# Package 'fastfurious'

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<b>Description</b> fast-furiuos gathers code (R, Matlab/Octave, Python), models and metamodels I needed in my Machine Learning Lab but I didn't found on the shelf.
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<pre>URL https://github.com/gtesei/fast-furious</pre>
<pre>BugReports https://github.com/gtesei/fast-furious/issues</pre>
VignetteBuilder knitr
<b>Suggests</b> knitr, lattice (>= 0.20), ggplot2 (>= 1.0.0), testthat, Cubist, arm, MASS, kknn, kernlab, ipred, randomForest, pROC
<b>Depends</b> R ( $>= 2.10$ ), caret
<b>Imports</b> parallel, subselect, plyr, xgboost, magrittr, stringr, e1071, glmnet, verification
R topics documented:
ff.bindPath  ff.blend  ff.corrFilter  ff.createEnsemble  ff.encodeCategoricalFeature  ff.extractDateFeature  ff.featureFilter  ff.getBestBlenderPerformance  ff.getBestBlenderTune  ff.getMaxCuncurrentThreads

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ff.bindPath

Bind an absolute path for a kind of resources.

# Description

Bind an absolute path for a kind of resources.

### Usage

```
ff.bindPath(type, sub_path, createDir = FALSE)
```

# Arguments

type the type of resource.

sub\_path the suffix to concatenate to the absolute path to get the absolute path of the kind

of resource.

createDir set to TRUE to create the directory if it does not exist

# **Examples**

```
ff.setBasePath(getwd())
if(! dir.exists("mydata") ) dir.create(mydata)
ff.bindPath(type = "data",sub_path = "mydata")
```

ff.blend

Given a tuned regression model, finds more performant tuning configurations using Nelder/Mead, quasi-Newton and conjugate-gradient algorithms.

# Description

Given a tuned regression model, finds more performant tuning configurations using Nelder/Mead, quasi-Newton and conjugate-gradient algorithms.

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#### **Usage**

```
ff.blend(bestTune, caretModelName, Xtrain, y, controlObject, max_secs = 10 *
60, seed = NULL, method = c("Nelder-Mead", "BFGS", "CG", "L-BFGS-B",
    "SANN"), useInteger = TRUE, parallelize = TRUE, verbose = TRUE)
```

#### **Arguments**

bestTune a data. frame with best tuned parameters of specified model.

caretModelName a string specifying which model to use. Possible values are lm, bayesglm,

glm, glmStepAIC, rlm, knn, pls, ridge, enet, svmRadial, treebag, gbm, rf,

cubist, avNNet, xgbTreeGTJ, xgbTree

Xtrain the encoded data. frame of train data. Must be a data. frame of numeric

y the output variable as numeric vector

controlObject a list of values that define how this function acts. Must be a caret trainControl

object

max\_secs the max number of seconds as time constraint

seed a user specified seed. Useful for replicable execution (e.g. passing the same

seed to the ff.verifyBlender function) if the control object involves random

steps for creating resamples.

method the method to use. Possible values are c(Nelder-Mead, BFGS, CG, L-BFGS-B, SANN).

useInteger TRUE if the tuning grid is composed of integers and not of continuous numbers.

parallelize TRUE to enable parallelization (require parallel).

verbose TRUE to enable verbose mode.

### Value

a list of lists (one for each specified optimization method) with components par (best set of parameters found), value (the value of fn corresponding to par), counts (a two-element integer vector giving the number of calls to fn and gr respectively; this excludes those calls needed to compute the Hessian, if requested, and any calls to fn to compute a finite-difference approximation to the gradient), convergence (an integer code. 0 indicates successful completion which is always the case for SANN and Brent), message (a character string giving any additional information returned by the optimizer, or NULL), seed (the used seed). For further details see optim.

#### References

```
https://stat.ethz.ch/pipermail/r-devel/2010-August/058081.html
```

```
## suppress warnings raised because there few obs
warn_def = getOption(warn)
options(warn=-1)

## data
Xtrain <- data.frame( a = rep(1:5 , each = 2), b = 10:1,
c = rep(as.Date(c("2007-06-22", "2004-02-13")),5) )
Xtest <- data.frame( a = rep(2:6 , each = 2), b = 1:10,
c = rep(as.Date(c("2007-03-01", "2004-05-23")),5) )
Ytrain = 1:10 + runif(nrow(Xtrain))</pre>
```

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```
## encode datasets
1 = ff.makeFeatureSet(Xtrain, Xtest, c("C", "N", "D"))
Xtrain = 1$traindata
Xtest = 1$testdata
## make a caret control object
controlObject <- trainControl(method = "repeatedcv",</pre>
repeats = 1, number = 2)
## train and predict
tp = ff.trainAndPredict.reg(Ytrain=Ytrain ,
                          Xtrain=Xtrain ,
                          Xtest=Xtest ,
                          model.label = "cubist" ,
                          controlObject=controlObject)
pred_test = tp*pred
model = tp\$model
secs = tp$secs
## blender
gBlender = ff.blend(bestTune = tp$model$bestTune,
                                caretModelName = "cubist" ,
                                Xtrain = Xtrain ,
                                 y = Ytrain, controlObject = tp$model$control,
                                max_secs = 3,
                                 seed = 123,
                                 method = c("Nelder-Mead", "BFGS", "CG", "L-BFGS-B", "SANN"),
                                 useInteger = TRUE,
                                 parallelize = TRUE,
                                 verbose = FALSE)
ff.summaryBlender(gBlender)
ff.getBestBlenderPerformance(gBlender)
bestTune = ff.getBestBlenderTune(gBlender)
ff.verifyBlender (gBlender,Xtrain=Xtrain,y=Ytrain,seed=123,
controlObject=tp$model$control,caretModelname = "cubist")
## restore warnings
options(warn=warn_def)
```

ff.corrFilter

 $\it Filter\ a\ data.$  frame of numeric according to a given threshold of correlation

# **Description**

Filter a data. frame of numeric according to a given threshold of correlation

### Usage

```
ff.corrFilter(Xtrain, Xtest, y, abs_th = NULL, rel_th = 1,
  method = "pearson")
```

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#### **Arguments**

Xtrain	a train set data.frame of numeric
Xtest	a test set data.frame of numeric
У	the output variable (as numeric vector)
abs_th	an absolute threshold (= number of data frame columns)
rel_th	a relative threshold (= percentage of data frame columns)
method	a character string indicating which correlation method is to be used for the test. One of "pearson", "kendall", or "spearman".

### Value

a list of filtered train set and test set with correlation test results

# **Examples**

```
Xtrain <- data.frame( a = rep(1:3 , each = 2), b = c(4:1,6,6), c = rep(1,6))
Xtest <- Xtrain + runif(nrow(Xtrain))
y = 1:6
l = ff.corrFilter(Xtrain=Xtrain, Xtest=Xtest, y=y, rel_th=0.5)
Xtrain.filtered = l$Xtrain
Xtest.filtered = l$Xtest</pre>
```

ff.createEnsemble

Create an ensemble of a tuned model

# **Description**

Create an ensemble of a tuned model

# Usage

```
ff.createEnsemble(Xtrain, Xtest, y, caretModelName, bestTune, predTest = NULL,
  removePredictorsMakingIllConditionedSquareMatrix_forLinearModels = TRUE,
  controlObject, parallelize = TRUE, verbose = TRUE, regression = TRUE,
  ...)
```

### **Arguments**

Xtrain	the encoded data.frame of train data. Must be a data.frame of numeric
Xtest	the encoded data. frame of test data. Must be a data.frame of numeric
у	the output variable as numeric vector
caretModelName	a string specifying which model to use. Possible values for regression are lm, bayesglm, glm, glmStepAIC, rlm, knn, pls, ridge, enet, svmRadial, treebag, gbm, rf, cubist, avNNet, xgbTreeGTJ, xgbTree.
bestTune	a data.frame with best tuned parameters of specified model.
predTest	test set prediction (numeric vector). If available, passing it through this paramter the function doesn't compute it again for creating the esemble.

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 $remove Predictors Making Ill Conditioned Square Matrix\_for Linear Models$ 

TRUE for removing predictors making ill-conditioned square matrices in case of

fragile linear models, i.e. c(rlm, pls, ridge, enet) for regression.

controlObject a list of values that define how this function acts. Must be a caret trainControl

object

parallelize TRUE to enable parallelization (require parallel).

verbose TRUE to enable verbose mode.

regression TRUE to create an ensemble of a tuned regression model and FALSE to create an

ensemble of a tuned classification model.

... arguments passed to the regression routine.

#### Value

a list of train and test predictions.

```
## suppress warnings raised because there few obs
warn_def = getOption(warn)
options(warn=-1)
Xtrain \leftarrow data.frame( a = rep(1:10 , each = 2), b = 20:1,
c = rep(as.Date(c("2007-06-22", "2004-02-13")),10))
Xtest \leftarrow data.frame( a = rep(2:11 , each = 2), b = 1:20,
c = rep(as.Date(c("2007-03-01", "2004-05-23")),10))
Ytrain = 1:20 + runif(nrow(Xtrain))
## encode datasets
1 = ff.makeFeatureSet(Xtrain, Xtest, c("C", "N", "D"))
Xtrain = l$traindata
Xtest = 1$testdata
## make a caret control object
controlObject <- trainControl(method = "repeatedcv", repeats = 1, number = 2)</pre>
tp = ff.trainAndPredict.reg(Ytrain=Ytrain ,
                          Xtrain=Xtrain ,
                          Xtest=Xtest ,
                           model.label = "cubist" ,
                           controlObject=controlObject)
pred_test = tp$pred
model = tp$model
secs = tp$secs
## create ensemble
en = ff.createEnsemble(Xtrain = Xtrain,
                      Xtest = Xtest,
                      y = Ytrain,
                      bestTune = tp$model$bestTune ,
                      caretModelName = "cubist" ,
                      parallelize = TRUE,
                      removePredictorsMakingIllConditionedSquareMatrix_forLinearModels = TRUE,
                      controlObject = tp$model$control)
```

```
predTrain = en$predTrain
predTest = en$predTest

## restore warnings
options(warn=warn_def)
```

#### ff.encodeCategoricalFeature

Encode a generic predictor as a categorical features using both observations of train set and test for levels. It's anyway possible to adopt more levels by using the parameter levels. Notice that modeling a generic vector, e.g. c(1,2,3,4,5,2,3) as a categorical predictor xor a numeric predictor is a modeling choice (eventually to be assessed by cross-validation).

# Description

Encode a generic predictor as a categorical features using both observations of train set and test for levels. It's anyway possible to adopt more levels by using the parameter levels. Notice that modeling a generic vector, e.g. c(1,2,3,4,5,2,3) as a categorical predictor xor a numeric predictor is a modeling choice (eventually to be assessed by cross-validation).

### Usage

```
ff.encodeCategoricalFeature(data.train, data.test, colname.prefix,
   asNumericSequence = F, replaceWhiteSpaceInLevelsWith = NULL,
   levels = NULL, remove1DummyVar = FALSE)
```

### **Arguments**

data.train the observations of the predictor in train set.
data.test the observations of the predictor in test set.
colname.prefix the prefix of output data frame.
asNumericSequence

set T if the predictor is a numeric sequence filling any possible hole between min and max in observations that could occour both in train set and test set.

replaceWhiteSpaceInLevelsWith

replace possible spaces in the train/test name of feature.

levels the levels of the categorical feature. Must be NULL if as Numeric Sequence is T. remove 1 Dummy Var

T to remove one dummy variable. Why? First, if you know the values of the first C - 1 dummy variables, you know the last one too and it is more economical to use C - 1. Secondly, if the model has slopes and intercepts (e.g. linear regression), the sum of all of the dummy variables wil add up to the intercept (usually encoded as a "1") and that is bad for the math involved. On the other hand, there are models like penalized methods (such as ridge regression) that seldom penalize the intercept, so a C-1 encoded variable could cause the other category effects to be penalized towards the reference category effect.

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#### Value

the list of trainset and testset after applying the specified filters

#### References

http://appliedpredictivemodeling.com/blog/2013/10/23/the-basics-of-encoding-categorical-data-fo

#### Examples

```
Xtrain \leftarrow data.frame( a = rep(1:3 , each = 2), b = 6:1, c = letters[1:6])
Xtest <- data.frame( a = rep(2:4 , each = 2), b = 1:6, c = letters[6:1])
print(Xtrain)
  a b c
# 1 1 6 a
# 2 1 5 b
# 3 2 4 c
# 4 2 3 d
# 5 3 2 e
#631f
1 = ff.encodeCategoricalFeature (Xtrain$c , Xtest$c , "c")
1$traindata
     c_1 c_2 c_3 c_4 c_5 c_6
# 7
      1 0 0
                0
# 8
      0 1
             0 0
# 9
      0 0 1
                0 0
                         0
# 10
     0 0 0 1
                     0 0
# 11
     0 0 0 0 1 0
# 12
     0 0 0
                        1
Xtrain[,c] = NULL
Xtest[,c] = NULL
Xtrain = cbind(Xtrain, 1$traindata)
Xtest = cbind(Xtest, 1$testdata)
```

ff.extractDateFeature Extracts a numerical feature from a date predictor. The feature is built as the difference in days from the oldest date in bothe train set and test set and any given observation.

### Description

Extracts a numerical feature from a date predictor. The feature is built as the difference in days from the oldest date in bothe train set and test set and any given observation.

#### Usage

```
ff.extractDateFeature(data.train, data.test)
```

### **Arguments**

```
data.train the observations of the predictor in train set.
data.test the observations of the predictor in test set.
```

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#### Value

the list of trainset and testset after applying the specified encoding and the related date range

#### **Examples**

```
Xtrain <- data.frame( a = rep(1:3 , each = 2), b = 6:1,
    c = rep(as.Date(c("2007-06-22", "2004-02-13")),3) )
Xtest <- data.frame( a = rep(2:4 , each = 2), b = 1:6,
    c = rep(as.Date(c("2007-03-01", "2004-05-23")),3) )
1 = ff.extractDateFeature(Xtrain$c,Xtest$c)
Xtrain[,c] = NULL
Xtest[,c] = NULL
Xtrain = cbind(Xtrain,c=1$traindata)
Xtest = cbind(Xtest,c=1$testdata)</pre>
```

ff.featureFilter

Filter predictors according to specified criteria.

#### **Description**

Filter predictors according to specified criteria.

featureScaling TRUE to perform feature scaling

TRUE to set verbose mode

# Usage

```
ff.featureFilter(traindata, testdata, y = NULL,
  removeOnlyZeroVariacePredictors = FALSE,
  performVarianceAnalysisOnTrainSetOnly = TRUE, correlationThreshold = NULL,
  removePredictorsMakingIllConditionedSquareMatrix = TRUE,
  removeIdenticalPredictors = TRUE, removeHighCorrelatedPredictors = TRUE,
  featureScaling = TRUE, verbose = TRUE)
```

#### **Arguments**

verbose

traindata the train set testdata the test set the response variable. Must be not NULL if correlationThreshold is not NULL. removeOnlyZeroVariacePredictors TRUE to remove only zero variance predictors performVarianceAnalysisOnTrainSetOnly TRUE to perform the variance analysis on the train set only correlationThreshold a correlation threshold above which keeping predictors (considered only if removeOnlyZeroVariacel is FALSE). remove Predictors Making Ill Conditioned Square MatrixTRUE to predictors making ill conditioned square matrices removeIdenticalPredictors TRUE to remove identical predictors (using base::identical function) remove High Correlated PredictorsTRUE to remove high correlared predictors

#### Value

the list of trainset and testset after applying the specified filters

### **Examples**

ff.getBestBlenderPerformance

Helper function that given a blender object returns the best optimization method.

#### **Description**

Helper function that given a blender object returns the best optimization method.

#### Usage

```
ff.getBestBlenderPerformance(blender)
```

### **Arguments**

blender a blender object

## Value

a numeric of best score and as object name the best performant method name.

# See Also

ff.blend for examples.

ff.getBestBlenderTune Helper function that given a blender object returns the best tuning parameters found by the blender.

### **Description**

Helper function that given a blender object returns the best tuning parameters found by the blender.

# Usage

```
ff.getBestBlenderTune(blender, truncate = TRUE)
```

## **Arguments**

blender a blender object

truncate TRUE to cut at the first tuning best configuration in case there are more than one

optimal tuning configurations.

#### Value

a data. frame of the best tuning parameters.

#### See Also

ff.blend for examples.

 ${\tt ff.getMaxCuncurrentThreads}$ 

Get the max number of cuncurrent threads.

### **Description**

Get the max number of cuncurrent threads.

#### Usage

```
ff.getMaxCuncurrentThreads()
```

### **Examples**

ff.getMaxCuncurrentThreads()

ff.getPath

Get the absolute path for a kind of resources.

#### **Description**

Get the absolute path for a kind of resources.

#### Usage

```
ff.getPath(type = "base")
```

# Arguments

type the type of resource.

# Value

the absolute path for a kind of resources (as character)

```
ff.setBasePath(./)
ff.getPath() ## equivalent to ff.getPath(type="base")
```

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ff.getPathBindings	Get the list of bindings, i.e. (type resource, absolute path) pairs as a list
--------------------	---

### **Description**

Get the list of bindings, i.e. (type resource, absolute path) pairs as a list

### Usage

```
ff.getPathBindings()
```

#### Value

the list of bindings

### **Examples**

```
ff.setBasePath(getwd())
if(! dir.exists("mydata") ) dir.create(mydata)
ff.bindPath(type = "data",sub_path = "mydata")
ff.getPathBindings()
```

ff.kmeans

An useful wrapper of kmeans performing k-means clustering on the given trainset / testset (Xtrain / Xtest) and assuming a number of cluster from 1:max\_centers. The best number of cluster is computed so that the variation in the within group sum of squares between two subsequent number of clusters is maximized in absolute value.

# Description

An useful wrapper of kmeans performing k-means clustering on the given trainset / testset (Xtrain / Xtest) and assuming a number of cluster from 1:max\_centers. The best number of cluster is computed so that the variation in the within group sum of squares between two subsequent number of clusters is maximized in absolute value.

#### Usage

```
ff.kmeans(Xtrain, Xtest, max_centers = 10, nstart = 5, iter.max = 10,
    doPlot = FALSE, verbose = FALSE)
```

### **Arguments**

Xtrain	the encoded data. frame of train data. Must be a data. frame of numeric
Xtest	the encoded data. frame of train data. Must be a data. frame of numeric
max_centers	the max number of clusters to be evaluated
nstart	how many random sets should be chosen? Such a parameter is passed to kmeans.

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iter.max the maximum number of iterations allowed. Such a parameter is passed to

kmeans.

doPlot a logical value indicating whether plotting.

verbose a logical value indicating whether verbose mode should be enabled.

varThreshold a threshold indicating the proportion of variance that should be explained. Must

be a numeric between 0 and 1.

#### Value

a list whose components are the max number of clusters evaluated (max\_centers), the best number of clusters (best\_n\_cluters), the sequence of within groups sum of squares across number of clusters (wss), assuming best\_n\_cluters as number of clusters a vector of integers (from 1:best\_n\_cluters) indicating the cluster to which each point is allocated in the train set (K.train) and test set (K.test).

#### **Examples**

ff.makeFeatureSet

*Encode the feature set according to meta data passed as input.* 

### **Description**

Encode the feature set according to meta data passed as input.

### Usage

```
ff.makeFeatureSet(data.train, data.test, meta, scaleNumericFeatures = FALSE,
    parallelize = FALSE, remove1DummyVarInCatPreds = FALSE)
```

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#### **Arguments**

data.train the observations of the predictor in train set.

data.test the observations of the predictor in test set.

meta the meata data. It should be a vector of the character C, N, D, e.g. c(N,C,D) of

the same length of the train set / test set columns

scaleNumericFeatures

seto to TRUE to center and scale numeric features

parallelize set to TRUE to enable parallelization (require parallel package)

remove1DummyVarInCatPreds

T to remove one dummy variable in encoding categorical predictors. For further details see ff.encodeCategoricalFeature.

#### Value

the list of trainset and testset after applying the specified encodings

#### **Examples**

```
Xtrain <- data.frame( a = rep(1:3 , each = 2), b = 6:1,
    c = rep(as.Date(c("2007-06-22", "2004-02-13")),3) )
Xtest <- data.frame( a = rep(2:4 , each = 2), b = 1:6,
    c = rep(as.Date(c("2007-03-01", "2004-05-23")),3) )
1 = ff.makeFeatureSet(Xtrain, Xtest, c(C, N, D))
Xtrain = l$traindata
Xtest = l$testdata</pre>
```

ff.pca

An useful wrapper of prcomp performing a principal components analysis on the given trainset / testset (Xtrain / Xtest).

# Description

An useful wrapper of prcomp performing a principal components analysis on the given trainset / testset (Xtrain / Xtest).

#### Usage

```
ff.pca(Xtrain, Xtest, center = TRUE, scale. = FALSE,
  removeZeroVarPredictors = TRUE, varThreshold = 0.95, doPlot = TRUE,
  verbose = FALSE)
```

# **Arguments**

Xtrain the encoded data.frame of train data. Must be a data.frame of numeric
the encoded data.frame of train data. Must be a data.frame of numeric
a logical value indicating whether the variables should be shifted to be zero
centered. Alternately, a vector of length equal the number of columns of data
can be supplied. The value is passed to scale.

ff.pca

scale.

a logical value indicating whether the variables should be scaled to have unit variance before the analysis takes place. The default is FALSE for consistency with S, but in general scaling is advisable. Alternatively, a vector of length equal the number of columns of data can be supplied. The value is passed to scale.

removeZeroVarPredictors

a logical value indicating whether removing zero variance predictors before calling prcomp preventing errors due to the fact that the latter cannot rescale a context of the context of th

stant/zero column to unit variance.

varThreshold a threshold indicating the proportion of variance that should be explained. Must

be a numeric between 0 and 1.

doPlot a logical value indicating whether plotting the proportion of variance explained

vs. principal components.

verbose a logical value indicating whether verbose mode should be enabled.

#### Value

a list whose components are the number of principal components (numComp), the number of principal components to hold so that the proportion of variance explained by each subsequent principal component drops off as an elbow in the screen plot (numComp.elbow), the number of principal components explaining a given (specified by the varThreshold input parameter) proportion of variance (numComp.threshold), the threshold indicating the proportion of variance that should be explained (varThreshold), the cumulative sum of proportion of variance explained by each principal component (cumVar), the proportion of variance explained by each principal components for train and test set (PC.train and PC.test)

```
Xtrain \leftarrow data.frame(a = rep(1:10 , each = 2), b = 20:1,
                     c = rep(as.Date(c("2007-06-22", "2004-02-13")), 10), d = 20:1)
Xtest \leftarrow data.frame(a = rep(2:11, each = 2), b = 1:20,
                    c = rep(as.Date(c("2007-03-01", "2004-05-23")),10), d = 1:20)
## encode data sets
1 = ff.makeFeatureSet(Xtrain, Xtest, c("C", "N", "D", "N"))
Xtrain = 1$traindata
Xtest = 1$testdata
ffPCA = ff.pca(Xtrain = Xtrain , Xtest = Xtest , center = TRUE , scale. = TRUE ,
                removeZeroVarPredictors = TRUE ,
                varThreshold = 0.95 , doPlot = FALSE , verbose = TRUE)
numComp <- ffPCA$numComp</pre>
numComp.elbow <- ffPCA$numComp.elbow</pre>
numComp.threshold <- ffPCA$numComp.threshold</pre>
PC_Xtrain_95Var = ffPCA$PC.train[1:numComp.threshold,,drop=FALSE]
PC_Xtest_95Var = ffPCA$PC.test[1:numComp.threshold,,drop=FALSE]
```

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```
ff.plotPerformance.reg
```

Plot predicted values vs. observed / residual values.

### **Description**

Plot predicted values vs. observed / residual values.

### Usage

```
ff.plotPerformance.reg(observed, predicted, main = NULL)
```

## Arguments

observed the observed output variables (numeric vector).

predicted the predicted values (numeric vector).

main a string as a title for the plot

# **Examples**

```
obs = 1:10
preds = obs + runif(length(obs))
ff.plotPerformance.reg(observed = obs , predicted = preds, main="Predicted vs. observed/residual")
```

ff.poly

Make polynomial terms of a data.frame

### **Description**

Make polynomial terms of a data. frame

# Usage

```
ff.poly(x, n, direction = 0)
```

### **Arguments**

x a data.frame of numeric n the polynomial degree

direction if set to 0 returns the terms  $x^{(1/n)}, x^{(1/(n-1))}, \dots, x, x^2, \dots, x^n$ . If

set to -1 returns the terms  $x^{(1/n)}, x^{(1/(n-1))}, \dots, x$ . If set to 1 returns the

terms  $x, x^2, \dots, x^n$ .

#### Value

the data. frame with the specified polynomial terms

ff.setBasePath

#### **Examples**

```
Xtrain <- data.frame( a = rep(1:3 , each = 2), b = c(4:1,6,6), c = rep(1,6))
Xtest <- Xtrain + runif(nrow(Xtrain))
data = rbind(Xtrain,Xtest)
data.poly = ff.poly(x=data,n=3)
Xtrain.poly = data.poly[1:nrow(Xtrain),]
Xtest.poly = data.poly[(nrow(Xtrain)+1):nrow(data),]</pre>
```

ff.setBasePath

Set base path

# Description

Set base path

# Usage

```
ff.setBasePath(path)
```

# Arguments

path

the absolute path.

### **Examples**

```
ff.setBasePath(./)
```

 $\verb"ff.setMaxCuncurrentThreads"$ 

Set the max number of cuncurrent threads.

# **Description**

Set the max number of cuncurrent threads.

# Usage

```
ff.setMaxCuncurrentThreads(nThreads = 2)
```

# **Arguments**

nThreads

max number of cuncurrent threads.

```
ff.setMaxCuncurrentThreads(4)
```

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ff.summaryBlender	Helper function that given a blender object returns a numeric vector
	of performances (one for each optimization method).

### **Description**

Helper function that given a blender object returns a numeric vector of performances (one for each optimization method).

### Usage

```
ff.summaryBlender(blender)
```

### **Arguments**

```
blender a blender object
```

### Value

a numeric vector of performances (one for each optimization method)

#### See Also

ff.blend for examples.

```
ff.trainAndPredict.class
```

Trains a specified classification model on the given train set and predicts on the given test set.

### **Description**

Trains a specified classification model on the given train set and predicts on the given test set.

# Usage

```
ff.trainAndPredict.class(Ytrain, Xtrain, Xtest, model.label, controlObject,
  best.tuning = FALSE, verbose = FALSE,
  removePredictorsMakingIllConditionedSquareMatrix_forLinearModels = TRUE,
  metric.label = "auc", xgb.metric.fun = NULL, xgb.maximize = FALSE,
  xgb.foldList = NULL, xgb.eta = NULL, xgb.max_depth = NULL,
  xgb.cv.default = TRUE, xgb.param = NULL, ...)
```

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#### **Arguments**

Ytrain	the output variable as numeric vector
Xtrain	the encoded data.frame of train data. Must be a data.frame of numeric
Xtest	the encoded data. frame of test data. Must be a data. frame of numeric
model.label	a string specifying which model to use.
controlObject	a list of values that define how this function acts. Must be a caret trainControl object for all models except that for xgbTreeGTJ.
best.tuning	TRUE to use more dense tuning grid or custom routine/tuning grid if available
verbose removePredictor	TRUE to enable verbose mode.  *SMakingIllConditionedSquareMatrix_forLinearModels  TRUE for removing predictors making ill-conditioned square matrices in case of fragile linear models.
metric.label	the label of function to optmize/minimize.
xgb.metric.fun	custom function to optmize/minimize for xgbTreeGTJ.
xgb.maximize	TRUE to maximize the specified xgb.metric.fun.
xgb.foldList	custom resampling folds list for xgbTreeGTJ.
xgb.eta	custom eta parameter for xgbTreeGTJ.
xgb.max_depth	custom max_depth parameter for xgbTreeGTJ.
xgb.cv.default	TRUE for using xgboost::xgb.cv function (mandatory in case of fix nrounds), FALSE for using the internal ff.xgb.cv function. The main advantage of the latter is that it doesn't need to restart nrounds in case for the specified nrounds cross validation error is still decreasing.
xgb.param	custom parameters for XGBoost.
	arguments passed to the regression routine.

# Value

a list of test predictions, model and number of excecuting seconds.

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ff.trainAndPredict.reg

Trains a specified model on the given train set and predicts on the given test set.

# Description

Trains a specified model on the given train set and predicts on the given test set.

## Usage

```
ff.trainAndPredict.reg(Ytrain, Xtrain, Xtest, model.label, controlObject,
  best.tuning = FALSE, verbose = FALSE,
  removePredictorsMakingIllConditionedSquareMatrix_forLinearModels = TRUE,
  xgb.metric.fun = RMSLE.xgb, xgb.maximize = FALSE,
  xgb.metric.label = "rmsle", xgb.foldList = NULL, xgb.eta = NULL,
  xgb.max_depth = NULL, xgb.cv.default = TRUE, xgb.param = NULL, ...)
```

### **Arguments**

Ytrain	the output variable as numeric vector
Xtrain	the encoded data.frame of train data. Must be a data.frame of numeric
Xtest	the encoded data. frame of test data. Must be a data. frame of numeric
model.label	a string specifying which model to use. Possible values are lm, bayesglm, glm, glmStepAIC, rlm, knn, pls, ridge, enet, svmRadial, treebag, gbm, rf, cubist, avNNet, xgbTreeGTJ, xgbTree
controlObject	a list of values that define how this function acts. Must be a caret trainControl object for all models except that for xgbTreeGTJ and xgbTree. In the latter case only if best.tuning is TRUE.
best.tuning	TRUE to use more dense tuning grid or custom routine if available
verbose	TRUE to enable verbose mode.

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 $remove Predictors Making Ill Conditioned Square Matrix\_for Linear Models$ TRUE for removing predictors making ill-conditioned square matrices in case of fragile linear models, i.e. c(rlm, pls, ridge, enet). xgb.metric.fun custom function to optmize/minimize for xgbTreeGTJ and xgbTree. In the latter case only if best. tuning is TRUE. xgb.maximize TRUE to maximize the specified xgb.metric.fun. Only for xgbTreeGTJ and xgbTree. In the latter case only if best.tuning is TRUE. xgb.metric.label custom label of function to optmize/minimize for xgbTreeGTJ and xgbTree. In the latter case only if best. tuning is TRUE. xgb.foldList custom resampling folds list for xgbTreeGTJ and xgbTree. In the latter case only if best. tuning is TRUE. custom eta parameter for xgbTreeGTJ and xgbTree. In the latter case only if xgb.eta best.tuning is TRUE. custom max\_depth parameter for xgbTreeGTJ and xgbTree. In the latter case xgb.max\_depth only if best.tuning is TRUE. xgb.cv.default TRUE for using xgboost::xgb.cv function (mandatory in case of fix nrounds), FALSE for using the internal ff.xgb.cv function. The main advantage of the latter is that it doesn't need to restart nrounds in case for the specified nrounds cross validation error is still decreasing. custom parameters for XGBoost. xgb.param arguments passed to the regression routine.

#### Value

a list of test predictions, model and number of excecuting seconds.

```
## suppress warnings raised because of few obs
warn_def = getOption(warn)
options(warn=-1)
## data
Xtrain \leftarrow data.frame( a = rep(1:10 , each = 2), b = 20:1,
c = rep(as.Date(c("2007-06-22", "2004-02-13")),10))
Xtest <- data.frame( a = rep(2:11 , each = 2), b = 1:20,
c = rep(as.Date(c("2007-03-01", "2004-05-23")),10))
Ytrain = 1:20 + runif(nrow(Xtrain))
## encode datasets
1 = ff.makeFeatureSet(Xtrain, Xtest, c("C", "N", "D"))
Xtrain = 1$traindata
Xtest = 1$testdata
## make a caret control object
controlObject <- trainControl(method = "repeatedcv", repeats = 1, number = 2)</pre>
tp = ff.trainAndPredict.reg(Ytrain=Ytrain ,
                          Xtrain=Xtrain ,
                           Xtest=Xtest ,
                           model.label = "cubist" ,
```

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#### controlObject=controlObject)

```
pred_test = tp$pred
model = tp$model
elapsed.secs = tp$secs
## restore warnings
options(warn=warn_def)
```

ff.verifyBlender

Helper function that given a blender object replicates the execution in order to verify performances.

### **Description**

Helper function that given a blender object replicates the execution in order to verify performances.

### Usage

```
ff.verifyBlender(blender, Xtrain, y, seed = NULL, controlObject,
  caretModelname)
```

### **Arguments**

blender a blender object
Xtrain the train set

y the output variable as numeric vector

seed the seed used by the blender, if applicable. If the blender used one, it is necessary

for replicating blender performances.

controlObject a list of values that define how this function acts. Must be a caret trainControl

object. It must be the same used by the blender.

caretModelname a string specifying which model to use. Possible values are lm, bayesglm,

glm, glmStepAIC, rlm, knn, pls, ridge, enet, svmRadial, treebag, gbm, rf, cubist, avNNet, xgbTreeGTJ, xgbTree. It must be the same model name used

by the blender.

# Value

a numeric as difference in performance between blender and replicated execution.

# See Also

ff.blend for examples.

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RMSE.xgb

Root mean square error

# **Description**

Root mean square error

### Usage

```
RMSE.xgb(preds, dtrain)
```

### **Arguments**

preds the predicted values (numeric vector).

dtrain the xgboost train set object.

# Value

a list of metric label / values

RMSLE.xgb

Root mean square logistic error

# Description

Root mean square logistic error

# Usage

```
RMSLE.xgb(preds, dtrain, th_err = 1.5)
```

# Arguments

preds the predicted values (numeric vector).

dtrain the xgboost train set object.

th\_err a threshold in case predictions are negative.

### Value

a list of metric label / values

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