## Package 'LedPred'

### April 26, 2024

Title Learning from DNA to Predict Enhancers

**Description** This package aims at creating a predictive model of regulatory sequences used to score unknown sequences based on the content of DNA motifs, next-generation sequencing (NGS) peaks and signals and other numerical scores of the sequences using supervised classification. The package contains a workflow based on the support vector machine (SVM) algorithm that maps features to sequences, optimize SVM parameters and feature number and creates a model that can be stored and used to score the regulatory potential of unknown sequences.

Version 1.37.0

**Date** 2016-08-13

Author Elodie Darbo, Denis Seyres, Aitor Gonzalez

Maintainer Aitor Gonzalez <aitor.gonzalez@univ-amu.fr>

**Depends** R (>= 3.2.0), e1071 (>= 1.6)

**Imports** akima, ggplot2, irr, jsonlite, parallel, plot3D, plyr, RCurl, ROCR, testthat

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LazyData true

**biocViews** SupportVectorMachine, Software, MotifAnnotation, ChIPSeq, Sequencing, Classification

NeedsCompilation no

BugReports https://github.com/aitgon/LedPred/issues

RoxygenNote 5.0.1

git\_url https://git.bioconductor.org/packages/LedPred

git\_branch devel

git\_last\_commit de57379

git\_last\_commit\_date 2023-10-24

**Repository** Bioconductor 3.19

Date/Publication 2024-04-26

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createModel

Create the model with the optimal features

### Description

createModel function creates a SVM model from the training data set with the selected features.

### Usage

```
createModel(data, cl = 1, kernel = "radial", cost = 1, gamma = 1,
valid.times = 10, feature.ranking = NULL, feature.nb = NULL,
file.prefix = NULL)
```

### Arguments

data	data.frame containing the training set	
cl	integer indicating the column number corresponding to the response vector that classify positive and negative regions (default = 1)	
kernel	SVM kernel, a character string: "linear" or "radial". (default = "radial")	
cost	The SVM cost parameter for both linear and radial kernels. If NULL (default), the function mcTune is run.	
gamma	The SVM gamma parameter for radial kernel. If radial kernel and NULL (default), the function mcTune is run.	
valid.times	Integer indicating how many times the training set will be split for the cross validation step (default = 10). This number must be smaller than positive and negative sets sizes.	
feature.ranking		
	List of ordered features.	
feature.nb	the optimal number of feature to use from the list of ordered features.	
file.prefix	A character string that will be used as a prefix followed by "_model.RData" for the resulting model file, if it is NULL (default), no model is saved	

### crm.features

### Value

the best SVM model

### Examples

crm. features This is data to be included in my package

### Description

This is data to be included in my package

evaluateModelPerformance

Evaluate model performances

#### Description

evaluateModelPerformance function computes the precision and recall measures to evaluate the model through cross validation steps using ROCR package.

### Usage

```
evaluateModelPerformance(data, cl = 1, valid.times = 10,
feature.ranking = NULL, feature.nb = NULL,
numcores = ifelse(.Platform$OS.type == "windows", 1, parallel::detectCores()
- 1), file.prefix = NULL, kernel = "linear", cost = NULL,
gamma = NULL)
```

### Arguments

data	data.frame containing the training set
cl	integer indicating the column number corresponding to the response vector that classify positive and negative regions (default = 1)
valid.times	Integer indicating how many times the training set will be split for the cross validation step (default = $10$ ). This number must be smaller than positive and negative sets sizes.

feature.ranking		
	List of ordered features.	
feature.nb	the optimal number of feature to use from the list of ordered features.	
numcores	Number of cores to use for parallel computing (default: the number of available cores in the machine - 1)	
file.prefix	A character string that will be used as a prefix followed by "_ROCR_perf.png" for the result plot file, if it is NULL (default), no plot is returned	
kernel	SVM kernel, a character string: "linear" or "radial". (default = "radial")	
cost	The SVM cost parameter for both linear and radial kernels. If NULL (default), the function mcTune is run.	
gamma	The SVM gamma parameter for radial kernel. If radial kernel and NULL (de-fault), the function mcTune is run.	

### Value

A list with two objects.

probs	The predictions computed by the model for each subset during the cross-validation
labels	The actual class for each subset

### Examples

feature.ranking	This is data to be included in my package
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### Description

This is data to be included in my package

### Description

The LedPred function computes the best SVM parameters, defines the optimal features for creating the SVM model by running sequentially mcTune, rankFeatures, tuneFeatureNb and createModel. The performances of this model are then computed usong evaluateModelPerformance.

### Usage

```
LedPred(data = NULL, cl = 1, ranges = list(gamma = c(1, 10), cost = c(1,
10)), cost = NULL, gamma = NULL, kernel = "linear", valid.times = 10,
file.prefix = NULL, numcores = ifelse(.Platform$0S.type == "windows", 1,
parallel::detectCores() - 1), step.nb = 10, halve.above = 100)
```

### Arguments

data	data.frame containing the training set
cl	integer indicating the column number corresponding to the response vector that classify positive and negative regions (default = 1)
ranges	list object containing one (linear kernel) or two (radial kernel) vectors of integers corresponding to SVM cost and SVM gamma parameters to test.
cost	The SVM cost parameter for both linear and radial kernels. If NULL (default), the function mcTune is run.
gamma	The SVM gamma parameter for radial kernel. If radial kernel and NULL (default), the function mcTune is run.
kernel	SVM kernel, a character string: "linear" or "radial". (default = "radial")
valid.times	Integer indicating how many times the training set will be split for the cross validation step (default = 10). This number must be smaller than positive and negative sets sizes.
file.prefix	A character string that will be used as a prefix for the result files. If it is NULL (default), no plot is returned
numcores	Number of cores to use for parallel computing (default: the number of available cores in the machine - 1)
step.nb	Number of features to add at each step (default = $10$ )
halve.above	During RFE, all the features are ranked at the first round and the half lowest ranked features (that contribute the least in the model) are removed for the next round. When the number of feature is lower or equal to halve.above, the features are removed one by one. (default=100)

### Value

A list of the object produced at each step

best.params	A list of the parameters giving the lowest misclassification error	
feature.ranking		
	List of ordered features from rankFeatures	
feature.nb	he optimal number of feature to use from the list of ordered features from ${\tt tuneFeatureNb}$	
model.svm	The best SVM model createModel	
probs.label.list		
	The cross-validation results from evaluateModelPerformance	

### Examples

```
data(crm.features)
#cost_vector <- c(1,3,10)
#gamma_vector <- c(1,3,10)
#ledpred.list=LedPred(data.granges=crm.features, cl=1, ranges = list(cost=cost_vector,
# gamma=gamma_vector), kernel="linear", halve.above=50)
#names(ledpred.list)</pre>
```

mapFeaturesToCRMs R interface to bed\_to\_matrix REST in server

### Description

The mapFeaturesToCRMs function allows the user to create a training set matrix to build a predictive model. The training set is composed of positive regions (known to be involved in the pathway of interest) and negative regions (randomly picked or known to not be involved in the pathway of interest) that will be described (scored) by features. Three types of features file format are accepted: Position specific scoring matrices modeling motifs recognised by transcription factors, bed files containing region coordinates for any discrete feature (NGS peaks, conservation blocks) and wig/bigWig files containing signal data. This script has been tested with version 0.99 of the online server. Go here to see current version of the server http://ifbprod.aitorgonzalezlab.org/map\_features\_to\_crms.php

### Usage

```
mapFeaturesToCRMs(URL = "http://ifbprod.aitorgonzalezlab.org/map_features_to_crms.php",
    positive.bed = NULL, genome = NULL, negative.bed = NULL,
    shuffling = NULL, background.seqs = NULL, genome.info = NULL,
    pssm = NULL, background.freqs = NULL, ngs = NULL, bed.overlap = NULL,
    my.values = NULL, feature.ranking = NULL, feature.nb = NULL,
    crm.feature.file = NULL, stderr.log.file = NULL, stdout.log.file = NULL)
```

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### Arguments

URL	URL of the server REST target	
positive.bed	Positive bed file path. Compulsory	
genome	Genome code, eg. dm3 for Drosophila Melanogaster. Compulsory	
negative.bed	Negative bed file path.	
shuffling	Integer with number of time shuffle background sequences (background.seqs). If negative.bed is NULL and shuffling is set at 0, the feature matrix does not contain negative sequences. It is useful to produce a test set matrix.	
background.seq	5	
	Background sequences used for shuffling. If shuffling = $0$ , set this parameter at $0$ .	
genome.info	File require for shuffling bed. If shuffling = $0$ , set this parameter at $0$ .	
pssm	Position specific scoring matrices	
background.free	ąs	
	Background frequencies of nucleotides in genome	
ngs	NGS (bed and wig) files	
bed.overlap	Minimal overlap as a fraction of query sequence with NGS bed peak. Equivalent with intersectBed -f argument. Default 1bp.	
my.values	Bed file where fourth column are values to append to the SVM matrix	
feature.ranking		
	File with ranked features (Output of rankFeatures). It is used for scoring a query bed file	
feature.nb	Integer with feature.nb	
crm.feature.file		
	Path to feature matrix file	
stderr.log.file		
Path to error log		
<pre>stdout.log.file</pre>	Path to standard output log	

### Value

A list	
--------	--

feature.matrix	a data frame where each row is a region and each column a feature, each cell carry a score, the first column is the response vector
stdout.log	Standard output log of mapFeaturesToCRMs script in server
stderr.log	Standard error log of mapFeaturesToCRMs script in server

### Examples

```
## Not run:
dirPath <- system.file("extdata", package="LedPred")</pre>
 file.list <- list.files(dirPath, full.names=TRUE)</pre>
 background.freqs <- file.list[grep("freq", file.list)]</pre>
 positive.regions <- file.list[grep("positive", file.list)]</pre>
 negative.regions <- file.list[grep("negative", file.list)]</pre>
 TF.matrices <- file.list[grep("tf", file.list)]</pre>
 ngs.path <- system.file("extdata/ngs", package="LedPred")</pre>
 ngs.files=list.files(ngs.path, full.names=TRUE)
 crm.features.list <- mapFeaturesToCRMs(positive.bed=positive.regions,</pre>
     negative.bed=negative.regions, background.freqs=background.freqs,
     pssm=TF.matrices, genome="dm3", ngs=ngs.files,
     crm.feature.file = "crm.features.tab",
     stderr.log.file = "stderr.log", stdout.log.file = "stdout.log")
 names(crm.features.list)
 class(crm.features.list$crm.features)
 crm.features.list$stdout.log
 crm.features.list$stderr.log
```

## End(Not run)

mcTune

Tuning the SVM parameters

### Description

The mcTune function is a modified version of the function tune from package e1071 [6]. It tests the different combinations of C and gamma parameters given as vectors in a list and will return the prediction error computed during the cross-validation step.

### Usage

```
mcTune(data, cl = 1, ranges = list(gamma = c(1, 10), cost = c(1, 10)),
kernel = "linear", valid.times = 10, file.prefix = NULL,
numcores = ifelse(.Platform$0S.type == "windows", 1, parallel::detectCores()
- 1))
```

#### Arguments

data	data.frame containing the training set
cl	integer indicating the column number corresponding to the response vector that classify positive and negative regions (default = $1$ )
ranges	list object containing one (linear kernel) or two (radial kernel) vectors of integers corresponding to SVM cost and SVM gamma parameters to test.
kernel	SVM kernel, a character string: "linear" or "radial". (default = "radial")

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### rankFeatures

valid.times	Integer indicating how many times the training set will be split for the cross validation step (default = 10). This number must be smaller than positive and negative sets sizes.
file.prefix	A character string that will be used as a prefix followed by "_c_g_eval.png" for result plot files, if it is NULL (default), no plot is returned
numcores	Number of cores to use for parallel computing (default: the number of available cores in the machine - 1)

### Value

A list of class tune		
best.parameters		
	A list of the parameters giving the lowest misclassification error	
best.performance		
	The lowest misclassification error	
method	The method used	
nparcomb	the number of tested parameter combinations	
train.ind	The indexes used to produce subsets during the cross validation step	
sampling	The cross-validation fold number	
performances	A matrix summarizing the cross-validation step with the error for each tested parameter at each round and the dispersion of these errors (regarding to the average error)	
best.model	The model produced by the best parameters	

### Examples

```
data(crm.features)
cost.vector <- c(1,3,10,30)
gamma.vector <- c(1,3,10,30)
#c.g.obj <- mcTune(data.granges= crm.features, ranges = list(cost=cost.vector,
# gamma=gamma.vector), kernel='linear', file.prefix = "test")
#names(c.g.obj)
# cost <- c.g.obj$best.parameters$cost
# gamma <- c.g.obj$best.parameters$gamma</pre>
```

rankFeatures

Ranking the features according to their importance

### Description

The rankFeatures function performs a Recursive Feature Elimination (RFE) on subsets of the feature matrix. For each subset the features are ranked according to the weight attributed by SVM at each round of elimination and the average rank of each feature over the subsets is returned. We recommand to save the object containing the ranked features for the following steps.

### Usage

```
rankFeatures(data, cl = 1, halve.above = 100, valid.times = 10,
kernel = "linear", cost = 1, gamma = 1,
numcores = ifelse(.Platform$0S.type == "windows", 1, parallel::detectCores()
- 1), file.prefix = NULL)
```

### Arguments

data	data.frame containing the training set
cl	integer indicating the column number corresponding to the response vector that classify positive and negative regions (default = 1)
halve.above	During RFE, all the features are ranked at the first round and the half lowest ranked features (that contribute the least in the model) are removed for the next round. When the number of feature is lower or equal to halve.above, the features are removed one by one. (default=100)
valid.times	Integer indicating how many times the training set will be split (default = 10). This number must be smaller than positive and negative sets sizes.
kernel	SVM kernel, a character string: "linear" or "radial". (default = "radial")
cost	The SVM cost parameter for both linear and radial kernels. If NULL (default), the function mcTune is run.
gamma	The SVM gamma parameter for radial kernel. If radial kernel and NULL (default), the function mcTune is run.
numcores	Number of cores to use for parallel computing (default: the number of available cores in the machine - 1)
file.prefix	A character string that will be used as a prefix for output file, if it is NULL (default), no file is writen.

### Value

A 3-columns data frame with ranked features. First column contains the feature names, the second the original position of the feature in the feature.matrix and the third the average rank over the subsets.

### Examples

scoreData

### Description

scoreData function predict new regulatory regions using SVM model from a test data set

#### Usage

scoreData(data, ledpred = NULL, model = NULL, score.file = NULL)

### Arguments

data	data.frame containing the test set. This test set must have the same descriptive features as the one that were used to build the model.
ledpred	Returned object from the LedPred function
model	Returned object of the createModel function
score.file	A character string that will be used as the file name for the output file, if it is NULL (default), no file is writen. The output file takes the form of two columns with object names and scores.

### Value

A 2-columns dataframe. First column containg the SVM model prediction probabilities and the second containing the corresponding regions

### Examples

```
data(crm.features)
data(svm.model)
#pred.test <- scoreData(data.granges=crm.features, model=svm.model,
# score.file="test_prediction.tab")</pre>
```

tuneFeatureNb	Selecting the optimal number of features

### Description

tuneFeatureNb iterates through increasing feature numbers to calculate kappa values which represents the performance of the model computed with the given features. We recommand to save the object containing the optimal number of features for the following steps.

```
tuneFeatureNb(data, cl = 1, feature.ranking, step.nb = 10,
valid.times = 10, cost = NULL, gamma = NULL, kernel = "linear",
numcores = ifelse(.Platform$0S.type == "windows", 1, parallel::detectCores()
- 1), file.prefix = NULL)
```

### Arguments

data	data.frame containing the training set	
cl	integer indicating the column number corresponding to the response vector that classify positive and negative regions (default = 1)	
feature.ranking		
	List of ordered features.	
step.nb	Number of features to add at each step (default = $10$ )	
valid.times	Integer indicating how many times the training set will be split for the cross validation step (default = 10). This number must be smaller than positive and negative sets sizes.	
cost	The SVM cost parameter for both linear and radial kernels. If NULL (default), the function mcTune is run.	
gamma	The SVM gamma parameter for radial kernel. If radial kernel and NULL (de-fault), the function mcTune is run.	
kernel	SVM kernel, a character string: "linear" or "radial". (default = "radial")	
numcores	Number of cores to use for parallel computing (default: the number of available cores in the machine - 1)	
file.prefix	A character string that will be used as a prefix followed by "_kappa_measures.png" for the result plot file. If it is NULL (default), no plot is returned	

### Value

A list with two objects.

performance	2-columns data frame. first column correspond to the number of tested features, second column contains the corresponding kappa value
best.feature.nb	
	Integer corresponding to the number of features producing the model with the highest kappa value

### Examples

```
data(crm.features)
data(feature.ranking)
cost <- 1
gamma <- 1
#feature.nb.obj <- tuneFeatureNb(data.granges=crm.features,
# feature.ranking=feature.ranking, kernel='linear', cost=cost,gamma=gamma,
# file.prefix = "test")
#names(feature.nb.obj)</pre>
```

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