(1))
Id string for object. Default is the name of the R variable passed to data.

data (data.frame)
A data frame containing the features and target variable(s).

target (character(1) | character(2) | character(n.classes))
Name(s) of the target variable(s). For survival analysis these are the names of
the survival time and event columns, so it has length 2. For multilabel classification it contains the names of the logical columns that encode whether a label is present or not and its length corresponds to the number of classes.

weights (numeric)
Optional, non-negative case weight vector to be used during fitting. Cannot be set for cost-sensitive learning. Default is NULL which means no (= equal) weights.

blocking (factor)
An optional factor of the same length as the number of observations. Observations with the same blocking level “belong together”. Specifically, they are either put all in the training or the test set during a resampling iteration. Default is NULL which means no blocking.

coordinates (data.frame)
Coordinates of a spatial data set that will be used for spatial partitioning of the data in a spatial cross-validation resampling setting. Coordinates have to be numeric values. Provided data.frame needs to have the same number of rows as data and consist of at least two dimensions.

fixup.data (character(1))
Should some basic cleaning up of data be performed? Currently this means removing empty factor levels for the columns. Possible choices are: “no” = Don’t do it. “warn” = Do it but warn about it. “quiet” = Do it but keep silent. Default is “warn”.

check.data (logical(1))
Should sanity of data be checked initially at task creation? You should have good reasons to turn this off (one might be speed). Default is TRUE.
makeOverBaggingWrapper

Details

For multilabel classification we assume that the presence of labels is encoded via logical columns in data. The name of the column specifies the name of the label. target is then a char vector that points to these columns.

Note

For multilabel classification we assume that the presence of labels is encoded via logical columns in data. The name of the column specifies the name of the label. target is then a char vector that points to these columns.

See Also

Task ClassifTask ClusterTask CostSensTask RegrTask SurvTask

makeOverBaggingWrapper

Fuse learner with the bagging technique and oversampling for imbalancy correction.

Description

Fuses a classification learner for binary classification with an over-bagging method for imbalancy correction when we have strongly unequal class sizes. Creates a learner object, which can be used like any other learner object. Models can easily be accessed via getLearnerModel.

OverBagging is implemented as follows: For each iteration a random data subset is sampled. Class examples are oversampled with replacement with a given rate. Members of the other class are either simply copied into each bag, or bootstrapped with replacement until we have as many majority class examples as in the original training data. Features are currently not changed or sampled.

Prediction works as follows: For classification we do majority voting to create a discrete label and probabilities are predicted by considering the proportions of all predicted labels.

Usage

makeOverBaggingWrapper(
  learner,
  obw.iters = 10L,
  obw.rate = 1,
  obw.maxcl = "boot",
  obw.cl = NULL
)
makePreprocWrapper

**Arguments**

- **learner**
  - Type: (Learner | character(1))
  - Description: The learner. If you pass a string the learner will be created via `makeLearner`.

- **obw.iters**
  - Type: integer(1)
  - Description: Number of fitted models in bagging. Default is 10.

- **obw.rate**
  - Type: numeric(1)
  - Description: Factor to upsample a class in each bag. Must be between 1 and Inf, where 1 means no oversampling and 2 would mean doubling the class size. Default is 1.

- **obw.maxcl**
  - Type: character(1)
  - Description: How should other class (usually larger class) be handled? “all” means every instance of the class gets in each bag, “boot” means the class instances are bootstrapped in each iteration. Default is “boot”.

- **obw.cl**
  - Type: character(1)
  - Description: Which class should be over- or undersampled. If NULL, `makeOverBaggingWrapper` will take the smaller class.

**Value**

- Type: Learner.

**See Also**

Other imbalancy: `makeUndersampleWrapper()`, `oversample()`, `smote()`

Other wrapper: `makeBaggingWrapper()`, `makeClassificationViaRegressionWrapper()`, `makeConstantClassWrapper()`, `makeCostSensClassifWrapper()`, `makeCostSensRegrWrapper()`, `makeDownsampleWrapper()`, `makeDummyFeaturesWrapper()`, `makeExtractFDAFeatsWrapper()`, `makeFeatSelWrapper()`, `makeFilterWrapper()`, `makeImputeWrapper()`, `makeMulticlassWrapper()`, `makeMultilabelBinaryRelevanceWrapper()`, `makeMultilabelClassifierChainsWrapper()`, `makeMultilabelDBRWrapper()`, `makeMultilabelNestedStackingWrapper()`, `makeMultilabelStackingWrapper()`, `makePreprocWrapperCaret()`

**Description**

Fuses a base learner with a preprocessing method. Creates a learner object, which can be used like any other learner object, but which internally preprocesses the data as requested. If the train or predict function is called on data / a task, the preprocessing is always performed automatically.
makePreprocWrapper

Usage

makePreprocWrapper(
  learner,
  train,
  predict,
  par.set = makeParamSet(),
  par.vals = list()
)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

train (function(data, target, args))
Function to preprocess the data before training. target is a string and denotes the target variable in data. args is a list of further arguments and parameters to influence the preprocessing. Must return a list(data,control), where data is the preprocessed data and control stores all information necessary to do the preprocessing before predictions.

predict (function(data, target, args, control))
Function to preprocess the data before prediction. target is a string and denotes the target variable in data. args are the args that were passed to train. control is the object you returned in train. Must return the processed data.

par.set (ParamHelpers::ParamSet)
Parameter set of ParamHelpers::LearnerParam objects to describe the parameters in args. Default is empty set.

par.vals (list)
Named list of default values for params in args respectively par.set. Default is empty list.

Value

(Learner).

See Also

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper())
**makePreprocWrapperCaret**

_Fuse learner with preprocessing._

**Description**

Fuses a learner with preprocessing methods provided by `caret::preProcess`. Before training the preprocessing will be performed and the preprocessing model will be stored. Before prediction the preprocessing model will transform the test data according to the trained model.

After being wrapped the learner will support missing values although this will only be the case if `ppc.knnImpute`, `ppc.bagImpute` or `ppc.medianImpute` is set to `TRUE`.

**Usage**

```r
makePreprocWrapperCaret(learner, ...)
```

**Arguments**

- `learner` *(Learner | character(1))*
  The learner. If you pass a string the learner will be created via `makeLearner`.

- `...` *(any)*
  See `caret::preProcess` for parameters not listed above. If you use them you might want to define them in the `add.par.set` so that they can be tuned.

**Value**

`Learner`.

**See Also**

Other wrapper: `makeBaggingWrapper()`, `makeClassificationViaRegressionWrapper()`, `makeConstantClassWrapper()`, `makeCostSensClassifWrapper()`, `makeCostSensRegrWrapper()`, `makeDownsampleWrapper()`, `makeDummyFeaturesWrapper()`, `makeExtractFDAFeatsWrapper()`, `makeFeatSelWrapper()`, `makeFilterWrapper()`, `makeImputeWrapper()`, `makeMulticlassWrapper()`, `makeMultilabelBinaryRelevanceWrapper()`, `makeMultilabelClassifierChainsWrapper()`, `makeMultilabelDBRWrapper()`, `makeMultilabelNestedStackingWrapper()`, `makeMultilabelStackingWrapper()`, `makeOverBaggingWrapper()`, `makePreprocWrapper()`, `makeRemoveConstantFeaturesWrapper()`, `makeSMOTEWrapper()`, `makeTuneWrapper()`, `makeUndersampleWrapper()`, `makeWeightedClassesWrapper()`
makeRegrTask

Create a regression task.

Description

Create a regression task.

Usage

makeRegrTask(
  id = deparse(substitute(data)),
  data,
  target,
  weights = NULL,
  blocking = NULL,
  coordinates = NULL,
  fixup.data = "warn",
  check.data = TRUE
)

Arguments

id (character(1))
Id string for object. Default is the name of the R variable passed to data.

data (data.frame)
A data frame containing the features and target variable(s).

target (character(1) | character(2) | character(n.classes))
Name(s) of the target variable(s). For survival analysis these are the names of the survival time and event columns, so it has length 2. For multilabel classification it contains the names of the logical columns that encode whether a label is present or not and its length corresponds to the number of classes.

weights (numeric)
Optional, non-negative case weight vector to be used during fitting. Cannot be set for cost-sensitive learning. Default is NULL which means no (= equal) weights.

blocking (factor)
An optional factor of the same length as the number of observations. Observations with the same blocking level “belong together”. Specifically, they are either put all in the training or the test set during a resampling iteration. Default is NULL which means no blocking.

coordinates (data.frame)
Coordinates of a spatial data set that will be used for spatial partitioning of the data in a spatial cross-validation resampling setting. Coordinates have to be numeric values. Provided data.frame needs to have the same number of rows as data and consist of at least two dimensions.
fixup.data (character(1))
Should some basic cleaning up of data be performed? Currently this means removing empty factor levels for the columns. Possible choices are: “no” = Don’t do it. “warn” = Do it but warn about it. “quiet” = Do it but keep silent. Default is “warn”.

check.data (logical(1))
Should sanity of data be checked initially at task creation? You should have good reasons to turn this off (one might be speed). Default is TRUE.

See Also
Task ClassifTask CostSensTask ClusterTask MultilabelTask SurvTask

makeRemoveConstantFeaturesWrapper
Fuses learner with removal of constant features preprocessing.

Description
Fuses a base learner with the preprocessing implemented in removeConstantFeatures.

Usage
makeRemoveConstantFeaturesWrapper(
  learner,
  perc = 0,
  dont.rm = character(0L),
  na.ignore = FALSE,
  tol = .Machine$double.eps^0.5
)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

perc (numeric(1))
The percentage of a feature values in [0, 1) that must differ from the mode value. Default is 0, which means only constant features with exactly one observed level are removed.

dont.rm (character)
Names of the columns which must not be deleted. Default is no columns.

na.ignore (logical(1))
Should NAs be ignored in the percentage calculation? (Or should they be treated as a single, extra level in the percentage calculation?) Note that if the feature has only missing values, it is always removed. Default is FALSE.

tol (numeric(1))
Numerical tolerance to treat two numbers as equal. Variables stored as double will get rounded accordingly before computing the mode. Default is \( \sqrt{\text{Machine}\text{double}\text{eps}} \).
makeResampleDesc

Create a description object for a resampling strategy.

Description

A description of a resampling algorithm contains all necessary information to create a ResampleInstance, when given the size of the data set.

Usage

makeResampleDesc(
  method, 
  predict = "test", 
  ...,
  stratify = FALSE, 
  stratify.cols = NULL, 
  fixed = FALSE, 
  blocking.cv = FALSE
)

Arguments

method (character(1))

predict (character(1))
  What to predict during resampling: “train”, “test” or “both” sets. Default is “test”.

... (any)
  Further parameters for strategies.
makenResampleDesc

**iters** (integer(1)) Number of iterations, for “CV”, “Subsample” and “Bootstrap”.

**split** (numeric(1)) Proportion of training cases for “Holdout” and “Subsample” between 0 and 1. Default is 2 / 3.

**reps** (integer(1)) Repeats for “RepCV”. Here iters = folds * reps. Default is 10.

**folds** (integer(1)) Folds in the repeated CV for RepCV. Here iters = folds * reps. Default is 10.

**horizon** (numeric(1)) Number of observations in the forecast test set for “GrowingWindowCV” and “FixedWindowCV”. When horizon > 1 this will be treated as the number of observations to forecast, else it will be a fraction of the initial window. IE, for 100 observations, initial window of .5, and horizon of .2, the test set will have 10 observations. Default is 1.

**initial.window** (numeric(1)) Fraction of observations to start with in the training set for “GrowingWindowCV” and “FixedWindowCV”. When initial.window > 1 this will be treated as the number of observations in the initial window, else it will be treated as the fraction of observations to have in the initial window. Default is 0.5.

**skip** (numeric(1)) How many resamples to skip to thin the total amount for “GrowingWindowCV” and “FixedWindowCV”. This is passed through as the “by” argument in seq(). When skip > 1 this will be treated as the increment of the sequence of resampling indices, else it will be a fraction of the total training indices. IE for 100 training sets and a value of .2, the increment of the resampling indices will be 20. Default is “horizon” which gives mutually exclusive chunks of test indices.

**stratify** (logical(1)) Should stratification be done for the target variable? For classification tasks, this means that the resampling strategy is applied to all classes individually and the resulting index sets are joined to make sure that the proportion of observations in each training set is as in the original data set. Useful for imbalanced class sizes. For survival tasks stratification is done on the events, resulting in training sets with comparable censoring rates.

**stratify.cols** (character) Stratify on specific columns referenced by name. All columns have to be factor or integer. Note that you have to ensure yourself that stratification is possible, i.e. that each strata contains enough observations. This argument and stratify are mutually exclusive.

**fixed** (logical(1)) Whether indices supplied via argument ‘blocking’ in the task should be used as fully pre-defined indices. Default is FALSE which means they will be used following the ‘blocking’ approach. fixed only works with ResampleDesc CV and the supplied indices must match the number of observations. When fixed = TRUE, the iters argument will be ignored and is interally set to the number of supplied factor levels in blocking.

**blocking.cv** (logical(1)) Should ‘blocking’ be used in CV? Default to FALSE. This is different to fixed
makeResampleDesc

= TRUE and cannot be combined. Please check the mlr online tutorial for more details.

Details

Some notes on some special strategies:

**Repeated cross-validation** Use “RepCV”. Then you have to set the aggregation function for your preferred performance measure to “testgroup.mean” via `setAggregation`.

**B632 bootstrap** Use “Bootstrap” for bootstrap and set `predict` to “both”. Then you have to set the aggregation function for your preferred performance measure to “b632” via `setAggregation`.

**B632+ bootstrap** Use “Bootstrap” for bootstrap and set `predict` to “both”. Then you have to set the aggregation function for your preferred performance measure to “b632plus” via `setAggregation`.

**Fixed Holdout set** Use `makeFixedHoldoutInstance`.

Object slots:

- **id** (character(1)) Name of resampling strategy.
- **iters** (integer(1)) Number of iterations. Note that this is always the complete number of generated train/test sets, so for a 10-times repeated 5fold cross-validation it would be 50.
- **predict** (character(1)) See argument.
- **stratify** (logical(1)) See argument.

All parameters passed in ... under the respective argument name See arguments.

Value

(ResampleDesc).

Standard ResampleDesc objects

For common resampling strategies you can save some typing by using the following description objects:

- **hout** holdout a.k.a. test sample estimation (two-thirds training set, one-third testing set)
- **cv2** 2-fold cross-validation
- **cv3** 3-fold cross-validation
- **cv5** 5-fold cross-validation
- **cv10** 10-fold cross-validation

See Also

Other resample: `ResamplePrediction`, `ResampleResult`, `addRRMeasure()`, `getRRPredictionList()`, `getRRPredictions()`, `getRRTaskDescription()`, `getRRTaskDesc()`, `makeResampleInstance()`, `resample()`


## Examples

### Bootstraping

```r
makeResampleDesc("Bootstrap", iters = 10)
makeResampleDesc("Bootstrap", iters = 10, predict = "both")
```

### Subsampling

```r
makeResampleDesc("Subsample", iters = 10, split = 3 / 4)
makeResampleDesc("Subsample", iters = 10)
```

### Holdout a.k.a. test sample estimation

```r
makeResampleDesc("Holdout")
```

## `makeResampleInstance`

**Instantiates a resampling strategy object.**

### Description

This class encapsulates training and test sets generated from the data set for a number of iterations. It mainly stores a set of integer vectors indicating the training and test examples for each iteration.

### Usage

```r
makeResampleInstance(desc, task, size, ...)
```

### Arguments

- **desc** *(ResampleDesc | character(1))*
  
  Resampling description object or name of resampling strategy. In the latter case `makeResampleDesc` will be called internally on the string.

- **task** *(Task)*
  
  Data of task to resample from. Prefer to pass this instead of `size`.

- **size** *(integer)*
  
  Size of the data set to resample. Can be used instead of `task`.

- **...** *(any)*
  
  Passed down to `makeResampleDesc` in case you passed a string in `desc`. Otherwise ignored.

### Details

Object slots:

- **desc** *(ResampleDesc)* See argument.
- **size** *(integer(1))* See argument.
- **train.inds** *(list of integer)* List of of training indices for all iterations.
- **test.inds** *(list of integer)* List of of test indices for all iterations.
- **group** *(factor)* Optional grouping of resampling iterations. This encodes whether specific iterations ‘belong together’ (e.g. repeated CV), and it can later be used to aggregate performance values accordingly. Default is ‘factor’.
Value

(ResampleInstance).

See Also

Other resample: ResamplePrediction, ResampleResult, addRRMeasure(), getRRPredictionList(), getRRPredictions(), getRRTaskDescription(), getRRTaskDesc(), makeResampleDesc(), resample().

Examples

rdesc = makeResampleDesc("Bootstrap", iters = 10)
in = makeResampleInstance(rdesc, task = iris.task)

rdesc = makeResampleDesc("CV", iters = 50)
in = makeResampleInstance(rdesc, size = nrow(iris))

in = makeResampleInstance("CV", iters = 10, task = iris.task)

makeRLearner.classif.fdausc.glm

Classification of functional data by Generalized Linear Models.

Description

Learner for classification using Generalized Linear Models.

Usage

## S3 method for class 'classif.fdausc.glm'
makeRLearner()

makeRLearner.classif.fdausc.kernel

Learner for kernel classification for functional data.

Description

Learner for kernel Classification.

Usage

## S3 method for class 'classif.fdausc.kernel'
makeRLearner()
makeR\texttt{Learner.classif.fdausc.np}

\textit{Learner for nonparametric classification for functional data.}

\section*{Description}

Learner for Nonparametric Supervised Classification.

\section*{Usage}

\begin{verbatim}
## S3 method for class 'classif.fdausc.np'
makeR\texttt{Learner}()
\end{verbatim}

\section*{make\texttt{SMOTEWrapper}}

\textit{Fuse learner with SMOTE oversampling for imbalancy correction in binary classification.}

\section*{Description}

Creates a learner object, which can be used like any other learner object. Internally uses \texttt{smote} before every model fit.

Note that observation weights do not influence the sampling and are simply passed down to the next learner.

\section*{Usage}

\begin{verbatim}
make\texttt{SMOTEWrapper}(  
  learner,  
  sw.rate = 1,  
  sw.nn = 5L,  
  sw.standardize = TRUE,  
  sw.alt.logic = FALSE  
)
\end{verbatim}

\section*{Arguments}

\begin{itemize}
  \item \texttt{learner} \hspace{1cm} \texttt{(Learner|character(1))}
    The learner. If you pass a string the learner will be created via \texttt{makeLearner}.
  \item \texttt{sw.rate} \hspace{1cm} \texttt{(numeric(1))}
    Factor to oversample the smaller class. Must be between 1 and Inf, where 1 means no oversampling and 2 would mean doubling the class size. Default is 1.
  \item \texttt{sw.nn} \hspace{1cm} \texttt{(integer(1))}
    Number of nearest neighbors to consider. Default is 5.
\end{itemize}
makeStackedLearner

Create a stacked learner object.

Description

A stacked learner uses predictions of several base learners and fits a super learner using these predictions as features in order to predict the outcome. The following stacking methods are available:

- **'average'** Averaging of base learner predictions without weights.
- **'stack.nocv'** Fits the super learner, where in-sample predictions of the base learners are used.
- **'stack.cv'** Fits the super learner, where the base learner predictions are computed by crossvalidated predictions (the resampling strategy can be set via the ‘resampling’ argument).
- **'hill.climb'** Select a subset of base learner predictions by hill climbing algorithm.
- **'compress'** Train a neural network to compress the model from a collection of base learners.

sw.standardize (logical(1))

Standardize input variables before calculating the nearest neighbors for data sets with numeric input variables only. For mixed variables (numeric and factor) the gower distance is used and variables are standardized anyway. Default is TRUE.

sw.alt.logic (logical(1))

Use an alternative logic for selection of minority class observations. Instead of sampling a minority class element AND one of its nearest neighbors, each minority class element is taken multiple times (depending on rate) for the interpolation and only the corresponding nearest neighbor is sampled. Default is FALSE.

Value

Learner.

See Also

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMulticlassBinaryRelevanceWrapper(), makeMulticlassClassifierChainsWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeTuneWrapper(), makeUndersampleWrapper(), makeWeightedClassesWrapper()
Usage

makeStackedLearner(
  base.learners,
  super.learner = NULL,
  predict.type = NULL,
  method = "stack.nocv",
  use.feat = FALSE,
  resampling = NULL,
  parset = list()
)

Arguments

base.learners ([list of) [Learner])
A list of learners created with 'makeLearner'.

super.learner [Learner | character(1)]
The super learner that makes the final prediction based on the base learners. If
you pass a string, the super learner will be created via 'makeLearner'. Not used
for 'method = 'average''. Default is 'NULL'.

predict.type ('character(1)')
Sets the type of the final prediction for 'method = 'average''. For other methods,
the predict type should be set within 'super.learner'. If the type of the base
learner prediction, which is set up within 'base.learners', is

"prob" then 'predict.type = 'prob' will use the average of all bease learner
predictions and 'predict.type = 'response' will use the class with highest
probability as final prediction.

"response" then, for classification tasks with 'predict.type = 'prob', the fi-
nal prediction will be the relative frequency based on the predicted base
learner classes and classification tasks with 'predict.type = 'response' will
use majority vote of the base learner predictions to determine the final pre-
diction. For regression tasks, the final prediction will be the average of the
base learner predictions.

method ('character(1)')
"average" for averaging the predictions of the base learners, "stack.nocv" for
building a super learner using the predictions of the base learners, "stack.cv" for
building a super learner using crossvalidated predictions of the base learners.
"hill.climb" for averaging the predictions of the base learners, with the weights
learned from hill climbing algorithm and "compress" for compressing the model
to mimic the predictions of a collection of base learners while speeding up the
predictions and reducing the size of the model. Default is "stack.nocv",

use.feat ('logical(1)')
Whether the original features should also be passed to the super learner. Not
used for 'method = 'average'. Default is 'FALSE'.

resampling ([ResampleDesc])
Resampling strategy for 'method = 'stack.cv'. Currently only CV is allowed
for resampling. The default 'NULL' uses 5-fold CV.
the parameters for ‘hill.climb’ method, including

‘replace’ Whether a base learner can be selected more than once.

‘init’ Number of best models being included before the selection algorithm.

‘bagprob’ The proportion of models being considered in one round of selection.

‘bagtime’ The number of rounds of the bagging selection.

‘metric’ The result evaluation metric function taking two parameters ‘pred’ and ‘true’, the smaller the score the better.

the parameters for ‘compress’ method, including

k the size multiplier of the generated data

prob the probability to exchange values

s the standard deviation of each numerical feature

Examples

# Classification
data(iris)
  tsk = makeClassifTask(data = iris, target = "Species")
  base = c("classif.rpart", "classif.lda", "classif.svm")
  lrns = lapply(base, makeLearner)
  lrns = lapply(lrns, setPredictType, "prob")
  m = makeStackedLearner(base.learners = lrns,
                         predict.type = "prob", method = "hill.climb")
  tmp = train(m, tsk)
  res = predict(tmp, tsk)

# Regression
data(BostonHousing, package = "mlbench")
  tsk = makeRegrTask(data = BostonHousing, target = "medv")
  base = c("regr.rpart", "regr.svm")
  lrns = lapply(base, makeLearner)
  m = makeStackedLearner(base.learners = lrns,
                         predict.type = "response", method = "compress")
  tmp = train(m, tsk)
  res = predict(tmp, tsk)
Usage

makeSurvTask(
  id = deparse(substitute(data)),
  data,
  target,
  weights = NULL,
  blocking = NULL,
  coordinates = NULL,
  fixup.data = "warn",
  check.data = TRUE
)

Arguments

id (character(1))
Id string for object. Default is the name of the R variable passed to data.

data (data.frame)
A data frame containing the features and target variable(s).

target (character(1) | character(2) | character(n.classes))
Name(s) of the target variable(s). For survival analysis these are the names of
the survival time and event columns, so it has length 2. For multilabel classifi-
cation it contains the names of the logical columns that encode whether a label
is present or not and its length corresponds to the number of classes.

weights (numeric)
Optional, non-negative case weight vector to be used during fitting. Cannot
be set for cost-sensitive learning. Default is NULL which means no (= equal)
weights.

blocking (factor)
An optional factor of the same length as the number of observations. Observ-
ations with the same blocking level “belong together”. Specifically, they are
either put all in the training or the test set during a resampling iteration. Default
is NULL which means no blocking.

coordinates (data.frame)
Coordinates of a spatial data set that will be used for spatial partitioning of the
data in a spatial cross-validation resampling setting. Coordinates have to be
numeric values. Provided data.frame needs to have the same number of rows as
data and consist of at least two dimensions.

fixup.data (character(1))
Should some basic cleaning up of data be performed? Currently this means
removing empty factor levels for the columns. Possible choices are: “no” =
Don’t do it. “warn” = Do it but warn about it. “quiet” = Do it but keep silent.
Default is “warn”.

check.data (logical(1))
Should sanity of data be checked initially at task creation? You should have
good reasons to turn this off (one might be speed). Default is TRUE.
makeTuneControlCMAES

Create control object for hyperparameter tuning with CMAES.

Description

CMA Evolution Strategy with method \texttt{cmaes::cma_es}. Can handle numeric(vector) and integer(vector) hyperparameters, but no dependencies. For integers the internally proposed numeric values are automatically rounded. The sigma variance parameter is initialized to 1/4 of the span of box-constraints per parameter dimension.

Usage

\begin{verbatim}
makeTuneControlCMAES(
    same.resampling.instance = TRUE,
    impute.val = NULL,
    start = NULL,
    tune.threshold = FALSE,
    tune.threshold.args = list(),
    log.fun = "default",
    final.dw.perc = NULL,
    budget = NULL,
    ...
)
\end{verbatim}

Arguments

\begin{itemize}
  \item \texttt{same.resampling.instance} (logical(1))
    Should the same resampling instance be used for all evaluations to reduce variance? Default is \texttt{TRUE}.
  \item \texttt{impute.val} (numeric)
    If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or \texttt{Inf} instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.
  \item \texttt{start} (list)
    Named list of initial parameter values.
\end{itemize}
tune.threshold (logical(1))
Should the threshold be tuned for the measure at hand, after each hyperparameter evaluation, via tuneThreshold? Only works for classification if the predict type is “prob”. Default is FALSE.

tune.threshold.args
(list)
Further arguments for threshold tuning that are passed down to tuneThreshold. Default is none.

log.fun (function | character(1))
Function used for logging. If set to “default” (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to “memory” the memory usage for each evaluation will also be displayed, with character(1) small increase in run time. Otherwise character(1) function with arguments learner, resampling, measures, par.set, control, opt.path, dob, x, y, remove.nas, stage and prev.stage is expected. The default displays the performance measures, the time needed for evaluating, the currently used memory and the max memory ever used before (the latter two both taken from gc). See the implementation for details.

final.dw.perc (boolean)
If a Learner wrapped by a makeDownsampleWrapper is used, you can define
the value of dw.perc which is used to train the Learner with the final parameter setting found by the tuning. Default is NULL which will not change anything.

budget (integer(1))
Maximum budget for tuning. This value restricts the number of function evaluations. The budget corresponds to the product of the number of generations (maxit) and the number of offsprings per generation (lambda).

...
(any)
Further control parameters passed to the control arguments of cmaes::cma_es or GenSA::GenSA, as well as towards the tunerConfig argument of irace::irace.

Value
(TuneControlCMAES)

See Also
Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(), getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(), makeTuneControlIrace(), makeTuneControlMBO(), makeTuneControlRandom(), makeTuner(), makeWrapper(), tuneParams(), tuneThreshold()
**Description**

Completely pre-specify a data.frame of design points to be evaluated during tuning. All kinds of parameter types can be handled.

**Usage**

```r
makeTuneControlDesign(
  same.resampling.instance = TRUE,
  impute.val = NULL,
  design = NULL,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default"
)
```

**Arguments**

- `same.resampling.instance` *(logical(1))*
  Should the same resampling instance be used for all evaluations to reduce variance? Default is `TRUE`.

- `impute.val` *(numeric)*
  If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or `Inf` instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.

- `design` *(data.frame)*
  data.frame containing the different parameter settings to be evaluated. The columns have to be named according to the ParamSet which will be used in `tune()`. Proper designs can be created with `ParamHelpers::generateDesign` for instance.

- `tune.threshold` *(logical(1))*
  Should the threshold be tuned for the measure at hand, after each hyperparameter evaluation, via `tuneThreshold`? Only works for classification if the predict type is “prob”. Default is `FALSE`.

- `tune.threshold.args` *(list)*
  Further arguments for threshold tuning that are passed down to `tuneThreshold`. Default is none.

- `log.fun` *(function | character(1))*
  Function used for logging. If set to “default” (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to
makeTuneControlGenSA

Create control object for hyperparameter tuning with GenSA.

Description

Generalized simulated annealing with method GenSA::GenSA. Can handle numeric(vector) and integer(vector) hyperparameters, but no dependencies. For integers the internally proposed numeric values are automatically rounded.

Usage

makeTuneControlGenSA(
  same.resampling.instance = TRUE,
  impute.val = NULL,
  start = NULL,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default",
  final.dw.perc = NULL,
  budget = NULL,
  ...
)

Arguments

same.resampling.instance

(logical(1))
Should the same resampling instance be used for all evaluations to reduce variance? Default is TRUE.
impute.val \textbf{(numeric)}
If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or Inf instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.

\textbf{start} \textbf{(list)}
Named list of initial parameter values.

\textbf{tune.threshold} \textbf{(logical(1))}
Should the threshold be tuned for the measure at hand, after each hyperparameter evaluation, via \texttt{tuneThreshold}? Only works for classification if the predict type is “prob”. Default is \texttt{FALSE}.

\textbf{tune.threshold.args} \textbf{(list)}
Further arguments for threshold tuning that are passed down to \texttt{tuneThreshold}. Default is none.

\textbf{log.fun} \textbf{(function | character(1))}
Function used for logging. If set to “default” (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to “memory” the memory usage for each evaluation will also be displayed, with \texttt{character(1)} small increase in run time. Otherwise \texttt{character(1)} function with arguments \texttt{learner, resampling, measures, par.set, control, opt.path, dob, x, y, remove.nas, stage and prev.stage} is expected. The default displays the performance measures, the time needed for evaluating, the currently used memory and the max memory ever used before (the latter two both taken from \texttt{go}). See the implementation for details.

\textbf{final.dw.perc} \textbf{(boolean)}
If a Learner wrapped by a \texttt{makeDownsampleWrapper} is used, you can define the value of \texttt{dw.perc} which is used to train the Learner with the final parameter setting found by the tuning. Default is \texttt{NULL} which will not change anything.

\textbf{budget} \textbf{(integer(1))}
Maximum budget for tuning. This value restricts the number of function evaluations. \texttt{GenSA::GenSA} defines the budget via the argument \texttt{max.call}. However, one should note that this algorithm does not stop its local search before its end. This behavior might lead to an extension of the defined budget and will result in a warning.

\textbf{...} \textbf{(any)}
Further control parameters passed to the \texttt{control} arguments of \texttt{cmaes::cma_es} or \texttt{GenSA::GenSA}, as well as towards the \texttt{tunerConfig} argument of \texttt{irace::irace}.

\textbf{Value}

\texttt{(TuneControlGenSA)}.
makeTuneControlGrid

Create control object for hyperparameter tuning with grid search.

Description

A basic grid search can handle all kinds of parameter types. You can either use their correct type and resolution, or discretize them yourself by always using `ParamHelpers::makeDiscreteParam` in the `par.set` passed to `tuneParams`.

Usage

```r
makeTuneControlGrid(
  same.resampling.instance = TRUE,
  impute.val = NULL,
  resolution = 10L,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default",
  final.dw.perc = NULL,
  budget = NULL
)
```

Arguments

same.resampling.instance

(logical(1))

Should the same resampling instance be used for all evaluations to reduce variance? Default is `TRUE`.

impute.val

(numeric)

If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or `Inf` instead. For multi-criteria optimization pass a vector of imputations, one for each of your measures, in the same order as your measures.
makeTuneControlGrid

resolution  (integer)
Resolution of the grid for each numeric/integer parameter in par.set. For vector
parameters, it is the resolution per dimension. Either pass one resolution
for all parameters, or a named vector. See ParamHelpers::generateGridDesign.
Default is 10.

tune.threshold  (logical(1))
Should the threshold be tuned for the measure at hand, after each hyperparam-
eter evaluation, via tuneThreshold? Only works for classification if the predict
type is “prob”. Default is FALSE.

tune.threshold.args
(list)
Further arguments for threshold tuning that are passed down to tuneThreshold.
Default is none.

log.fun  (function | character(1))
Function used for logging. If set to “default” (the default), the evaluated design
points, the resulting performances, and the runtime will be reported. If set to
“memory” the memory usage for each evaluation will also be displayed, with
character(1) small increase in run time. Otherwise character(1) function
with arguments learner, resampling, measures, par.set, control, opt.path,
dob, x, y, remove.nas, stage and prev.stage is expected. The default displays
the performance measures, the time needed for evaluating, the currently
used memory and the max memory ever used before (the latter two both taken
from gc). See the implementation for details.

final.dw.perc  (boolean)
If a Learner wrapped by a makeDownsampleWrapper is used, you can define
the value of dw.perc which is used to train the Learner with the final parameter
setting found by the tuning. Default is NULL which will not change anything.

budget  (integer(1))
Maximum budget for tuning. This value restricts the number of function evalua-
tions. If set, must equal the size of the grid.

Value

(TuneControlGrid)

See Also

Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(),
getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(),
makeTuneControlCMAES(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlIrace(),
makeTuneControlMBO(), makeTuneControlRandom(), makeTuneWrapper(), tuneParams(), tuneThreshold()
makeTuneControlIrace  

Create control object for hyperparameter tuning with Irace.

Description

Tuning with iterated F-Racing with method irace::irace. All kinds of parameter types can be handled. We return the best of the final elite candidates found by irace in the last race. Its estimated performance is the mean of all evaluations ever done for that candidate. More information on irace can be found in the TR at http://iridia.ulb.ac.be/IridiaTrSeries/link/IridiaTr2011-004.pdf.

For resampling you have to pass a ResampleDesc, not a ResampleInstance. The resampling strategy is randomly instantiated n.instances times and these are the instances in the sense of irace (instances element of tunerConfig in irace::irace). Also note that irace will always store its tuning results in a file on disk, see the package documentation for details on this and how to change the file path.

Usage

makeTuneControlIrace(
  impute.val = NULL,
  n.instances = 100L,
  show.irace.output = FALSE,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default",
  final.dw.perc = NULL,
  budget = NULL,
  ...
)

Arguments

impute.val  (numeric)
If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or Inf instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.

n.instances  (integer(1))
Number of random resampling instances for irace, see details. Default is 100.

show.irace.output  (logical(1))
Show console output of irace while tuning? Default is FALSE.
**makeTuneControlMBO**

Create control object for hyperparameter tuning with MBO.

---

**Description**

Model-based / Bayesian optimization with the function `mlrMBO::mbo` from the `mlrMBO` package. Please refer to [https://github.com/mlr-org/mlrMBO](https://github.com/mlr-org/mlrMBO) for further info.

---

**tune.threshold** (logical(1))

Should the threshold be tuned for the measure at hand, after each hyperparameter evaluation, via `tuneThreshold`? Only works for classification if the predict type is "prob". Default is FALSE.

**tune.threshold.args** (list)

Further arguments for threshold tuning that are passed down to `tuneThreshold`. Default is none.

**log.fun** (function | character(1))

Function used for logging. If set to "default" (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to "memory" the memory usage for each evaluation will also be displayed, with character(1) small increase in run time. Otherwise character(1) function with arguments learner, resampling, measures, par.set, control, opt.path, dob, x, y, remove.nas, stage and prev.stage is expected. The default displays the performance measures, the time needed for evaluating, the currently used memory and the max memory ever used before (the latter two both taken from `gc`). See the implementation for details.

**final.dw.perc** (boolean)

If a Learner wrapped by a `makeDownsampleWrapper` is used, you can define the value of dw.perc which is used to train the Learner with the final parameter setting found by the tuning. Default is NULL which will not change anything.

**budget** (integer(1))

Maximum budget for tuning. This value restricts the number of function evaluations. It is passed to `maxExperiments`.

... (any)

Further control parameters passed to the control arguments of `cmaes::cma_es` or `GenSA::GenSA`, as well as towards the `tunerConfig` argument of `irace::irace`.

**Value**

(TuneControlIrace)

**See Also**

Other tune: `TuneControl`, `getNestedTuneResultsOptPathDf()`, `getNestedTuneResultsX()`, `getResamplingIndices()`, `getTuneResult()`, `makeModelMultiplexerParamSet()`, `makeModelMultiplexer()`, `makeTuneControlCMAES()`, `makeTuneControlDesign()`, `makeTuneControlGenSA()`, `makeTuneControlGrid()`, `makeTuneControlMBO()`, `makeTuneControlRandom()`, `makeTuneWrapper()`, `tuneParams()`, `tuneThreshold()`
Usage

```r
makeTuneControlMBO(
  same.resampling.instance = TRUE,
  impute.val = NULL,
  learner = NULL,
  mbo.control = NULL,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  continue = FALSE,
  log.fun = "default",
  final.dw.perc = NULL,
  budget = NULL,
  mbo.design = NULL
)
```

Arguments

- `same.resampling.instance` (logical(1))
  Should the same resampling instance be used for all evaluations to reduce variance? Default is `TRUE`.

- `impute.val` (numeric)
  If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or `Inf` instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.

- `learner` (Learner | NULL)
  The surrogate learner: A regression learner to model performance landscape. For the default, NULL, `mlrMBO` will automatically create a suitable learner based on the rules described in `mlrMBO::makeMBOLearner`.

- `mbo.control` (mlrMBO::MBOControl | NULL)
  Control object for model-based optimization tuning. For the default, NULL, the control object will be created with all the defaults as described in `mlrMBO::makeMBOControl`.

- `tune.threshold` (logical(1))
  Should the threshold be tuned for the measure at hand, after each hyperparameter evaluation, via `tuneThreshold`? Only works for classification if the predict type is “prob”. Default is `FALSE`.

- `tune.threshold.args` (list)
  Further arguments for threshold tuning that are passed down to `tuneThreshold`. Default is none.
makeTuneControlMBO

continue (logical(1))
Resume calculation from previous run using mlrMBO::mboContinue? Requires “save.file.path” to be set. Note that the ParamHelpers::OptPath in the mlrMBO::OptResult will only include the evaluations after the continuation. The complete OptPath will be found in the slot $mbo.result$opt.path.

log.fun (function | character(1))
Function used for logging. If set to “default” (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to “memory” the memory usage for each evaluation will also be displayed, with character(1) small increase in run time. Otherwise character(1) function with arguments learner, resampling, measures, par.set, control, opt.path, dob, x, y, remove.nas, stage and prev.stage is expected. The default displays the performance measures, the time needed for evaluating, the currently used memory and the max memory ever used before (the latter two both taken from gc). See the implementation for details.

final.dw.perc (boolean)
If a Learner wrapped by a makeDownsampleWrapper is used, you can define the value of dw.perc which is used to train the Learner with the final parameter setting found by the tuning. Default is NULL which will not change anything.

budget (integer(1))
Maximum budget for tuning. This value restricts the number of function evaluations.

mbo.design (data.frame | NULL)
Initial design as data frame. If the parameters have corresponding trafo functions, the design must not be transformed before it is passed! For the default, NULL, a default design is created like described in mlrMBO::mbo.

Value

(TuneControlMBO)

References


See Also

Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(), getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(), makeTuneControlCMAES(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(), makeTuneControlIrace(), makeTuneControlRandom(), makeTuneWrapper(), tuneParams(), tuneThreshold()
makeTuneControlRandom  
*Create control object for hyperparameter tuning with random search.*

**Description**

Random search. All kinds of parameter types can be handled.

**Usage**

```r
makeTuneControlRandom(
  same.resampling.instance = TRUE,
  maxit = NULL,
  tune.threshold = FALSE,
  tune.threshold.args = list(),
  log.fun = "default",
  final.dw.perc = NULL,
  budget = NULL
)
```

**Arguments**

- **same.resampling.instance**  
  (logical(1))  
  Should the same resampling instance be used for all evaluations to reduce variance? Default is TRUE.

- **maxit**  
  (integer(1) | NULL)  
  Number of iterations for random search. Default is 100.

- **tune.threshold**  
  (logical(1))  
  Should the threshold be tuned for the measure at hand, after each hyperparameter evaluation, via `tuneThreshold`? Only works for classification if the predict type is “prob”. Default is FALSE.

- **tune.threshold.args**  
  (list)  
  Further arguments for threshold tuning that are passed down to `tuneThreshold`. Default is none.

- **log.fun**  
  (function | character(1))  
  Function used for logging. If set to “default” (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to “memory” the memory usage for each evaluation will also be displayed, with character(1) small increase in run time. Otherwise character(1) function with arguments learner, resampling, measures, par.set, control, opt.path, dob, x, y, remove.nas, stage and prev.stage is expected. The default displays the performance measures, the time needed for evaluating, the currently used memory and the max memory ever used before (the latter two both taken from gc). See the implementation for details.
If a Learner wrapped by a `makeDownsampleWrapper` is used, you can define the value of `dw.perc` which is used to train the Learner with the final parameter setting found by the tuning. Default is `NULL` which will not change anything.

Maximum budget for tuning. This value restricts the number of function evaluations. The budget equals the number of iterations (`maxit`) performed by the random search algorithm.

Value

(TuneControlRandom)

See Also

Other tune: `TuneControl`, `getNestedTuneResultsOptPathDf()`, `getNestedTuneResultsX()`, `getResamplingIndices()`, `getTuneResult()`, `makeModelMultiplexerParamSet()`, `makeModelMultiplexer()`, `makeTuneControlCMAES()`, `makeTuneControlDesign()`, `makeTuneControlGenSA()`, `makeTuneControlGrid()`, `makeTuneControlIrace()`, `makeTuneControlMBO()`, `makeTuneWrapper()`, `tuneParams()`, `tuneThreshold()`

makeTuneWrapper

Fuse learner with tuning.

Description

Fuses a base learner with a search strategy to select its hyperparameters. Creates a learner object, which can be used like any other learner object, but which internally uses `tuneParams`. If the train function is called on it, the search strategy and resampling are invoked to select an optimal set of hyperparameter values. Finally, a model is fitted on the complete training data with these optimal hyperparameters and returned. See `tuneParams` for more details.

After training, the optimal hyperparameters (and other related information) can be retrieved with `getTuneResult`.

Usage

```r
makeTuneWrapper(
  learner,
  resampling,
  measures,
  par.set,
  control,
  show.info = getMlrOption("show.info")
)
```
Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

resampling (ResampleInstance | ResampleDesc)
Resampling strategy to evaluate points in hyperparameter space. If you pass a
description, it is instantiated once at the beginning by default, so all points are
evaluated on the same training/test sets. If you want to change that behavior,
look at TuneControl.

measures (list of Measure | Measure)
Performance measures to evaluate. The first measure, aggregated by the first
aggregation function is optimized, others are simply evaluated. Default is the
default measure for the task, see here getDefaultMeasure.

par.set (ParamHelpers::ParamSet)
Collection of parameters and their constraints for optimization. Dependent pa-
rameters with a requires field must use quote and not expression to define it.

color (TuneControl)
Control object for search method. Also selects the optimization algorithm for
tuning.

show.info (logical(1))
Print verbose output on console? Default is set via configureMlr.

Value

Learner.

See Also

Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(),
getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(),
makeTuneControlCMAES(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(),
makeTuneControlIrace(), makeTuneControlMBO(), makeTuneControlRandom(), tuneParams(),
tuneThreshold()

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(),
makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(),
makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(),
makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(),
makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(),
makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(),
makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeUndersampleWrapper(),
makeWeightedClassesWrapper()

Examples

task = makeClassifTask(data = iris, target = "Species")
lrn = makeLearner("classif.rpart")
### makeUndersampleWrapper

_Fuse learner with simple over/undersampling for imbalancy correction in binary classification._

**Description**

Creates a learner object, which can be used like any other learner object. Internally uses `oversample` or `undersample` before every model fit.

Note that observation weights do not influence the sampling and are simply passed down to the next learner.

**Usage**

```r
makeUndersampleWrapper(learner, usw.rate = 1, usw.cl = NULL)
makeOversampleWrapper(learner, osw.rate = 1, osw.cl = NULL)
```

**Arguments**

- **learner** (_Learner | character(1))
  The learner. If you pass a string the learner will be created via `makeLearner`.

- **usw.rate** (_numeric(1))
  Factor to downsample a class. Must be between 0 and 1, where 1 means no downsampling, 0.5 implies reduction to 50 percent and 0 would imply reduction to 0 observations. Default is 1.

- **usw.cl** (_character(1))
  Class that should be undersampled. Default is NULL, which means the larger one.
makeWeightedClassesWrapper

osw.rate (numeric(1))
Factor to oversample a class. Must be between 1 and Inf, where 1 means no oversampling and 2 would mean doubling the class size. Default is 1.

osw.cl (character(1))
Class that should be oversampled. Default is NULL, which means the smaller one.

Value
Learner.

See Also
Other imbalancy: makeOverBaggingWrapper(), oversample(), smote()
Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(), makeConstantClassWrapper(), makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(), makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(), makeFilterWrapper(), makeImputeWrapper(), makeMulticlassWrapper(), makeMultilabelBinaryRelevanceWrapper(), makeMultilabelClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(), makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(), makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(), makeWeightedClassesWrapper()

makeWeightedClassesWrapper
Wraps a classifier for weighted fitting where each class receives a weight.

Description
Creates a wrapper, which can be used like any other learner object.
Fitting is performed in a weighted fashion where each observation receives a weight, depending on the class it belongs to, see wcw.weight. This might help to mitigate problems caused by imbalanced class distributions.

This weighted fitting can be achieved in two ways:
a) The learner already has a parameter for class weighting, so one weight can directly be defined per class. Example: “classif.ksvm” and parameter class.weights. In this case we don't really do anything fancy. We convert wcw.weight a bit, but basically simply bind its value to the class weighting param. The wrapper in this case simply offers a convenient, consistent fashion for class weighting - and tuning! See example below.
b) The learner does not have a direct parameter to support class weighting, but supports observation weights, so hasLearnerProperties(learner, 'weights') is TRUE. This means that an individual, arbitrary weight can be set per observation during training. We set this weight depending on the class internally in the wrapper. Basically we introduce something like a new “class.weights” parameter for the learner via observation weights.
makeWeightedClassesWrapper

Usage

makeWeightedClassesWrapper(learner, wcw.param = NULL, wcw.weight = 1)

Arguments

learner (Learner | character(1))
The classification learner. If you pass a string the learner will be created via
makeLearner.

wcw.param (character(1))
Name of already existing learner parameter, which allows class weighting. The
default (wcw.param = NULL) will use the parameter defined in the learner (class.weights.param).
During training, the parameter must accept a named vector of class weights,
where length equals the number of classes.

wcw.weight (numeric)
Weight for each class. Must be a vector of the same number of elements as
classes are in task, and must also be in the same order as the class levels are in
getTaskDesc(task)$class.levels. For convenience, one must pass a single
number in case of binary classification, which is then taken as the weight of the
positive class, while the negative class receives a weight of 1. Default is 1.

Value

Learner.

See Also

Other wrapper: makeBaggingWrapper(), makeClassificationViaRegressionWrapper(),
makeCostSensClassifWrapper(), makeCostSensRegrWrapper(), makeDownsampleWrapper(),
makeDummyFeaturesWrapper(), makeExtractFDAFeatsWrapper(), makeFeatSelWrapper(),
makeImputeWrapper(), makeMulticlassWrapper(), makeMulticlassBinaryRelevanceWrapper(),
makeMulticlassClassifierChainsWrapper(), makeMultilabelDBRWrapper(), makeMultilabelNestedStackingWrapper(),
makeMultilabelStackingWrapper(), makeOverBaggingWrapper(), makePreprocWrapperCaret(),
makePreprocWrapper(), makeRemoveConstantFeaturesWrapper(), makeSMOTEWrapper(), makeTuneWrapper(),
makeUndersampleWrapper()

Examples

set.seed(123)
# using the direct parameter of the SVM (which is already defined in the learner)
lrn = makeWeightedClassesWrapper(“classif.ksvm”, wcw.weight = 0.01)
res = holdout(lrn, sonar.task)
print(calculateConfusionMatrix(res$pred))

# using the observation weights of logreg
lrn = makeWeightedClassesWrapper(“classif.logreg”, wcw.weight = 0.01)
res = holdout(lrn, sonar.task)
print(calculateConfusionMatrix(res$pred))

# tuning the imbalance param and the SVM param in one go
makeWrappedModel

Induced model of learner.

Description

Result from train.

It internally stores the underlying fitted model, the subset used for training, features used for training, levels of factors in the data set and computation time that was spent for training.

Object members: See arguments.

The constructor makeWrappedModel is mainly for internal use.

Usage

makeWrappedModel(
  learner,
  learner.model,
  task.desc,
  subset,
  features,
  factor.levels,
  time
)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

learner.model (any)
Underlying model.

task.desc TaskDesc
Task description object.

subset (integer | logical | NULL)
Selected cases. Either a logical or an index vector. By default NULL if all observations are used.
**MeasureProperties**

- **features**: (character)
  - Features used for training.

- **factor.levels**: (named list of character)
  - Levels of factor variables (features and potentially target) in training data. Named by variable name, non-factors do not occur in the list.

- **time**: (numeric(1))
  - Computation time for model fit in seconds.

**Value**

- **WrappedModel**.

---

**MeasureProperties**

- **Query properties of measures.**

**Description**

Properties can be accessed with `getMeasureProperties(measure)`, which returns a character vector.

The measure properties are defined in **Measure**.

**Usage**

```
getMeasureProperties(measure)
```

```
hasMeasureProperties(measure, props)
```

**Arguments**

- **measure**: (Measure)
  - Performance measure. Default is the first measure used in the benchmark experiment.

- **props**: (character)
  - Vector of properties to query.

**Value**

- `getMeasureProperties` returns a character vector with measure properties. `hasMeasureProperties` returns a logical vector of the same length as `props`. 
Performance measures.

Description

A performance measure is evaluated after a single train/predict step and returns a single number to assess the quality of the prediction (or maybe only the model, think AIC). The measure itself knows whether it wants to be minimized or maximized and for what tasks it is applicable.

All supported measures can be found by listMeasures or as a table in the tutorial appendix: https://mlr.mlr-org.com/articles/tutorial/measures.html.

If you want a measure for a misclassification cost matrix, look at makeCostMeasure. If you want to implement your own measure, look at makeMeasure.

Most measures can directly be accessed via the function named after the scheme measureX (e.g. measureSSE).

For clustering measures, we compact the predicted cluster IDs such that they form a continuous series starting with 1. If this is not the case, some of the measures will generate warnings.

Some measure have parameters. Their defaults are set in the constructor makeMeasure and can be overwritten using setMeasurePars.

Usage

featperc
timetrain
timepredict
timeboth
sse
measureSSE(truth, response)
mse
measureMSE(truth, response)
rmse
measureRMSE(truth, response)
medse
measureMEDSE(truth, response)
sae
measureSAE(truth, response)

mae
measureMAE(truth, response)

medae
measureMEDAE(truth, response)

rsq
measureRSQ(truth, response)

expvar
measureEXPVAR(truth, response)

arsq

rrse
measureRRSE(truth, response)

rae
measureRAE(truth, response)

mape
measureMAPE(truth, response)

msle
measureMSLE(truth, response)

rmsle
measureRMSLE(truth, response)

kendalltau
measureKendallTau(truth, response)

spearmanrho
measuresSpearmanRho(truth, response)
mnce
measuresMMCE(truth, response)
acc
measuresACC(truth, response)
ber
measuresBER(truth, response)
multiclass.aunu
measuresAUNU(probabilities, truth)
multiclass.aump
measuresAUNP(probabilities, truth)
multiclass.au1u
measuresAU1U(probabilities, truth)
multiclass.au1p
measuresAU1P(probabilities, truth)
multiclass.brier
measuresMulticlassBrier(probabilities, truth)
logloss
measuresLogloss(probabilities, truth)
ssr
measuresSSR(probabilities, truth)
qsr
measuresQSR(probabilities, truth)
lsr
measures

measureLSR(probabilities, truth)

kappa

measureKAPPA(truth, response)

wkappa

measureWKAPPA(truth, response)

auc

measureAUC(probabilities, truth, negative, positive)

brier

measureBrier(probabilities, truth, negative, positive)

brier.scaled

measureBrierScaled(probabilities, truth, negative, positive)

bac

measureBAC(truth, response)

tp

measureTP(truth, response, positive)

tn

measureTN(truth, response, negative)

fp

measureFP(truth, response, positive)

fn

measureFN(truth, response, negative)

tpr

measureTPR(truth, response, positive)

tnr
measureTNR(truth, response, negative)

fpr

measureFPR(truth, response, negative, positive)

fnr

measureFNR(truth, response, negative, positive)

ppv

measurePPV(truth, response, positive, probabilities = NULL)

npv

measureNPV(truth, response, negative)

fdr

measureFDR(truth, response, positive)

mcc

measureMCC(truth, response, negative, positive)

f1

measureF1(truth, response, positive)

gmean

measureGMEAN(truth, response, negative, positive)

gpr

measureGPR(truth, response, positive)

multilabel.hamloss

measureMultilabelHamloss(truth, response)

multilabel.subset01

measureMultilabelSubset01(truth, response)

multilabel.f1
measureMultilabelF1(truth, response)
multilabel.acc
measureMultilabelACC(truth, response)
multilabel.ppv
measureMultilabelPPV(truth, response)
multilabel.tpr
measureMultilabelTPR(truth, response)
cindex
cindex.uno
iauc.uno
ibrier
meancosts
mcp
db
dunn
G1
G2
silhouette

Arguments

truth
(factor)
Vector of the true class.
response
(factor)
Vector of the predicted class.
probabilities
(numeric | matrix)
a) For purely binary classification measures: The predicted probabilities for the positive class as a numeric vector. b) For multiclass classification measures: The predicted probabilities for all classes, always as a numeric matrix, where columns are named with class labels.
negative (character(1))
The name of the negative class.

positive (character(1))
The name of the positive class.

Format
none

References


See Also
Other performance: ConfusionMatrix, calculateConfusionMatrix(), calculateROCMeasures(), estimateRelativeOverfitting(), makeCostMeasure(), makeCustomResampledMeasure(), makeMeasure(), performance(), setAggregation(), setMeasurePars()

mergeBenchmarkResults Merge different BenchmarkResult objects.

Description
The function automatically combines a list of [BenchmarkResult] objects into a single [BenchmarkResult] object as long as the full crossproduct of all task-learner combinations are available.

Usage
mergeBenchmarkResults(bmrs)

Arguments
bmrs [list of [BenchmarkResult]]
‘BenchmarkResult’ objects that should be merged.

Details
Note that if you want to merge several [BenchmarkResult] objects, you must ensure that all possible learner and task combinations will be contained in the returned object. Otherwise, the user will be notified which task-learner combinations are missing or duplicated. When merging [BenchmarkResult] objects with different measures, all missing measures will automatically be recomputed.
**mergeSmallFactorLevels**

*Merges small levels of factors into new level.*

**Description**

Merges factor levels that occur only infrequently into combined levels with a higher frequency.

**Usage**

```r
mergeSmallFactorLevels(
  task, 
  cols = NULL, 
  min.perc = 0.01, 
  new.level = ".merged"
)
```

**Arguments**

- **task** *(Task)*
  The task.

- **cols** *(character)*
  Which columns to convert. Default is all factor and character columns.

- **min.perc** *(numeric(1))*
  The smallest levels of a factor are merged until their combined proportion w.r.t. the length of the factor exceeds `min.perc`. Must be between 0 and 1. Default is 0.01.

- **new.level** *(character(1))*
  New name of merged level. Default is ".merged"

**Value**

Task, where merged levels are combined into a new level of name `new.level`.

**See Also**

Other eda_and_preprocess: `capLargeValues()`, `createDummyFeatures()`, `dropFeatures()`, `normalizeFeatures()`, `removeConstantFeatures()`, `summarizeColumns()`, `summarizeLevels()`
mlrFamilies

mlr documentation families

Description
List of all mlr documentation families with members.

Arguments
benchmark
  batchmark, reduceBatchmarkResults, benchmark, benchmarkParallel, getBMRTaskIds, getBMR Learners, getBMR LearnerIds, getBMR Learner ShortNames, getBMR Measures, getBMR MeasureIds, getBMR Predictions, getBMR Performances, getBMR Aggr Performances, getBMR Tune Results, getBMR FeatureSel Results, getBMR Filtered Features, getBMR Models, getBMR TaskDescs, convertBMRToRankMatrix, friedmanPostHocTestBMR, friedmanTestBMR, plotBMRBoxplots, plotBMR Ranks As Bar Chart, generateCritDifferencesData, plotCritDifferences

calibration
  generateCalibrationData, plotCalibration

configure
  configureMlr, getMlrOptions

costsens
  makeCostSensTask, makeCostSens Weighted Pairs Wrapper

debug
  predictFailureModel, getPredictionDump, getRRDump, print.ResampleResult
downsample
downsampling
downsampling
downsampling
eda_and_preprocess
capLargeValues, createDummyFeatures, dropFeatures, mergeSmallFactorLevels, normalizeFeatures, removeConstantFeatures, summarizeColumns, summarizeLevels

extractFDAFeatures
  reextractFDAFeatures

defda_featextractor
  extractFDA Fourier, extractFDA Wavelets, extractFDA FPCA, extractFDA Multi Res Features

defda
  makeExtractFDAFeatMethod, extractFDAFeatures

dfeatsel
  analyzeFeatSelResult, makeFeatSelControl, getFeatSelResult, selectFeatures

filter
  filterFeatures, makeFilter, listFilterMethods, getFilteredFeatures, generateFilterValuesData, getFilterValues

generate_plot_data
generateFeatureImportanceData, plotFilterValues, generatePartialDependenceData

help
  help Learner, help Learner Param

imbalance
  oversample, smote

impute
  makeImputeMethod, impute Constant, impute, reimpute
### mtcars.task

**Description**

Contains the task (mtcars.task).
References

See datasets::mtcars.

normalizeFeatures Normalize features.

Description

Normalize features by different methods. Internally BBmisc::normalize is used for every feature column. Non numerical features will be left untouched and passed to the result. For constant features most methods fail, special behaviour for this case is implemented.

Usage

normalizeFeatures(
  obj, 
  target = character(0L), 
  method = "standardize", 
  cols = NULL, 
  range = c(0, 1), 
  on.constant = "quiet"
)

Arguments

obj (data.frame | Task)
Input data.

target (character(1) | character(2) | character(n.classes))
Name(s) of the target variable(s). Only used when obj is a data.frame, otherwise ignored. If survival analysis is applicable, these are the names of the survival time and event columns, so it has length 2. For multilabel classification these are the names of logical columns that indicate whether a class label is present and the number of target variables corresponds to the number of classes.

method (character(1))
Normalizing method. Available are:
“center”: Subtract mean.
“scale”: Divide by standard deviation.
“standardize”: Center and scale.
“range”: Scale to a given range.

cols (character)
Columns to normalize. Default is to use all numeric columns.

range (numeric(2))
Range for method “range”. Default is c(0,1).
How should constant vectors be treated? Only used, of “method != center”, since this method does not fail for constant vectors. Possible actions are: “quiet”: Depending on the method, treat them quietly: “scale”: No division by standard deviation is done, input values. will be returned untouched. “standardize”: Only the mean is subtracted, no division is done. “range”: All values are mapped to the mean of the given range. “warn”: Same behaviour as “quiet”, but print a warning message. “stop”: Stop with an error.

Value

data.frame | Task. Same type as obj.

See Also

BBmisc::normalize

Other eda_and_preprocess: capLargeValues(), createDummyFeatures(), dropFeatures(), mergeSmallFactorLevels(), removeConstantFeatures(), summarizeColumns(), summarizeLevels()

Description

Oversampling: For a given class (usually the smaller one) all existing observations are taken and copied and extra observations are added by randomly sampling with replacement from this class. Undersampling: For a given class (usually the larger one) the number of observations is reduced (downsampled) by randomly sampling without replacement from this class.

Usage

oversample(task, rate, cl = NULL)

undersample(task, rate, cl = NULL)

Arguments

<table>
<thead>
<tr>
<th>task</th>
<th>(Task)</th>
</tr>
</thead>
<tbody>
<tr>
<td>The task.</td>
<td></td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>rate</th>
<th>(numeric(1))</th>
</tr>
</thead>
<tbody>
<tr>
<td>Factor to upsample or downsample a class. For undersampling: Must be between 0 and 1, where 1 means no downsampling, 0.5 implies reduction to 50 percent and 0 would imply reduction to 0 observations. For oversampling: Must</td>
<td></td>
</tr>
</tbody>
</table>
be between 1 and Inf, where 1 means no oversampling and 2 would mean doubling the class size.

\[ \text{cl} \ (\text{character}(1)) \]

Which class should be over- or undersampled. If NULL, oversample will select the smaller and undersample the larger class.

**Value**

Task.

**See Also**

Other imbalancy: `makeOverBaggingWrapper()`, `makeUndersampleWrapper()`, `smote()`

---

### parallelization

**Supported parallelization methods**

**Description**

mlr supports different methods to activate parallel computing capabilities through the integration of the `parallelMap::parallelMap` package, which supports all major parallelization backends for R. You can start parallelization with `parallelStart*`, where * should be replaced with the chosen backend. `parallelMap::parallelStop` is used to stop all parallelization backends.

Parallelization is divided into different levels and will automatically be carried out for the first level that occurs, e.g. if you call `resample()` after `parallelMap::parallelStart`, each resampling iteration is a parallel job and possible underlying calls like parameter tuning won’t be parallelized further.

The supported levels of parallelization are:

- "mlr.resample" Each resampling iteration (a train/test step) is a parallel job.
- "mlr.benchmark" Each experiment "run this learner on this data set" is a parallel job.
- "mlr.tuneParams" Each evaluation in hyperparameter space "resample with these parameter settings" is a parallel job. How many of these can be run independently in parallel depends on the tuning algorithm. For grid search or random search there is no limit, but for other tuners it depends on how many points to evaluate are produced in each iteration of the optimization. If a tuner works in a purely sequential fashion, we cannot work magic and the hyperparameter evaluation will also run sequentially. But note that you can still parallelize the underlying resampling.
- "mlr.selectFeatures" Each evaluation in feature space "resample with this feature subset" is a parallel job. The same comments as for "mlr.tuneParams" apply here.
- "mlr.ensemble" For all ensemble methods, the training and prediction of each individual learner is a parallel job. Supported ensemble methods are the `makeBaggingWrapper`, `makeCostSensitiveRegrWrapper`, `makeMulticlassWrapper`, `makeMultilabelBinaryRelevanceWrapper` and the `makeOverBaggingWrapper`. 
Description

Measures the quality of a prediction w.r.t. some performance measure.

Usage

performance(
  pred, 
  measures, 
  task = NULL, 
  model = NULL, 
  feats = NULL, 
  simpleaggr = FALSE 
)

Arguments

pred (Prediction)  Prediction object.
measures (Measure | list of Measure)  Performance measure(s) to evaluate. Default is the default measure for the task, see here getDefaultMeasure.
task (Task)  Learning task, might be requested by performance measure, usually not needed except for clustering or survival.
model (WrappedModel)  Model built on training data, might be requested by performance measure, usually not needed except for survival.
feats (data.frame)  Features of predicted data, usually not needed except for clustering. If the prediction was generated from a task, you can also pass this instead and the features are extracted from it.
simpleaggr (logical)  If TRUE, aggregation of ResamplePrediction objects is skipped. This is used internally for threshold tuning. Default is FALSE.

Value

(named numeric). Performance value(s), named by measure(s).
See Also

Other performance: `ConfusionMatrix`, `calculateConfusionMatrix()`, `calculateROCMeasures()`, `estimateRelativeOverfitting()`, `makeCostMeasure()`, `makeCustomResampledMeasure()`, `makeMeasure()`, `measures`, `setAggregation()`, `setMeasurePars()`

Examples

```r
training.set = seq(1, nrow(iris), by = 2)
test.set = seq(2, nrow(iris), by = 2)

task = makeClassifTask(data = iris, target = "Species")
ln = makeLearner("classif.lda")
mod = train(ln, task, subset = training.set)
pred = predict(mod, newdata = iris[test.set, ])
performance(pred, measures = mmce)

# Compute multiple performance measures at once
ms = list("mmce" = mmce, "acc" = acc, "timetrain" = timetrain)
performance(pred, measures = ms, task, mod)
```

---

**phoneme.task**  
*Phoneme functional data multilabel classification task.*

**Description**

Contains the task (phoneme.task). The task contains a single functional covariate and 5 equally big classes (aa, ao, dcl, iy, sh). The aim is to predict the class of the phoneme in the functional. The dataset is contained in the package fda.usc.

**References**


---

**pid.task**  
*PimaIndiansDiabetes classification task.*

**Description**

Contains the task (pid.task).

**References**

See `mlbench::PimaIndiansDiabetes`. Note that this is the uncorrected version from mlbench.
plotBMRBoxplots

Create box or violin plots for a BenchmarkResult.

Description

Plots box or violin plots for a selected measure across all iterations of the resampling strategy, faceted by the task.id.

Usage

plotBMRBoxplots(
  bmr,
  measure = NULL,
  style = "box",
  order.lrns = NULL,
  order.tsks = NULL,
  pretty.names = TRUE,
  facet.wrap.nrow = NULL,
  facet.wrap.ncol = NULL
)

Arguments

bmr (BenchmarkResult)
Benchmark result.

measure (Measure)
Performance measure. Default is the first measure used in the benchmark experiment.

style (character(1))
Type of plot, can be “box” for a boxplot or “violin” for a violin plot. Default is “box”.

order.lrns (character(n.learners))
Character vector with learner.ids in new order.

order.tsks (character(n.tasks))
Character vector with task.ids in new order.

pretty.names (logical(1))
Whether to use the Measure name and the Learner short name instead of the id.
Default is TRUE.

facet.wrap.nrow, facet.wrap.ncol (integer)
Number of rows and columns for facetting. Default for both is NULL. In this case ggplot's facet_wrap will choose the layout itself.

Value

ggplot2 plot object.
See Also

Other plot: createSpatialResamplingPlots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCalibration(), plotCritDifferences(), plotLearningCurve(), plotPartialDependence(), plotROCCurves(), plotResiduals(), plotThreshVsPerf()

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRfilteredFeatures(), getBMRlearnerIds(), getBMRlearnerShortNames(), getBMRlearners(), getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMRtuneResults(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

Examples

# see benchmark

```
plotBMRRanksAsBarChart

Create a bar chart for ranks in a BenchmarkResult.
```

Description

Plots a bar chart from the ranks of algorithms. Alternatively, tiles can be plotted for every rank-task combination, see pos for details. In all plot variants the ranks of the learning algorithms are displayed on the x-axis. Areas are always colored according to the learner.id.

Usage

```
plotBMRRanksAsBarChart(
  bmr,
  measure = NULL,
  ties.method = "average",
  aggregation = "default",
  pos = "stack",
  order.lrns = NULL,
  order.tsks = NULL,
  pretty.names = TRUE
)
```

Arguments

- **bmr** *(BenchmarkResult)*
  Benchmark result.
- **measure** *(Measure)*
  Performance measure. Default is the first measure used in the benchmark experiment.
plotBMRSummary

```r
ties.method (character(1))
See rank for details.

aggregation (character(1))
"mean" or "default". See getBMRAggrPerformances for details on "default".

pos (character(1))
Optionally set how the bars are positioned in ggplot2. Ranks are plotted on the x-axis. "tile" plots a heat map with task as the y-axis. Allows identification of the performance in a special task. "stack" plots a stacked bar plot. Allows for comparison of learners within and across ranks. "dodge" plots a bar plot with bars next to each other instead of stacked bars.

order.lrns (character(n.learners))
Character vector with learner.ids in new order.

order.tsks (character(n.tasks))
Character vector with task.ids in new order.

pretty.names (logical(1))
Whether to use the short name of the learner instead of its ID in labels. Defaults to TRUE.
```

Value
ggplot2 plot object.

See Also

Other plot: createSpatialResamplingPlots(), plotBMRBoxplots(), plotBMRSummary(), plotCalibration(), plotCritDifferences(), plotLearningCurve(), plotPartialDependence(), plotROCCurves(), plotResiduals(), plotThreshVsPerf()

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMR LearnerIds(), getBMR Learner Short Names(), getBMR Learners(), getBMR Measure Ids(), getBMR Measures(), getBMR Models(), getBMR Performances(), getBMR Predictions(), getBMR Task Descs(), getBMR Task Ids(), getBMR Tune Results(), plotBMRBoxplots(), plotBMRSummary(), plotCritDifferences(), reduceBatchmarkResults()

Examples

```r
# see benchmark
```
Usage

```r
plotBMRSummary(
  bmr,
  measure = NULL,
  trafo = "none",
  order.tsks = NULL,
  pointsize = 4L,
  jitter = 0.05,
  pretty.names = TRUE
)
```

Arguments

- `bmr` *(BenchmarkResult)*
  Benchmark result.
- `measure` *(Measure)*
  Performance measure. Default is the first measure used in the benchmark experiment.
- `trafo` *(character(1))*
  Currently either “none” or “rank”, the latter performing a rank transformation (with average handling of ties) of the scores per task. NB: You can add always add `ggplot2::scale_x_log10` to the result to put scores on a log scale. Default is “none”.
- `order.tsks` *(character(n.tasks))*
  Character vector with task.ids in new order.
- `pointsize` *(numeric(1))*
  Point size for `ggplot2 ggplot2::geom_point` for data points. Default is 4.
- `jitter` *(numeric(1))*
  Small vertical jitter to deal with overplotting in case of equal scores. Default is 0.05.
- `pretty.names` *(logical(1))*
  Whether to use the short name of the learner instead of its ID in labels. Defaults to TRUE.

Value

`ggplot2` plot object.

See Also

Other benchmark: `BenchmarkResult`, `benchmark()`, `convertBMRToRankMatrix()`, `friedmanPostHocTestBMR()`, `friedmanTestBMR()`, `generateCritDifferencesData()`, `getBMRaggrPerformances()`, `getBMRFeatSelResults()`, `getBMRFilteredFeatures()`, `getBMR LearnerIds()`, `getBMR Learner ShortNames()`, `getBMR Learners()`, `getBMRMeasureIds()`, `getBMRMeasures()`, `getBMRModels()`, `getBMR Performances()`, `getBMR Predictions()`, `getBMR TaskDescs()`, `getBMR TaskIds()`, `getBMR Tune Results()`, `plotBMR Boxplots()`, `plotBMR Ranks As Bar Chart()`, `plot Crit Differences()`, `reduceBatchmarkResults()`
plotCalibration

Other plot: createSpatialResamplingPlots(), plotBMRRankAsBarChart(), plotCalibration(), plotCritDifferences(), plotLearningCurve(), plotPartialDependence(), plotROCCurves(), plotResiduals(), plotThreshVsPerf()

Examples

# see benchmark

---

plotCalibration  Plot calibration data using ggplot2.

Description

Plots calibration data from generateCalibrationData.

Usage

plotCalibration(
  obj,
  smooth = FALSE,
  reference = TRUE,
  rag = TRUE,
  facet.wrap.nrow = NULL,
  facet.wrap.ncol = NULL
)

Arguments

obj  (CalibrationData)
  Result of generateCalibrationData.
smooth  (logical(1))
  Whether to use a loess smoother. Default is FALSE.
reference  (logical(1))
  Whether to plot a reference line showing perfect calibration. Default is TRUE.
rag  (logical(1))
  Whether to include a rag plot which shows a rug plot on the top which pertains to positive cases and on the bottom which pertains to negative cases. Default is TRUE.
facet.wrap.nrow, facet.wrap.ncol  (integer)
  Number of rows and columns for facetting. Default for both is NULL. In this case ggplot's `facet_wrap` will choose the layout itself.

Value

ggplot2 plot object.
See Also

Other plot: `createSpatialResamplingPlots()`, `plotBMRRanksAsBarChart()`, `plotBMRSummary()`, `plotCritDifferences()`, `plotLearningCurve()`, `plotPartialDependence()`, `plotROCCurves()`, `plotResiduals()`, `plotThreshVsPerf()

Other calibration: `generateCalibrationData()`

Examples

```r
## Not run:
lrns = list(makeLearner("classif.rpart", predict.type = "prob"),
            makeLearner("classif.nnet", predict.type = "prob"))
fit = lapply(lrns, train, task = iris.task)
pred = lapply(fit, predict, task = iris.task)
names(pred) = c("rpart", "nnet")
out = generateCalibrationData(pred, groups = 3)
plotCalibration(out)

fit = lapply(lrns, train, task = sonar.task)
pred = lapply(fit, predict, task = sonar.task)
names(pred) = c("rpart", "lda")
out = generateCalibrationData(pred)
plotCalibration(out)

## End(Not run)
```

```r
plotCritDifferences  Plot critical differences for a selected measure.
```

Description

Plots a critical-differences diagram for all classifiers and a selected measure. If a baseline is selected for the Bonferroni-Dunn test, the critical difference interval will be positioned around the baseline. If not, the best performing algorithm will be chosen as baseline. The positioning of some descriptive elements can be moved by modifying the generated data.

Usage

```r
plotCritDifferences(obj, baseline = NULL, pretty.names = TRUE)
```

Arguments

- `obj` ([critDifferencesData]) Result of `generateCritDifferencesData` function.
- `baseline` (`character(1)`): ([learner.id])
  Overwrites baseline from `generateCritDifferencesData`!
  Select a `[learner.id` as baseline for the critical difference diagram, the critical difference will be positioned around this learner. Defaults to best performing algorithm.
plotFilterValues

pretty.names (logical(1))
Whether to use the short name of the learner instead of its ID in labels. Defaults to TRUE.

Value

ggplot2 plot object.

References

Janez Demsar, Statistical Comparisons of Classifiers over Multiple Data Sets, JMLR, 2006

See Also

Other plot: createSpatialResamplingPlots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCalibration(), plotLearningCurve(), plotPartialDependence(), plotROCCurves(), plotResiduals(), plotThreshVsPerf()

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRLearners(), getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMR TuneResults(), plotBMROxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), reduceBatchmarkResults()

Examples

# see benchmark

plotFilterValues

Plot filter values using ggplot2.

Description

Plot filter values using ggplot2.

Usage

plotFilterValues(
  fvalues,
  sort = "dec",
  n.show = nrow(fvalues$data),
  filter = NULL,
  feat.type.cols = FALSE
)
**Arguments**

- `fvalues` *(FilterValues)*
  Filter values.

- `sort` *(character(1))*
  Available options are:
  - "dec" -> descending
  - "inc" -> increasing
  - "none" -> no sorting
  Default is decreasing.

- `n.show` *(integer(1))*
  Number of features (maximal) to show. Default is to plot all features.

- `filter` *(character(1))*
  In case `fvalues` contains multiple filter methods, which method should be plotted?

- `feat.type.cols` *(logical(1))*
  Whether to color different feature types (e.g. numeric | factor). Default is to use no colors (`feat.type.cols = FALSE`).

**Value**

`ggplot2` plot object.

**See Also**

Other filter: `filterFeatures()`, `generateFilterValuesData()`, `getFilteredFeatures()`, `listFilterEnsembleMethods()`, `makeFilterEnsemble()`, `makeFilterWrapper()`, `makeFilter()`

Other generate_plot_data: `generateCalibrationData()`, `generateCritDifferencesData()`, `generateFeatureImportanceData()`, `generateFilterValuesData()`, `generateLearningCurveData()`, `generatePartialDependenceData()`, `generateThreshVsPerfData()`

**Examples**

```r
fv = generateFilterValuesData(iris.task, method = "variance")
plotFilterValues(fv)
```

---

**Description**

Plot hyperparameter validation path. Automated plotting method for `HyperParsEffectData` object. Useful for determining the importance or effect of a particular hyperparameter on some performance measure and/or optimizer.
Usage

plotHyperParsEffect(
  hyperpars.effect.data,
  x = NULL,
  y = NULL,
  z = NULL,
  plot.type = "scatter",
  loess.smooth = FALSE,
  facet = NULL,
  global.only = TRUE,
  interpolate = NULL,
  show.experiments = FALSE,
  show.interpolated = FALSE,
  nested agg = mean,
  partial.dep.learn = NULL
)

Arguments

hyperpars.effect.data
  (HyperParsEffectData)
  Result of generateHyperParsEffectData

x
  (character(1))
  Specify what should be plotted on the x axis. Must be a column from HyperParsEffectData$data. For partial dependence, this is assumed to be a hyperparameter.

y
  (character(1))
  Specify what should be plotted on the y axis. Must be a column from HyperParsEffectData$data

z
  (character(1))
  Specify what should be used as the extra axis for a particular geom. This could be the fill on a heatmap or color aesthetic for a line. Must be a column from HyperParsEffectData$data. Default is NULL.

plot.type
  (character(1))
  Specify the type of plot: “scatter” for a scatterplot, “heatmap” for a heatmap, “line” for a scatterplot with a connecting line, or “contour” for a contour plot layered on top of a heatmap. Default is “scatter”.

loess.smooth
  (logical(1))
  If TRUE, will add loess smoothing line to plots where possible. Note that this is probably only useful when plot.type is set to either “scatter” or “line”. Must be a column from HyperParsEffectData$data. Not used with partial dependence. Default is FALSE.

facet
  (character(1))
  Specify what should be used as the facet axis for a particular geom. When using nested cross validation, set this to “nested_cv_run” to obtain a facet for each outer loop. Must be a column from HyperParsEffectData$data. Please note that facetting is not supported with partial dependence plots! Default is NULL.

global.only
  (logical(1))
  If TRUE, will only plot the current global optima when setting x = "iteration" and
y as a performance measure from `HyperParsEffectData$measures`. Set this to FALSE to always plot the performance of every iteration, even if it is not an improvement. Not used with partial dependence. Default is TRUE.

```r
interpolate (Learner | character(1))
```

If not NULL, will interpolate non-complete grids in order to visualize a more complete path. Only meaningful when attempting to plot a heatmap or contour. This will fill in “empty” cells in the heatmap or contour plot. Note that cases of irregular hyperparameter paths, you will most likely need to use this to have a meaningful visualization. Accepts either a regression Learner object or the learner as a string for interpolation. This cannot be used with partial dependence. Default is NULL.

```r
show.experiments (logical(1))
```

If TRUE, will overlay the plot with points indicating where an experiment ran. This is only useful when creating a heatmap or contour plot with interpolation so that you can see which points were actually on the original path. Note: if any learner crashes occurred within the path, this will become TRUE. Not used with partial dependence. Default is FALSE.

```r
show.interpolated (logical(1))
```

If TRUE, will overlay the plot with points indicating where interpolation ran. This is only useful when creating a heatmap or contour plot with interpolation so that you can see which points were interpolated. Not used with partial dependence. Default is FALSE.

```r
nested.agg (function)
```

The function used to aggregate nested cross validation runs when plotting 2 hyperparameters. This is also used for nested aggregation in partial dependence. Default is mean.

```r
partial.dep.learn (Learner | character(1))
```

The regression learner used to learn partial dependence. Must be specified if “partial.dep” is set to TRUE in `generateHyperParsEffectData`. Accepts either a Learner object or the learner as a string for learning partial dependence. Default is NULL.

**Value**

`ggplot2` plot object.

**Note**

Any NAs incurred from learning algorithm crashes will be indicated in the plot (except in the case of partial dependence) and the NA values will be replaced with the column min/max depending on the optimal values for the respective measure. Execution time will be replaced with the max. Interpolation by its nature will result in predicted values for the performance measure. Use interpolation with caution. If “partial.dep” is set to TRUE in `generateHyperParsEffectData`, only partial dependence will be plotted.

Since a `ggplot2` plot object is returned, the user can change the axis labels and other aspects of the plot using the appropriate `ggplot2` syntax.
plotLearnerPrediction

Visualizes a learning algorithm on a 1D or 2D data set.

Description

Trains the model for 1 or 2 selected features, then displays it via `ggplot2::ggplot`. Good for teaching or exploring models.

For classification and clustering, only 2D plots are supported. The data points, the classification and potentially through color alpha blending the posterior probabilities are shown.

For regression, 1D and 2D plots are supported. 1D shows the data, the estimated mean and potentially the estimated standard error. 2D does not show estimated standard error, but only the estimated mean via background color.

The plot title displays the model id, its parameters, the training performance and the cross-validation performance.

Usage

```
plotLearnerPrediction(
  learner,       
  task,          
  features = NULL, 
  measures,      
  cv = 10L,       
  ...,           
  gridsize,      
  pointsize = 2, 
  prob.alpha = TRUE,  
  se.band = TRUE,   
  err.mark = "train",  
  bg.cols = c("darkblue", "green", "darkred"),  
  err.col = "white",  
  err.size = pointsize,  
  greyscale = FALSE,  
  pretty.names = TRUE  
)
```

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via `makeLearner`.

task (Task)
The task.
plotLearnerPrediction

features (character)
Selected features for model. By default the first 2 features are used.

measures (Measure | list of Measure)
Performance measure(s) to evaluate. Default is the default measure for the task, see here getDefaultMeasure.

cv (integer(1))
Do cross-validation and display in plot title? Number of folds. 0 means no CV. Default is 10.

... (any)
Parameters for learner.

gridsize (integer(1))
Grid resolution per axis for background predictions. Default is 500 for 1D and 100 for 2D.

pointsize (numeric(1))
Pointsize for ggplot2 ggplot2::geom_point for data points. Default is 2.

prob.alpha (logical(1))
For classification: Set alpha value of background to probability for predicted class? Allows visualization of “confidence” for prediction. If not, only a constant color is displayed in the background for the predicted label. Default is TRUE.

se.band (logical(1))
For regression in 1D: Show band for standard error estimation? Default is TRUE.

err.mark (character(1)): For classification: Either mark error of the model on the training data (“train”) or during cross-validation (“cv”) or not at all with “none”. Default is “train”.

bg.cols (character(3))
Background colors for classification and regression. Sorted from low, medium to high. Default is TRUE.

err.col (character(1))
For classification: Color of misclassified data points. Default is “white”

err.size (integer(1))
For classification: Size of misclassified data points. Default is pointsize.

greyscale (logical(1))
Should the plot be greyscale completely? Default is FALSE.

pretty.names (logical(1))
Whether to use the short name of the learner instead of its ID in labels. Defaults to TRUE.

Value
The ggplot2 object.
plotLearningCurve  

Plot learning curve data using ggplot2.

Description
Visualizes data size (percentage used for model) vs. performance measure(s).

Usage

plotLearningCurve(
  obj,  
  facet = "measure",  
  pretty.names = TRUE,  
  facet.wrap.nrow = NULL,  
  facet.wrap.ncol = NULL
)

Arguments

obj (LearningCurveData)
Result of generateLearningCurveData, with class LearningCurveData.

facet (character(1))
Selects “measure” or “learner” to be the facetting variable. The variable mapped to facet must have more than one unique value, otherwise it will be ignored. The variable not chosen is mapped to color if it has more than one unique value. The default is “measure”.

pretty.names (logical(1))
Whether to use the Measure name instead of the id in the plot. Default is TRUE.

facet.wrap.nrow, facet.wrap.ncol (integer)
Number of rows and columns for facetting. Default for both is NULL. In this case ggplot’s facet_wrap will choose the layout itself.

Value

ggplot2 plot object.

See Also

Other learning_curve: generateLearningCurveData()

Other plot: createSpatialResamplingPlots(), plotBMRBoxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCalibration(), plotCritDifferences(), plotPartialDependence(), plotROCCurves(), plotResiduals(), plotThreshVsPerf()
plotPartialDependence  Plot a partial dependence with ggplot2.

Description
Plot a partial dependence from generatePartialDependenceData using ggplot2.

Usage
plotPartialDependence(
  obj,
  geom = "line",
  facet = NULL,
  facet.wrap.nrow = NULL,
  facet.wrap.ncol = NULL,
  p = 1,
  data = NULL
)

Arguments

obj  PartialDependenceData
     Generated by generatePartialDependenceData.

geom  (character(1))
The type of geom to use to display the data. Can be “line” or “tile”. For tiling
at least two features must be used with interaction = TRUE in the call to generatePartialDependenceData. This may be used in conjunction with the facet argument if three features are specified in the call to generatePartialDependenceData. Default is “line”.

facet  (character(1))
The name of a feature to be used for facetting. This feature must have been an element of the features argument to generatePartialDependenceData and is only applicable when said argument had length greater than 1. The feature must be a factor or an integer. If generatePartialDependenceData is called with the interaction argument FALSE (the default) with argument features of length greater than one, then facet is ignored and each feature is plotted in its own facet. Default is NULL.

facet.wrap.nrow, facet.wrap.ncol  (integer)
Number of rows and columns for facetting. Default for both is NULL. In this case ggplot’s facet_wrap will choose the layout itself.

p  (numeric(1))
If individual = TRUE then sample allows the user to sample without replacement from the output to make the display more readable. Each row is sampled with probability p. Default is 1.
Create residual plots for prediction objects or benchmark results.

Description
Plots for model diagnostics. Provides scatterplots of true vs. predicted values and histograms of the model’s residuals.

Usage
plotResiduals(
  obj,
  type = "scatterplot",
  loess.smooth = TRUE,
  rug = TRUE,
  pretty.names = TRUE
)

Arguments

obj (Prediction | BenchmarkResult)
Input data.

type
Type of plot. Can be "scatterplot", the default. Or "hist", for a histogram, or in case of classification problems a barplot, displaying the residuals.

loess.smooth (logical(1))
Should a loess smoother be added to the plot? Defaults to TRUE. Only applicable for regression tasks and if type is set to scatterplot.
Plots a ROC curve from predictions.

Usage

plotROCCurves(
    obj,  # (ThreshVsPerfData)
    measures,  # ([list(2)] of Measure)
    diagonal = TRUE,  # Whether to plot a dashed diagonal line. Default is TRUE.
    pretty.names = TRUE,  # Whether to use the Measure name instead of the id in the plot. Default is TRUE.
    facet.learner = FALSE
)

Arguments

obj (ThreshVsPerfData)
Result of generateThreshVsPerfData.

measures ([list(2)] of Measure)
Default is the first 2 measures passed to generateThreshVsPerfData.

diagonal (logical(1))

pretty.names (logical(1))

Value

ggplot2 plot object.

See Also

Other plot: createSpatialResamplingPlots(), plotBMRBoxplots(), plotBMRtrendsAsBarChart(), plotBenchmarkResults(), plotCalibration(), plotCritDifferences(), plotLearningCurve(), plotPartialDependence(), plotROCCurves(), plotThresholdVsPerf()
plotThreshVsPerf

facet.learner  (logical(1))
Weather to use facetting or different colors to compare multiple learners. Default is FALSE.

Value

ggplot2 plot object.

See Also

Other plot: createSpatialResamplingPlots(), plotBMRRboxplots(), plotBMRRanksAsBarChart(),
plotBMRSsummary(), plotCalibration(), plotCritDifferences(), plotLearningCurve(), plotPartialDependence(),
plotResiduals(), plotThreshVsPerf()

Other thresh_vs_perf: generateThreshVsPerfData(), plotThreshVsPerf()

Examples

```r
lrn = makeLearner("classif.rpart", predict.type = "prob")
fit = train(lrn, sonar.task)
pred = predict(fit, task = sonar.task)
roc = generateThreshVsPerfData(pred, list(fpr, tpr))
plotROCCurves(roc)

r = bootstrapB632plus(lrn, sonar.task, iters = 3)
roc_r = generateThreshVsPerfData(r, list(fpr, tpr), aggregate = FALSE)
plotROCCurves(roc_r)

r2 = crossval(lrn, sonar.task, iters = 3)
roc_l = generateThreshVsPerfData(list(boot = r, cv = r2), list(fpr, tpr), aggregate = FALSE)
plotROCCurves(roc_l)
```

Description

Plots threshold vs. performance(s) data that has been generated with generateThreshVsPerfData.

Usage

```r
plotThreshVsPerf(
  obj,
  measures = obj$measures,
  facet = "measure",
  mark.th = NA_real_,
  ...)
Arguments

obj (ThreshVsPerfData)
Result of generateThreshVsPerfData.

measures (Measure | list of Measure)
Performance measure(s) to plot. Must be a subset of those used in generateThreshVsPerfData. Default is all the measures stored in obj generated by generateThreshVsPerfData.

facet (character(1))
Selects “measure” or “learner” to be the facetting variable. The variable mapped to facet must have more than one unique value, otherwise it will be ignored. The variable not chosen is mapped to color if it has more than one unique value. The default is “measure”.

mark.th (numeric(1))
Mark given threshold with vertical line? Default is NA which means not to do it.

pretty.names (logical(1))
Whether to use the Measure name instead of the id in the plot. Default is TRUE.

facet.wrap.nrow, facet.wrap.ncol (integer)
Number of rows and columns for facetting. Default for both is NULL. In this case ggplot’s facet_wrap will choose the layout itself.

Value
ggplot2 plot object.

See Also

Other plot: createSpatialResamplingPlots(), plotBMRBoxplots(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCalibration(), plotCritDifferences(), plotLearningCurve(), plotPartialDependence(), plotROCCurves(), plotResiduals()

Other thresh_vs_perf: generateThreshVsPerfData(), plotROCCurves()

Examples

lrn = makeLearner("classif.rpart", predict.type = "prob")
mod = train(lrn, sonar.task)
pred = predict(mod, sonar.task)
pvs = generateThreshVsPerfData(pred, list(acc, setAggregation(acc, train.mean)))
plotThreshVsPerf(pvs)
plotTuneMultiCritResult

Plots multi-criteria results after tuning using ggplot2.

Description

Visualizes the pareto front and possibly the dominated points.

Usage

plotTuneMultiCritResult(
  res,
  path = TRUE,
  col = NULL,
  shape = NULL,
  pointsize = 2,
  pretty.names = TRUE
)

Arguments

res (TuneMultiCritResult)
Result of tuneParamsMultiCrit.

path (logical(1))
Visualize all evaluated points (or only the non-dominated pareto front)? For the full path, the size of the points on the front is slightly increased. Default is TRUE.

col (character(1))
Which column of res$opt.path should be mapped to ggplot2 color? Default is NULL, which means none.

shape (character(1))
Which column of res$opt.path should be mapped to ggplot2 shape? Default is NULL, which means none.

pointsize (numeric(1))
Point size for ggplot2 ggplot2::geom_point for data points. Default is 2.

pretty.names (logical(1))
Whether to use the ID of the measures instead of their name in labels. Defaults to TRUE.

Value

ggplot2 plot object.

See Also

Other tune_multicrit: TuneMultiCritControl, tuneParamsMultiCrit()
predict.WrappedModel  

**Examples**

```r
# see tuneParamsMultiCrit
```

**Description**

Predict the target variable of new data using a fitted model. What is stored exactly in the (Prediction) object depends on the predict.type setting of the Learner. If predict.type was set to “prob” probability thresholding can be done calling the setThreshold function on the prediction object.

The row names of the input task or newdata are preserved in the output.

**Usage**

```r
## S3 method for class 'WrappedModel'
predict(object, task, newdata, subset = NULL, ...)
```

**Arguments**

- `object` (WrappedModel)
  - Wrapped model, result of train.
- `task` (Task)
  - The task. If this is passed, data from this task is predicted.
- `newdata` (data.frame)
  - New observations which should be predicted. Pass this alternatively instead of task.
- `subset` (integer | logical | NULL)
  - Selected cases. Either a logical or an index vector. By default NULL if all observations are used.
- `...` (any)
  - Currently ignored.

**Value**

(Prediction).

**See Also**

Other predict: asROCRPrediction(), getPredictionProbabilities(), getPredictionResponse(), getPredictionTaskDesc(), setPredictThreshold(), setPredictType()
Examples

# train and predict
train.set = seq(1, 150, 2)
test.set = seq(2, 150, 2)
model = train("classif.lda", iris.task, subset = train.set)
p = predict(model, newdata = iris, subset = test.set)
print(p)
predict(model, task = iris.task, subset = test.set)

# predict now probabilities instead of class labels
lrn = makeLearner("classif.lda", predict.type = "prob")
model = train(lrn, iris.task, subset = train.set)
p = predict(model, task = iris.task, subset = test.set)
print(p)
getPredictionProbabilities(p)

predictLearner

Predict new data with an R learner.

Description

Mainly for internal use. Predict new data with a fitted model. You have to implement this method if you want to add another learner to this package.

Usage

predictLearner(.learner, .model, .newdata, ...)

Arguments

.learner (RLearner)
Wrapped learner.

.model (WrappedModel)
Model produced by training.

.newdata (data.frame)
New data to predict. Does not include target column.

...(any)
Additional parameters, which need to be passed to the underlying predict function.

Details

Your implementation must adhere to the following: Predictions for the observations in .newdata must be made based on the fitted model (.model$learner.model). All parameters in ... must be passed to the underlying predict function.
reduceBatchmarkResults

Reduce results of a batch-distributed benchmark.

Description

This creates a BenchmarkResult from a batchtools::ExperimentRegistry. To setup the benchmark have a look at batchmark.

Usage

reduceBatchmarkResults(
  ids = NULL,
  keep.pred = TRUE,
  keep.extract = FALSE,
  show.info = getMlrOption("show.info"),
  reg = batchtools::getDefaultRegistry()
)

Arguments

ids (data.frame or integer)
A base::data.frame (or data.table::data.table) with a column named “job.id”. Alternatively, you may also pass a vector of integerish job ids. If not set, defaults to all successfully terminated jobs (return value of batchtools::findDone).

keep.pred (logical(1))
Keep the prediction data in the pred slot of the result object. If you do many experiments (on larger data sets) these objects might unnecessarily increase object size / mem usage, if you do not really need them. The default is set to TRUE.
```r
keep.extract (logical(1))
Keep the extract slot of the result object. When creating a lot of benchmark results with extensive tuning, the resulting R objects can become very large in size. That is why the tuning results stored in the extract slot are removed by default (keep.extract = FALSE). Note that when keep.extract = FALSE you will not be able to conduct analysis in the tuning results.

show.info (logical(1))
Print verbose output on console? Default is set via configureMlr.

reg (batchtools::ExperimentRegistry)
Registry, created by batchtools::makeExperimentRegistry. If not explicitly passed, uses the last created registry.

Value

(BenchmarkResult).

See Also

Other benchmark: BenchmarkResult, batchmark(), benchmark(), convertBMRToRankMatrix(), friedmanPostHocTestBMR(), friedmanTestBMR(), generateCritDifferencesData(), getBMRAggrPerformances(), getBMRFeatSelResults(), getBMRFilteredFeatures(), getBMRLearnerIds(), getBMRLearnerShortNames(), getBMRLearners(), getBMRMeasureIds(), getBMRMeasures(), getBMRModels(), getBMRPerformances(), getBMRPredictions(), getBMRTaskDescs(), getBMRTaskIds(), getBMRTuneResults(), plotBMRRanksAsBarChart(), plotBMRSummary(), plotCritDifferences()
```

**Description**

This function accepts a data frame or a task and an extractFDAFeatDesc (a FDA feature extraction description) as returned by extractFDAFeatures to extract features from previously unseen data.

**Usage**

```r
reextractFDAFeatures(obj, desc, ...)
```

**Arguments**

- **obj** 
  (Task | data.frame) 
  Task or data.frame to extract functional features from. Must contain functional features as matrix columns.

- **desc** 
  (extractFDAFeatDesc) 
  FDAFeature extraction description as returned by extractFDAFeatures

- **...** 
  (any) 
  Further args passed on to methods.
reimpute

Value
data.frame or Task containing the extracted Features

Description
This function accepts a data frame or a task and an imputation description as returned by impute to perform the following actions:

1. Restore dropped columns, setting them to NA
2. Add dummy variables for columns as specified in impute
3. Optionally check factors for new levels to treat them as NAs
4. Reorder factor levels to ensure identical integer representation as before
5. Impute missing values using previously collected data

Usage
reimpute(obj, desc)

Arguments

obj (data.frame | Task)
Input data.

desc (ImputationDesc)
Imputation description as returned by impute.

Value
Imputated data.frame or task with imputed data.

See Also
Other impute: imputations, impute(), makeImputeMethod(), makeImputeWrapper()
removeConstantFeatures

Remove constant features from a data set.

Description

Constant features can lead to errors in some models and obviously provide no information in the training set that can be learned from. With the argument “perc”, there is a possibility to also remove features for which less than “perc” percent of the observations differ from the mode value.

Usage

removeConstantFeatures(
  obj,               # (data.frame | Task) Input data.
  perc = 0,         # (numeric(1)) The percentage of a feature values in [0, 1) that must differ from the mode value. Default is 0, which means only constant features with exactly one observed level are removed.
  dont.rm = character(0L),  # (character) Names of the columns which must not be deleted. Default is no columns.
  na.ignore = FALSE,  # (logical(1)) Should NAs be ignored in the percentage calculation? (Or should they be treated as a single, extra level in the percentage calculation?) Note that if the feature has only missing values, it is always removed. Default is FALSE.
  tol = .Machine$double.eps^0.5,  # (numeric(1)) Numerical tolerance to treat two numbers as equal. Variables stored as double will get rounded accordingly before computing the mode. Default is sqrt(.Machine$double.eps).
  show.info = getMlrOption("show.info")  # (logical(1)) Print verbose output on console? Default is set via configureMlr.
)

Arguments

- **obj** *(data.frame | Task)*: Input data.
- **perc** *(numeric(1))*: The percentage of a feature values in [0, 1) that must differ from the mode value. Default is 0, which means only constant features with exactly one observed level are removed.
- **dont.rm** *(character)*: Names of the columns which must not be deleted. Default is no columns.
- **na.ignore** *(logical(1))*: Should NAs be ignored in the percentage calculation? (Or should they be treated as a single, extra level in the percentage calculation?) Note that if the feature has only missing values, it is always removed. Default is FALSE.
- **tol** *(numeric(1))*: Numerical tolerance to treat two numbers as equal. Variables stored as double will get rounded accordingly before computing the mode. Default is sqrt(.Machine$double.eps).
- **show.info** *(logical(1))*: Print verbose output on console? Default is set via configureMlr.

Value

- **data.frame | Task**: Same type as obj.
removeHyperPars

See Also

Other eda_and_preprocess: capLargeValues(), createDummyFeatures(), dropFeatures(), mergeSmallFactorLevels(), normalizeFeatures(), summarizeColumns(), summarizeLevels()

---

removeHyperPars  Remove hyperparameters settings of a learner.

Description

Remove settings (previously set through mlr) for some parameters. Which means that the default behavior for that param will now be used.

Usage

removeHyperPars(learner, ids = character(0L))

Arguments

learner  (Learner | character(1))
  The learner. If you pass a string the learner will be created via makeLearner.

ids  (character)
  Parameter names to remove settings for. Default is character(0L).

Value

Learner.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), makeLearner(), setHyperPars(), setId(), setLearnerId(), setPredictThreshold(), setPredictType()
resample

Fit models according to a resampling strategy.

Description

The function resample fits a model specified by Learner on a Task and calculates predictions and performance measures for all training and all test sets specified by a either a resampling description (ResampleDesc) or resampling instance (ResampleInstance).

You are able to return all fitted models (parameter models) or extract specific parts of the models (parameter extract) as returning all of them completely might be memory intensive.

The remaining functions on this page are convenience wrappers for the various existing resampling strategies. Note that if you need to work with precomputed training and test splits (i.e., resampling instances), you have to stick with resample.

Usage

resample(
  learner,
  task,
  resampling,
  measures,
  weights = NULL,
  models = FALSE,
  extract,
  keep.pred = TRUE,
  ..., 
  show.info = getMlrOption("show.info")
)

crossval(
  learner,
  task,
  iters = 10L,
  stratify = FALSE,
  measures,
  models = FALSE,
  keep.pred = TRUE,
  ..., 
  show.info = getMlrOption("show.info")
)

crepcv(
  learner,
  task,
  folds = 10L,
  reps = 10L,
resample

stratify = FALSE,
measures,
models = FALSE,
keep.pred = TRUE,
...
show.info = getMlrOption("show.info")
)

holdout(
learner,
task,
split = 2/3,
stratify = FALSE,
measures,
models = FALSE,
keep.pred = TRUE,
...
show.info = getMlrOption("show.info")
)

subsample(
learner,
task,
iters = 30,
split = 2/3,
stratify = FALSE,
measures,
models = FALSE,
keep.pred = TRUE,
...
show.info = getMlrOption("show.info")
)

bootstrapOOB(
learner,
task,
iters = 30,
stratify = FALSE,
measures,
models = FALSE,
keep.pred = TRUE,
...
show.info = getMlrOption("show.info")
)

bootstrapB632(
learner,
task,
resample

```r
iters = 30,
stratify = FALSE,
measures,
models = FALSE,
keep.pred = TRUE,
...
show.info = getMlrOption("show.info")
)

bootstrapB632plus(
  learner,
  task,
  iters = 30,
  stratify = FALSE,
  measures,
  models = FALSE,
  keep.pred = TRUE,
  ...
  show.info = getMlrOption("show.info")
)

growingcv(
  learner,
  task,
  horizon = 1,
  initial.window = 0.5,
  skip = 0,
  measures,
  models = FALSE,
  keep.pred = TRUE,
  ...
  show.info = getMlrOption("show.info")
)

fixedcv(
  learner,
  task,
  horizon = 1L,
  initial.window = 0.5,
  skip = 0,
  measures,
  models = FALSE,
  keep.pred = TRUE,
  ...
  show.info = getMlrOption("show.info")
)
Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

task (Task)
The task.

resampling (ResampleDesc or ResampleInstance)
Resampling strategy. If a description is passed, it is instantiated automatically.

measures (Measure | list of Measure)
Performance measure(s) to evaluate. Default is the default measure for the task, see here getDefaultMeasure.

weights (numeric)
Optional, non-negative case weight vector to be used during fitting. If given, must be of same length as observations in task and in corresponding order. Overwrites weights specified in the task. By default NULL which means no weights are used unless specified in the task.

models (logical(1))
Should all fitted models be returned? Default is FALSE.

extract (function)
Function used to extract information from a fitted model during resampling. Is applied to every WrappedModel resulting from calls to train during resampling. Default is to extract nothing.

keep.pred (logical(1))
Keep the prediction data in the pred slot of the result object. If you do many experiments (on larger data sets) these objects might unnecessarily increase object size / mem usage, if you do not really need them. The default is set to TRUE.

... (any)
Further hyperparameters passed to learner.

show.info (logical(1))
Print verbose output on console? Default is set via configureMlr.

iters (integer(1))
See ResampleDesc.

stratify (logical(1))
See ResampleDesc.

folds (integer(1))
See ResampleDesc.

reps (integer(1))
See ResampleDesc.

split (numeric(1))
See ResampleDesc.

horizon (numeric(1))
See ResampleDesc.

initial.window (numeric(1))
See ResampleDesc.

skip (integer(1))
See ResampleDesc.
Value

(ResampleResult).

Note

If you would like to include results from the training data set, make sure to appropriately adjust the resampling strategy and the aggregation for the measure. See example code below.

See Also

Other resample: ResamplePrediction, ResampleResult, addRRMeasure(), getRRPredictionList(), getRRPredictions(), getRRTaskDescription(), getRRTaskDesc(), makeResampleDesc(), makeResampleInstance().

Examples

```r
task = makeClassifTask(data = iris, target = "Species")
rdesc = makeResampleDesc("CV", iters = 2)
r = resample(makeLearner("classif.qda"), task, rdesc)
print(r$aggr)
print(r$measures.test)
print(r$pred)

# include the training set performance as well
r = resample(makeLearner("classif.qda"), task, rdesc,
    measures = list(mmce, setAggregation(mmce, train.mean)))
print(r$aggr)
```

ResamplePrediction  Prediction from resampling.

Description

Contains predictions from resampling, returned (among other stuff) by function resample. Can basically be used in the same way as Prediction, its super class. The main differences are: (a) The internal data.frame (member data) contains an additional column iter, specifying the iteration of the resampling strategy, and additional columns set, specifying whether the prediction was from an observation in the “train” or “test” set. (b) The prediction time is a numeric vector, its length equals the number of iterations.

See Also

Other resample: ResampleResult, addRRMeasure(), getRRPredictionList(), getRRPredictions(), getRRTaskDescription(), getRRTaskDesc(), makeResampleDesc(), makeResampleInstance(), resample()
ResampleResult

Description

A container for resample results.

Details

Resample Result:
A resample result is created by resample and contains the following object members:

- **task.id** (character(1)): Name of the Task.
- **learner.id** (character(1)): Name of the Learner.
- **measures.test** (data.frame): Gives you access to performance measurements on the individual test sets. Rows correspond to sets in resampling iterations, columns to performance measures.
- **measures.train** (data.frame): Gives you access to performance measurements on the individual training sets. Rows correspond to sets in resampling iterations, columns to performance measures. Usually not available, only if specifically requested, see general description above.
- **aggr** (numeric): Named vector of aggregated performance values. Names are coded like this `<measure>.<aggregation>`.
- **err.msgs** (data.frame): Number of rows equals resampling iterations and columns are: `iter`, `train`, `predict`. Stores error messages generated during train or predict, if these were caught via `configureMlr`.
- **err.dumps** (list of list of dump.frames): List with length equal to number of resampling iterations. Contains lists of `dump.frames` objects that can be fed to `debugger()` to inspect error dumps generated on learner errors. One iteration can generate more than one error dump depending on which of training, prediction on training set, or prediction on test set, operations fail. Therefore the lists have named slots `$train`, `$predict.train`, or `$predict.test` if relevant. The error dumps are only saved when option `on.error.dump` is TRUE.
- **pred** (**ResamplePrediction**): Container for all predictions during resampling.
- **models** [list of **WrappedModel**]: List of fitted models or NULL.
- **extract** (list): List of extracted parts from fitted models or NULL.
- **runtime** (numeric(1)): Time in seconds it took to execute the resampling.

The print method of this object gives a short overview, including task and learner ids, aggregated measures and runtime for the resampling.

See Also

Other resample: **ResamplePrediction.addRRMeasure()**, **getRRPredictionList()**, **getRRPredictions()**, **getRRTaskDescription()**, **getRRTaskDesc()**, **makeResampleDesc()**, **makeResampleInstance()**, **resample()**

Other debug: **FailureModel**, **getPredictionDump()**, **getRRDump()**
Description

Wraps an already implemented learning method from R to make it accessible to mlr. Call this method in your constructor. You have to pass an id (name), the required package(s), a description object for all changeable parameters (you do not have to do this for the learner to work, but it is strongly recommended), and use property tags to define features of the learner.

For a general overview on how to integrate a learning algorithm into mlr’s system, please read the section in the online tutorial: https://mlr.mlr-org.com/articles/tutorial/create_learner.html

To see all possible properties of a learner, go to: LearnerProperties.

Usage

makeRLearner()

makeRLearnerClassif(
  cl, package, par.set, par.vals = list(), properties = character(0L), name = cl, short.name = cl, note = "", class.weights.param = NULL, callees = character(0L)
)

makeRLearnerMultilabel(
  cl, package, par.set, par.vals = list(), properties = character(0L), name = cl, short.name = cl, note = "", callees = character(0L)
)

makeRLearnerRegr(
  cl, package,
par.set,
par.vals = list(),
properties = character(0L),
name = cl,
short.name = cl,
note = "",
callees = character(0L)
)

makeRLearnerSurv(
  cl,
  package,
  par.set,
  par.vals = list(),
  properties = character(0L),
  name = cl,
  short.name = cl,
  note = "",
callees = character(0L)
)

makeRLearnerCluster(
  cl,
  package,
  par.set,
  par.vals = list(),
  properties = character(0L),
  name = cl,
  short.name = cl,
  note = "",
callees = character(0L)
)

makeRLearnerCostSens(
  cl,
  package,
  par.set,
  par.vals = list(),
  properties = character(0L),
  name = cl,
  short.name = cl,
  note = "",
callees = character(0L)
)

Arguments

cl (character(1))
Class of learner. By convention, all classification learners start with “classif.”
all regression learners with “regr.” all survival learners start with “surv.” all clustering learners with “cluster.” and all multilabel classification learners start with “multilabel.”. A list of all integrated learners is available on the learners help page.

package (character)
Package(s) to load for the implementation of the learner.

par.set (ParamHelpers::ParamSet)
Parameter set of (hyper)parameters and their constraints. Dependent parameters with a requires field must use quote and not expression to define it.

par.vals (list)
Always set hyperparameters to these values when the object is constructed. Useful when default values are missing in the underlying function. The values can later be overwritten when the user sets hyperparameters. Default is empty list.

properties (character)
Set of learner properties. See above. Default is character(0).

name (character(1))
Meaningful name for learner. Default is id.

short.name (character(1))
Short name for learner. Should only be a few characters so it can be used in plots and tables. Default is id.

note (character(1))
Additional notes regarding the learner and its integration in mlr. Default is “”.

class.weights.param (character(1))
Name of the parameter, which can be used for providing class weights.

callees (character)
Character vector naming all functions of the learner’s package being called which have a relevant R help page. Default is character(0).

Value

(RLearner). The specific subclass is one of RLearnerClassif, RLearnerCluster, RLearnerMultilabel, RLearnerRegr, RLearnerSurv.

selectFeatures Feature selection by wrapper approach.

Description

Optimizes the features for a classification or regression problem by choosing a variable selection wrapper approach. Allows for different optimization methods, such as forward search or a genetic algorithm. You can select such an algorithm (and its settings) by passing a corresponding control object. For a complete list of implemented algorithms look at the subclasses of ([FeatSelControl]).

All algorithms operate on a 0-1-bit encoding of candidate solutions. Per default a single bit corresponds to a single feature, but you are able to change this by using the arguments ‘bit.names’ and ‘bits.to.features’. Thus allowing you to switch on whole groups of features with a single bit.
selectFeatures

Usage

selectFeatures(
  learner,
  task,
  resampling,
  measures,
  bit.names,
  bits.to.features,
  control,
  show.info = getMlrOption("show.info")
)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

task (Task)
The task.

resampling ([ResampleInstance] | [ResampleDesc])
Resampling strategy for feature selection. If you pass a description, it is instantiated once at the beginning by default, so all points are evaluated on the same training/test sets. If you want to change that behaviour, look at [FeatSelControl].

measures (list of Measure | Measure)
Performance measures to evaluate. The first measure, aggregated by the first aggregation function is optimized, others are simply evaluated. Default is the default measure for the task, see here getDefaultMeasure.

bit.names [character]
Names of bits encoding the solutions. Also defines the total number of bits in the encoding. Per default these are the feature names of the task. Has to be used together with ‘bits.to.features’.

bits.to.features [function(x, task)]
Function which transforms an integer-0-1 vector into a character vector of selected features. Per default a value of 1 in the ith bit selects the ith feature to be in the candidate solution. The vector ‘x’ will correspond to the ‘bit.names’ and has to be of the same length.

control [see [FeatSelControl]] Control object for search method. Also selects the optimization algorithm for feature selection.

show.info (logical(1))
Print verbose output on console? Default is set via configureMlr.

Value

([FeatSelResult]).

See Also

Other featsel: FeatSelControl, analyzeFeatSelResult(), getFeatSelResult(), makeFeatSelWrapper()
**setAggregation**

**Set aggregation function of measure.**

**Description**

Set how this measure will be aggregated after resampling. To see possible aggregation functions: `aggregations`.

**Usage**

```
setAggregation(measure, aggr)
```

**Arguments**

- `measure` *(Measure)*
  Performance measure.

- `aggr` *(Aggregation)*
  Aggregation function.

**Value**

*(Measure)* with changed aggregation behaviour.

**See Also**

Other performance: `ConfusionMatrix, calculateConfusionMatrix(), calculateROCMeasures(), estimateRelativeOverfitting(), makeCostMeasure(), makeCustomResampledMeasure(), makeMeasure(), measures, performance(), setMeasurePars()`
setHyperPars

Set the hyperparameters of a learner object.

Description

Set the hyperparameters of a learner object.

Usage

setHyperPars(learner, ..., par.vals = list())

Arguments

learner

(Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

...

(any)
Optional named (hyper)parameters. If you want to set specific hyperparameters for a learner during model creation, these should go here. You can get a list of available hyperparameters using getParamSet(<learner>). Alternatively hyperparameters can be given using the par.vals argument but ... should be preferred!

par.vals

(list)
Optional list of named (hyper)parameters. The arguments in ... take precedence over values in this list. We strongly encourage you to use ... for passing hyperparameters.

Value

Learner.

Note

If a named (hyper)parameter can’t be found for the given learner, the 3 closest (hyper)parameter names will be output in case the user mistyped.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setId(), setLearnerId(), setPredictThreshold(), setPredictType()
Examples

```r
c1 = makeLearner("classif.ksvm", sigma = 1)
c2 = setHyperPars(c1, sigma = 10, par.vals = list(C = 2))
print(c1)
# note the now set and altered hyperparameters:
print(c2)
```

setHyperPars2

*Only exported for internal use.*

**Description**

Only exported for internal use.

**Usage**

```r
setHyperPars2(learner, par.vals)
```

**Arguments**

- `learner` *(Learner)*
  
The learner.
- `par.vals` *(list)*
  
  List of named (hyper)parameter settings.

setId

*Set the id of a learner object.*

**Description**

Deprecated, use `setLearnerId` instead.

**Usage**

```r
setId(learner, id)
```

**Arguments**

- `learner` *(Learner | character(1))*
  
  The learner. If you pass a string the learner will be created via `makeLearner`.
- `id` *(character(1))*
  
  New id for learner.

**Value**

`Learner`.
setLearnerId

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(),
getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(),
helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setLearnerId(),
setPredictThreshold(), setPredictType()

---

setLearnerId

Set the ID of a learner object.

Description

Set the ID of the learner.

Usage

setLearnerId(learner, id)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

id (character(1))
New ID for learner.

Value

Learner.

See Also

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(),
getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(),
getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(),
helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setLearnerId(),
setPredictThreshold(), setPredictType()
**setMeasurePars**  
*Set parameters of performance measures*

**Description**  
Sets hyperparameters of measures.

**Usage**

```r
setMeasurePars(measure, ..., par.vals = list())
```

**Arguments**

- `measure` *(Measure)*  
  Performance measure.
- `...` *(any)*  
  Named (hyper)parameters with new settings. Alternatively these can be passed using the `par.vals` argument.
- `par.vals` *(list)*  
  Optional list of named (hyper)parameter settings. The arguments in `...` take precedence over values in this list.

**Value**

*Measure.*

**See Also**

Other performance: `ConfusionMatrix`, `calculateConfusionMatrix()`, `calculateROCMeasures()`, `estimateRelativeOverfitting()`, `makeCostMeasure()`, `makeCustomResampledMeasure()`, `makeMeasure()`, `measures`, `performance()`, `setAggregation()`

---

**setPredictThreshold**  
*Set the probability threshold the learner should use.*

**Description**

See `predict.threshold` in `makeLearner` and `setThreshold`.

For complex wrappers only the top-level `predict.type` is currently set.

**Usage**

```r
setPredictThreshold(learner, predict.threshold)
```
Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

predict.threshold (numeric)
Threshold to produce class labels. Has to be a named vector, where names correspond to class labels. Only for binary classification it can be a single numerical threshold for the positive class. See setThreshold for details on how it is applied. Default is NULL which means 0.5 / an equal threshold for each class.

Value

Learner.

See Also

Other predict: asROCRPrediction(), getPredictionProbabilities(), getPredictionResponse(), getPredictionTaskDesc(), predict.WrappedModel(), setPredictType()

Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(),getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(), setPredictType()

Description

Possible prediction types are: Classification: Labels or class probabilities (including labels). Regression: Numeric or response or standard errors (including numeric response). Survival: Linear predictor or survival probability.

For complex wrappers the predict type is usually also passed down the encapsulated learner in a recursive fashion.

Usage

setPredictType(learner, predict.type)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

predict.type (character(1))
Classification: “response” or “prob”. Regression: “response” or “se”. Survival: “response” (linear predictor) or “prob”. Clustering: “response” or “prob”. Default is “response”.

setPredictType  Set the type of predictions the learner should return.
setThreshold

Value
Learner.

See Also
Other predict: asROCRPrediction(), getPredictionProbabilities(), getPredictionResponse(), getPredictionTaskDesc(), predict.WrappedModel(), setPredictThreshold()
Other learner: LearnerProperties, getClassWeightParam(), getHyperPars(), getLearnerId(), getLearnerNote(), getLearnerPackages(), getLearnerParVals(), getLearnerParamSet(), getLearnerPredictType(), getLearnerShortName(), getLearnerType(), getParamSet(), helpLearnerParam(), helpLearner(), makeLearners(), makeLearner(), removeHyperPars(), setHyperPars(), setId(), setLearnerId(), setPredictThreshold()

---

setThreshold (Prediction) Set threshold of prediction object.

Description
Set threshold of prediction object for classification or multilabel classification. Creates corresponding discrete class response for the newly set threshold. For binary classification: The positive class is predicted if the probability value exceeds the threshold. For multiclass: Probabilities are divided by corresponding thresholds and the class with maximum resulting value is selected. The result of both are equivalent if in the multi-threshold case the values are greater than 0 and sum to 1. For multilabel classification: A label is predicted (with entry TRUE) if a probability matrix entry exceeds the threshold of the corresponding label.

Usage
setThreshold(pred, threshold)

Arguments
- pred (Prediction) Prediction object.
- threshold (numeric) Threshold to produce class labels. Has to be a named vector, where names correspond to class labels. Only for binary classification it can be a single numerical threshold for the positive class.

Value
(Prediction) with changed threshold and corresponding response.

See Also
predict.WrappedModel
Examples

```r
# create task and train learner (LDA)
task = makeClassifTask(data = iris, target = "Species")
lrn = makeLearner("classif.lda", predict.type = "prob")
mod = train(lrn, task)

# predict probabilities and compute performance
pred = predict(mod, newdata = iris)
performance(pred, measures = mmce)
head(as.data.frame(pred))

# adjust threshold and predict probabilities again
threshold = c(setosa = 0.4, versicolor = 0.3, virginica = 0.3)
pred = setThreshold(pred, threshold = threshold)
performance(pred, measures = mmce)
head(as.data.frame(pred))
```

simplifyMeasureNames  
Simplify measure names.

Description

Clips aggregation names from character vector. E.g: 'mmce.test.mean' becomes 'mmce'. Elements that don’t contain a measure name are ignored and returned unchanged.

Usage

`simplifyMeasureNames(xs)`

Arguments

- `xs`  
  (character)  
  Character vector that (possibly) contains aggregated measure names.

Value

(character).
Smote

Description

In each iteration, samples one minority class element x1, then one of x1’s nearest neighbors: x2. Both points are now interpolated / convex-combined, resulting in a new virtual data point x3 for the minority class.

The method handles factor features, too. The gower distance is used for nearest neighbor calculation, see cluster::daisy. For interpolation, the new factor level for x3 is sampled from the two given levels of x1 and x2 per feature.

Usage

smote(task, rate, nn = 5L, standardize = TRUE, alt.logic = FALSE)

Arguments

- **task** (Task): The task.
- **rate** (numeric(1)): Factor to upsample the smaller class. Must be between 1 and Inf, where 1 means no oversampling and 2 would mean doubling the class size.
- **nn** (integer(1)): Number of nearest neighbors to consider. Default is 5.
- **standardize** (integer(1)): Standardize input variables before calculating the nearest neighbors for data sets with numeric input variables only. For mixed variables (numeric and factor) the gower distance is used and variables are standardized anyway. Default is TRUE.
- **alt.logic** (integer(1)): Use an alternative logic for selection of minority class observations. Instead of sampling a minority class element AND one of its nearest neighbors, each minority class element is taken multiple times (depending on rate) for the interpolation and only the corresponding nearest neighbor is sampled. Default is FALSE.

Value

Task.

References

See Also

Other imbalancy: `makeOverBaggingWrapper()`, `makeUndersampleWrapper()`, `oversample()`

---

**sonar.task**

*Sonar classification task.*

**Description**

Contains the task (`sonar.task`).

**References**

See `mlbench::Sonar`.

---

**spam.task**

*Spam classification task.*

**Description**

Contains the task (`spam.task`).

**References**

See `kernlab::spam`.

---

**spatial.task**

*J. Muenchow’s Ecuador landslide data set*

**Description**

Data set created by Jannes Muenchow, University of Erlangen-Nuremberg, Germany. These data should be cited as Muenchow et al. (2012) (see reference below). This publication also contains additional information on data collection and the geomorphology of the area. The data set provided here is (a subset of) the one from the 'natural' part of the RBSF area and corresponds to landslide distribution in the year 2000.

**Format**

a `data.frame` with point samples of landslide and non-landslide locations in a study area in the Andes of southern Ecuador.

**References**


subsetTask  Subset data in task.

Description
See title.

Usage
subsetTask(task, subset = NULL, features)

Arguments

- **task** *(Task)*
  The task.

- **subset** *(integer | logical | NULL)*
  Selected cases. Either a logical or an index vector. By default NULL if all observations are used.

- **features** *(character | integer | logical)*
  Vector of selected inputs. You can either pass a character vector with the feature names, a vector of indices, or a logical vector.
  In case of an index vector each element denotes the position of the feature name returned by `getTaskFeatureNames`.
  Note that the target feature is always included in the resulting task, you should not pass it here. Default is to use all features.

Value

*(Task)*. Task with subsetted data.

See Also

Other task: `getTaskClassLevels()`, `getTaskCosts()`, `getTaskData()`, `getTaskDesc()`, `getTaskFeatureNames()`, `getTaskFormula()`, `getTaskId()`, `getTaskNFeats()`, `getTaskSize()`, `getTaskTargetNames()`, `getTaskTargets()`, `getTaskType()`

Examples

```r
task = makeClassifTask(data = iris, target = "Species")
subsetTask(task, subset = 1:100)
```
summarizeColumns(Summarize columns of data.frame or task.

Description

Summarizes a data.frame, somewhat differently than the normal summary function of R. The function is mainly useful as a basic EDA tool on data.frames before they are converted to tasks, but can be used on tasks as well.

Columns can be of type numeric, integer, logical, factor, or character. Characters and logicals will be treated as factors.

Usage

summarizeColumns(obj)

Arguments

obj (data.frame | Task) Input data.

Value

(data.frame). With columns:

name Name of column.
type Data type of column.
na Number of NAs in column.
disp Measure of dispersion, for numerics and integers sd is used, for categorical columns the qualitative variation.
mean Mean value of column, NA for categorical columns.
median Median value of column, NA for categorical columns.
mad MAD of column, NA for categorical columns.
min Minimal value of column, for categorical columns the size of the smallest category.
max Maximal value of column, for categorical columns the size of the largest category.
nlevs For categorical columns, the number of factor levels, NA else.

See Also

Other eda_and_preprocess: capLargeValues(), createDummyFeatures(), dropFeatures(), mergeSmallFactorLevels(), normalizeFeatures(), removeConstantFeatures(), summarizeLevels()

Examples

summarizeColumns(iris)
summarizeLevels

**summarizeLevels**

_Summarizes factors of a data.frame by tabling them._

**Description**

Characters and logicals will be treated as factors.

**Usage**

```r
summarizeLevels(obj, cols = NULL)
```

**Arguments**

- **obj**
  - (data.frame | Task)
  - Input data.

- **cols**
  - (character)
  - Restrict result to columns in cols. Default is all factor, character and logical columns of obj.

**Value**

(list). Named list of tables.

**See Also**

Other eda_and_preprocess: `capLargeValues()`, `createDummyFeatures()`, `dropFeatures()`, `mergeSmallFactorLevels()`, `normalizeFeatures()`, `removeConstantFeatures()`, `summarizeColumns()`

**Examples**

```r
summarizeLevels(iris)
```

---

**Task**

Create a classification, regression, survival, cluster, cost-sensitive classification or multilabel task.

**Description**

The task encapsulates the data and specifies - through its subclasses - the type of the task. It also contains a description object detailing further aspects of the data.

Useful operators are:

- `getTaskFormula`,
- `getTaskFeatureNames`,
- `getTaskData`,
- `getTaskDescription`,
- `getTaskTarget`. 
- `getTaskTargets`, and
- `subsetTask`.

Object members:

- **env** *(environment)* Environment where data for the task are stored. Use `getTaskData` in order to access it.
- **weights** *(numeric)* See argument. NULL if not present.
- **blocking** *(factor)* See argument. NULL if not present.
- **task.desc** *(TaskDesc)* Encapsulates further information about the task.

Functional data can be added to a task via matrix columns. For more information refer to `make-FunctionalData`.

**Arguments**

- **id** *(character(1))*
  Id string for object. Default is the name of the R variable passed to `data`.

- **data** *(data.frame)*
  A data frame containing the features and target variable(s).

- **target** *(character(1) | character(2) | character(n.classes))*
  Name(s) of the target variable(s). For survival analysis these are the names of the survival time and event columns, so it has length 2. For multilabel classification it contains the names of the logical columns that encode whether a label is present or not and its length corresponds to the number of classes.

- **costs** *(data.frame)*
  A numeric matrix or data frame containing the costs of misclassification. We assume the general case of observation specific costs. This means we have n rows, corresponding to the observations, in the same order as `data`. The columns correspond to classes and their names are the class labels (if unnamed we use y1 to yk as labels). Each entry (i,j) of the matrix specifies the cost of predicting class j for observation i.

- **weights** *(numeric)*
  Optional, non-negative case weight vector to be used during fitting. Cannot be set for cost-sensitive learning. Default is NULL which means no (= equal) weights.

- **blocking** *(factor)*
  An optional factor of the same length as the number of observations. Observations with the same blocking level “belong together”. Specifically, they are either put all in the training or the test set during a resampling iteration. Default is NULL which means no blocking.

- **positive** *(character(1))*
  Positive class for binary classification (otherwise ignored and set to NA). Default is the first factor level of the target attribute.

- **fixup.data** *(character(1))*
  Should some basic cleaning up of data be performed? Currently this means
removing empty factor levels for the columns. Possible choices are: “no” = Don’t do it. “warn” = Do it but warn about it. “quiet” = Do it but keep silent. Default is “warn”.

check.data (logical(1))
Should sanity of data be checked initially at task creation? You should have good reasons to turn this off (one might be speed). Default is TRUE.

coordinates (data.frame)
Coordinates of a spatial data set that will be used for spatial partitioning of the data in a spatial cross-validation resampling setting. Coordinates have to be numeric values. Provided data.frame needs to have the same number of rows as data and consist of at least two dimensions.

Value
Task.

See Also
ClassifTask ClusterTask CostSensTask MultilabelTask RegrTask SurvTask

Examples
if (requireNamespace("mlbench")) {
  library(mlbench)
  data(BostonHousing)
  data(Ionosphere)

  makeClassifTask(data = iris, target = "Species")
  makeRegrTask(data = BostonHousing, target = "medv")
  # an example of a classification task with more than those standard arguments:
  blocking = factor(c(rep(1, 51), rep(2, 300)))
  makeClassifTask(id = "myIonosphere", data = Ionosphere, target = "Class",
                  positive = "good", blocking = blocking)
  makeClusterTask(data = iris[, -5L])
}

TaskDesc

Description object for task.

Description

Description object for task, encapsulates basic properties of the task without having to store the complete data set.
Details

Object members:

- **id** (character(1)) Id string of task.
- **type** (character(1)) Type of task, “classif” for classification, “regr” for regression, “surv” for survival and “cluster” for cluster analysis, “costsens” for cost-sensitive classification, and “multilabel” for multilabel classification.
- **target** (character(0) | character(1) | character(2) | character(n.classes)) Name(s) of the target variable(s). For “surv” these are the names of the survival time and event columns, so it has length 2. For “costsens” it has length 0, as there is no target column, but a cost matrix instead. For “multilabel” these are the names of logical columns that indicate whether a class label is present and the number of target variables corresponds to the number of classes.
- **size** (integer(1)) Number of cases in data set.
- **n.feat** (integer(2)) Number of features, named vector with entries: “numerics”, “factors”, “ordered”, “functionals”.
- **has.missings** (logical(1)) Are missing values present?
- **has.weights** (logical(1)) Are weights specified for each observation?
- **has.blocking** (logical(1)) Is a blocking factor for cases available in the task?
- **class.levels** (character) All possible classes. Only present for “classif”, “costsens”, and “multilabel”.
- **positive** (character(1)) Positive class label for binary classification. Only present for “classif”, NA for multiclass.
- **negative** (character(1)) Negative class label for binary classification. Only present for “classif”, NA for multiclass.

---

**train**

*Train a learning algorithm.*

---

Description

Given a Task, creates a model for the learning machine which can be used for predictions on new data.

Usage

```r
train(learner, task, subset = NULL, weights = NULL)
```

Arguments

- **learner** (Learner | character(1))
The learner. If you pass a string the learner will be created via `makeLearner`.
- **task** (Task)
The task.
trainLearner

**Description**

Mainly for internal use. Trains a wrapped learner on a given training set. You have to implement this method if you want to add another learner to this package.

**Usage**

`trainLearner(.learner, .task, .subset, .weights = NULL, ...)`

**subset**

(integer | logical | NULL)
Selected cases. Either a logical or an index vector. By default NULL if all observations are used.

**weights**

(numeric)
Optional, non-negative case weight vector to be used during fitting. If given, must be of same length as subset and in corresponding order. By default NULL which means no weights are used unless specified in the task (Task). Weights from the task will be overwritten.

**Value**

(WrappedModel).

**See Also**

predict.WrappedModel

**Examples**

```r
training.set = sample(seq_len(nrow(iris)), nrow(iris) / 2)
## use linear discriminant analysis to classify iris data
task = makeClassifTask(data = iris, target = "Species")
learner = makeLearner("classif.lda", method = "mle")
mod = train(learner, task, subset = training.set)
print(mod)

## use random forest to classify iris data
task = makeClassifTask(data = iris, target = "Species")
learner = makeLearner("classif.rpart", minsplit = 7, predict.type = "prob")
mod = train(learner, task, subset = training.set)
print(mod)
```
TuneControl

Arguments

- **.learner** (RLearner)
  Wrapped learner.
- **.task** (Task)
  Task to train learner on.
- **.subset** (integer)
  Subset of cases for training set, index the task with this. You probably want to use getTaskData for this purpose.
- **.weights** (numeric)
  Weights for each observation.
- **...**
  Additional (hyper)parameters, which need to be passed to the underlying train function.

Details

Your implementation must adhere to the following: The model must be fitted on the subset of .task given by .subset. All parameters in ... must be passed to the underlying training function.

Value

(any). Model of the underlying learner.

---

TuneControl  Control object for tuning

Description

General tune control object.

Arguments

- **same.resampling.instance** (logical(1))
  Should the same resampling instance be used for all evaluations to reduce variance? Default is TRUE.
- **impute.val** (numeric)
  If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or Inf instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.
TuneMultiCritControl

Create control structures for multi-criteria tuning.

Description

The following tuners are available:

makeTuneMultiCritControlGrid Grid search. All kinds of parameter types can be handled. You can either use their correct param type and resolution, or discretize them yourself by always using ParamHelpers::makeDiscreteParam in the par.set passed to tuneParams.
**makeTuneMultiCritControlRandom** Random search. All kinds of parameter types can be handled.

**makeTuneMultiCritControlNSGA2** Evolutionary method `mco::nsga2`. Can handle numeric(vector) and integer(vector) hyperparameters, but no dependencies. For integers the internally proposed numeric values are automatically rounded.

**makeTuneMultiCritControlMBO** Model-based/ Bayesian optimization. All kinds of parameter types can be handled.

### Usage

```r
makeTuneMultiCritControlGrid(
    same.resampling.instance = TRUE,
    resolution = 10L,
    log.fun = "default",
    final.dw.perc = NULL,
    budget = NULL
)

makeTuneMultiCritControlMBO(
    n.objectives = mbo.control$n.objectives,
    same.resampling.instance = TRUE,
    impute.val = NULL,
    learner = NULL,
    mbo.control = NULL,
    tune.threshold = FALSE,
    tune.threshold.args = list(),
    continue = FALSE,
    log.fun = "default",
    final.dw.perc = NULL,
    budget = NULL,
    mbo.design = NULL
)

makeTuneMultiCritControlNSGA2(
    same.resampling.instance = TRUE,
    impute.val = NULL,
    log.fun = "default",
    final.dw.perc = NULL,
    budget = NULL,
    ...
)

makeTuneMultiCritControlRandom(
    same.resampling.instance = TRUE,
    maxit = 100L,
    log.fun = "default",
    final.dw.perc = NULL,
    budget = NULL
)
Arguments

same.resampling.instance (logical(1))
Should the same resampling instance be used for all evaluations to reduce variance? Default is TRUE.

resolution (integer)
Resolution of the grid for each numeric/integer parameter in par.set. For vector parameters, it is the resolution per dimension. Either pass one resolution for all parameters, or a named vector. See ParamHelpers::generateGridDesign. Default is 10.

log.fun (function | character(1))
Function used for logging. If set to “default” (the default), the evaluated design points, the resulting performances, and the runtime will be reported. If set to “memory” the memory usage for each evaluation will also be displayed, with character(1) small increase in run time. Otherwise character(1) function with arguments learner, resampling, measures, par.set, control, opt.path, dob, x, y, remove.nas, stage and prev.stage is expected. The default displays the performance measures, the time needed for evaluating, the currently used memory and the max memory ever used before (the latter two both taken from gc). See the implementation for details.

final.dw.perc (boolean)
If a Learner wrapped by a makeDownsampleWrapper is used, you can define the value of dw.perc which is used to train the Learner with the final parameter setting found by the tuning. Default is NULL which will not change anything.

budget (integer(1))
Maximum budget for tuning. This value restricts the number of function evaluations. In case of makeTuneMultiCritControlGrid this number must be identical to the size of the grid. For makeTuneMultiCritControlRandom the budget equals the number of iterations (maxit) performed by the random search algorithm. In case of makeTuneMultiCritControlNSGA2 the budget corresponds to the product of the maximum number of generations (max(generations)) + 1 (for the initial population) and the size of the population (popsize). For makeTuneMultiCritControlMBO the budget equals the number of objective function evaluations, i.e. the number of MBO iterations + the size of the initial design. If not NULL, this will overwrite existing stopping conditions in mbo.control.

n.objectives (integer(1))
Number of objectives, i.e. number of Measures to optimize.

impute.val (numeric)
If something goes wrong during optimization (e.g. the learner crashes), this value is fed back to the tuner, so the tuning algorithm does not abort. It is not stored in the optimization path, an NA and a corresponding error message are logged instead. Note that this value is later multiplied by -1 for maximization measures internally, so you need to enter a larger positive value for maximization.
here as well. Default is the worst obtainable value of the performance measure you optimize for when you aggregate by mean value, or Inf instead. For multi-criteria optimization pass a vector of imputation values, one for each of your measures, in the same order as your measures.

**learner**

(Learner | NULL)

The surrogate learner: A regression learner to model performance landscape. For the default, NULL, mlrMBO will automatically create a suitable learner based on the rules described in mlrMBO::makeMBOlearner.

**mbo.control**

(mlrMBO::MBOControl | NULL)

Control object for model-based optimization tuning. For the default, NULL, the control object will be created with all the defaults as described in mlrMBO::makeMBOControl.

**tune.threshold**

(logical(1))

Should the threshold be tuned for the measure at hand, after each hyperparameter evaluation, via tuneThreshold? Only works for classification if the predict type is “prob”. Default is FALSE.

**tune.threshold.args**

(list)

Further arguments for threshold tuning that are passed down to tuneThreshold. Default is none.

**continue**

(logical(1))

Resume calculation from previous run using mlrMBO::mboContinue? Requires “save.file.path” to be set. Note that the ParamHelpers::OptPath in the mlrMBO::OptResult will only include the evaluations after the continuation. The complete OptPath will be found in the slot $mbo.result$opt.path.

**mbo.design**

(data.frame | NULL)

Initial design as data frame. If the parameters have corresponding trafo functions, the design must not be transformed before it is passed! For the default, NULL, a default design is created like described in mlrMBO::mbo.

**maxit**

(integer(1))

Number of iterations for random search. Default is 100.

**Value**

(TuneMultiCritControl). The specific subclass is one of TuneMultiCritControlGrid, TuneMultiCritControlRandom, TuneMultiCritControlNSGA2, TuneMultiCritControlMBO.

**See Also**

Other tune_multicrit: plotTuneMultiCritResult(), tuneParamsMultiCrit()
TuneMultiCritResult

Result of multi-criteria tuning.

Description

Container for results of hyperparameter tuning. Contains the obtained pareto set and front and the optimization path which lead there.

Object members:

learner (Learner) Learner that was optimized.
control (TuneControl) Control object from tuning.
x (list) List of lists of non-dominated hyperparameter settings in pareto set. Note that when you have trafo's on some of your params, x will always be on the TRANSFORMED scale so you directly use it.
y (matrix) Pareto front for x.
threshold Currently NULL.
opt.path (ParamHelpers::OptPath) Optimization path which lead to x. Note that when you have trafo's on some of your params, the opt.path always contains the UNTRANSFORMED values on the original scale. You can simply call trafoOptPath(opt.path) to transform them, or, as.data.frame{trafoOptPath(opt.path)}
ind (integer(n)) Indices of Pareto optimal params in opt.path.
measures [list of] Measure Performance measures.

tuneParams

Hyperparameter tuning.

Description

Optimizes the hyperparameters of a learner. Allows for different optimization methods, such as grid search, evolutionary strategies, iterated F-race, etc. You can select such an algorithm (and its settings) by passing a corresponding control object. For a complete list of implemented algorithms look at TuneControl.

Multi-criteria tuning can be done with tuneParamsMultiCrit.

Usage

tuneParams(
    learner,
    task,
    resampling,
    measures,
    par.set,
    control,
    show.info = getMlrOption("show.info"),
    resample.fun = resample
)
Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

task (Task)
The task.

resampling (ResampleInstance | ResampleDesc)
Resampling strategy to evaluate points in hyperparameter space. If you pass a
description, it is instantiated once at the beginning by default, so all points are
evaluated on the same training/test sets. If you want to change that behavior,
look at TuneControl.

measures (list of Measure | Measure)
Performance measures to evaluate. The first measure, aggregated by the first
aggregation function is optimized, others are simply evaluated. Default is the
default measure for the task, see here getDefaultMeasure.

par.set (ParamHelpers::ParamSet)
Collection of parameters and their constraints for optimization. Dependent pa-
rameters with a requires field must use quote and not expression to define it.

control (TuneControl)
Control object for search method. Also selects the optimization algorithm for
tuning.

show.info (logical(1))
Print verbose output on console? Default is set via configureMlr.

resample.fun (closure)
The function to use for resampling. Defaults to resample. If a user-given func-
tion is to be used instead, it should take the arguments “learner”, “task”, “re-
sampling”, “measures”, and “show.info”; see resample. Within this function, it
is easiest to call resample and possibly modify the result. However, it is pos-
sible to return a list with only the following essential slots: the “aggr” slot for
general tuning, additionally the “pred” slot if threshold tuning is performed (see
TuneControl), and the “err.msgs” and “err.dumps” slots for error reporting. This
parameter must be the default when mbo tuning is performed.

Value

(TuneResult).

Note

If you would like to include results from the training data set, make sure to appropriately adjust the
resampling strategy and the aggregation for the measure. See example code below.

See Also

generateHyperParsEffectData

Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(),
getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(),
Examples

```r
set.seed(123)
# a grid search for an SVM (with a tiny number of points...)
# note how easily we can optimize on a log-scale
ps = makeParamSet(
  makeNumericParam("C", lower = -12, upper = 12, trafo = function(x) 2^x),
  makeNumericParam("sigma", lower = -12, upper = 12, trafo = function(x) 2^x)
)
ctrl = makeTuneControlGrid(resolution = 2L)
rdesc = makeResampleDesc("CV", iters = 2L)
res = tuneParams("classif.ksvm", iris.task, rdesc, par.set = ps, control = ctrl)
print(res)
# access data for all evaluated points
df = as.data.frame(res$opt.path)
df1 = as.data.frame(res$opt.path, trafo = TRUE)
print(head(df[, -ncol(df)]))
print(head(df1[, -ncol(df1)]))
# access data for all evaluated points - alternative
df2 = generateHyperParsEffectData(res)
df3 = generateHyperParsEffectData(res, trafo = TRUE)
print(head(df2$data[, -ncol(df2$data)]))
print(head(df3$data[, -ncol(df3$data)]))
## Not run:
# we optimize the SVM over 3 kernels simultaneously
# note how we use dependent params (requires = ...) and iterated F-racing here
ps = makeParamSet(
  makeNumericParam("C", lower = -12, upper = 12, trafo = function(x) 2^x),
  makeDiscreteParam("kernel", values = c("vanilladot", "polydot", "rbfdot")),
  makeNumericParam("sigma", lower = -12, upper = 12, trafo = function(x) 2^x,
    requires = quote(kernel == "rbfdot")),
  makeIntegerParam("degree", lower = 2L, upper = 5L,
    requires = quote(kernel == "polydot"))
)
print(ps)
ctrl = makeTuneControlIrace(maxExperiments = 5, nbIterations = 1, minNbSurvival = 1)
rdesc = makeResampleDesc("Holdout")
res = tuneParams("classif.ksvm", iris.task, rdesc, par.set = ps, control = ctrl)
print(res)
df = as.data.frame(res$opt.path)
print(head(df[, -ncol(df)]))
# include the training set performance as well
rdesc = makeResampleDesc("Holdout", predict = "both")
res = tuneParams("classif.ksvm", iris.task, rdesc, par.set = ps,
  control = ctrl, measures = list(mmce, setAggregation(mmce, train.mean)))
print(res)
df2 = as.data.frame(res$opt.path)
print(head(df2[, -ncol(df2)]))
```
tuneParamsMultiCrit

Hyperparameter tuning for multiple measures at once.

Description

Optimizes the hyperparameters of a learner in a multi-criteria fashion. Allows for different optimization methods, such as grid search, evolutionary strategies, etc. You can select such an algorithm (and its settings) by passing a corresponding control object. For a complete list of implemented algorithms look at [TuneMultiCritControl].

Usage

tuneParamsMultiCrit(
  learner,
  task,
  resampling,
  measures,
  par.set,
  control,
  show.info = getMlrOption("show.info"),
  resample.fun = resample
)

Arguments

learner (Learner | character(1))
The learner. If you pass a string the learner will be created via makeLearner.

task (Task)
The task.

resampling ([ResampleInstance] | [ResampleDesc])
Resampling strategy to evaluate points in hyperparameter space. If you pass a description, it is instantiated once at the beginning by default, so all points are evaluated on the same training/test sets. If you want to change that behavior, look at [TuneMultiCritControl].

measures ([list of [Measure]])
Performance measures to optimize simultaneously.

par.set ([ParamHelpers::ParamSet])
Collection of parameters and their constraints for optimization. Dependent parameters with a ‘requires’ field must use ‘quote’ and not ‘expression’ to define it.

test (TuneMultiCritControl)
Control object for search method. Also selects the optimization algorithm for tuning.
show.info  (logical(1))
Print verbose output on console? Default is set via configureMlr.

resample.fun  ([closure])
The function to use for resampling. Defaults to [resample] and should take
the same arguments as, and return the same result type as, [resample].

Value
([TuneMultiCritResult]).

See Also
Other tune_multicrit: TuneMultiCritControl, plotTuneMultiCritResult()

Examples

```r
# multi-criteria optimization of (tpr, fpr) with NGSA-II
lrn = makeLearner("classif.ksvm")
rdesc = makeResampleDesc("Holdout")
ps = makeParamSet(
  makeNumericParam("C", lower = -12, upper = 12, trafo = function(x) 2^x),
  makeNumericParam("sigma", lower = -12, upper = 12, trafo = function(x) 2^x)
)
ctrl = makeTuneMultiCritControlNSGA2(popsize = 4L, generations = 1L)
res = tuneParamsMultiCrit(lrn, sonar.task, rdesc, par.set = ps,
  measures = list(tpr, fpr), control = ctrl)
plotTuneMultiCritResult(res, path = TRUE)
```

---

**TuneResult**

*Result of tuning.*

**Description**

Container for results of hyperparameter tuning. Contains the obtained point in search space, its
performance values and the optimization path which lead there.

Object members:

- **learner** *(Learner)* Learner that was optimized.
- **control** *(TuneControl)* Control object from tuning.
- **x** *(list)* Named list of hyperparameter values identified as optimal. Note that when you have trafos
  on some of your params, x will always be on the TRANSFORMED scale so you directly use it.
- **y** *(numeric)* Performance values for optimal x.
- **threshold** *(numeric)* Vector of finally found and used thresholds if tune.threshold was enabled
  in TuneControl, otherwise not present and hence NULL.
tuneThreshold

**opt.path (ParamHelpers::OptPath)** Optimization path which lead to x. Note that when you have trafos on some of your params, the opt.path always contains the UNTRANSFORMED values on the original scale. You can simply call trafoOptPath(opt.path) to transform them, or, as.data.frame(trafoOptPath(opt.path)). If mlr option on.error.dump is TRUE, OptPath will have a .dump object in its extra column which contains error dump traces from failed optimization evaluations. It can be accessed by getOptPathEl(opt.path)$extra$.dump.

---

**tuneThreshold**

Tune prediction threshold.

**Description**

Optimizes the threshold of predictions based on probabilities. Works for classification and multilabel tasks. Uses BBmisc::optimizeSubInts for normal binary class problems and GenSA::GenSA for multiclass and multilabel problems.

**Usage**

tuneThreshold(pred, measure, task, model, nsub = 20L, control = list())

**Arguments**

- **pred** (Prediction)
  - Prediction object.
- **measure** (Measure)
  - Performance measure to optimize. Default is the default measure for the task.
- **task** (Task)
  - Learning task. Rarely needed, only when required for the performance measure.
- **model** (WrappedModel)
  - Fitted model. Rarely needed, only when required for the performance measure.
- **nsub** (integer(1))
  - Passed to BBmisc::optimizeSubInts for 2class problems. Default is 20.
- **control** (list)
  - Control object for GenSA::GenSA when used. Default is empty list.

**Value**

(list). A named list with with the following components: th is the optimal threshold, perf the performance value.

**See Also**

Other tune: TuneControl, getNestedTuneResultsOptPathDf(), getNestedTuneResultsX(), getResamplingIndices(), getTuneResult(), makeModelMultiplexerParamSet(), makeModelMultiplexer(), makeTuneControlCMAES(), makeTuneControlDesign(), makeTuneControlGenSA(), makeTuneControlGrid(), makeTuneControlIrace(), makeTuneControlMBO(), makeTuneControlRandom(), makeTuneWrapper(), tuneParams()
wpbc.task

Wisconsin Prognostic Breast Cancer (WPBC) survival task.

Description

Contains the task (wpbc.task).

References

See TH.data::wpbc. Incomplete cases have been removed from the task.

yeast.task

Yeast multilabel classification task.

Description

Contains the task (yeast.task).

Source

https://archive.ics.uci.edu/ml/datasets/Yeast (In long instead of wide format)

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