

# Package ‘HDPenReg’

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**Copyright** inria

**Depends** R (>= 3.0.2), rtkore (>= 1.5.5)

**Imports** methods, Matrix

**Description** Algorithms for lasso and fused-lasso problems: implementation of the lars algorithm for lasso and fusion penalization and EM-based algorithms for (logistic) lasso and fused-lasso penalization.

**License** GPL (>= 2)

**LinkingTo** rtkore, Rcpp

**SystemRequirements** GNU make

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**NeedsCompilation** yes

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## R topics documented:

HDPenReg-package . . . . .	2
coef.LarsPath . . . . .	3
coeff . . . . .	4
computeCoefficients . . . . .	4
EMcvfusedlasso . . . . .	5
EMcvlasso . . . . .	6
EMfusedlasso . . . . .	8

EMlasso . . . . .	9
HDcvlars . . . . .	10
HDfusion . . . . .	11
HDlars . . . . .	12
LarsPath-class . . . . .	13
listToMatrix . . . . .	14
plot-methods . . . . .	15
plot.HDcvlars . . . . .	15
plotCoefficient . . . . .	16
predict.LarsPath . . . . .	17
simul . . . . .	18

## Index 19

---

HDPenReg-package	<i>Algorithms for lasso and fused-lasso problems.</i>
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## Description

This package contains algorithms for lasso and fused-lasso problems. It contains an implementation of the lars algorithm [1], for the lasso and fusion penalization and EM-based algorithms for (logistic) lasso and fused-lasso.

## Details

Package: HDPenReg  
 Type: Package  
 Version: 0.94.5  
 Date: 2019-03-29  
 License: GPL (>=2)

The main function is [HDlars](#).

## Author(s)

Maintainer: Quentin Grimonprez <quentin.grimonprez@inria.fr>

## See Also

[HDlars](#) [HDcvlars](#)

## Examples

```
## Not run:
#see vignette
vignette("HDPenReg")
```

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

---

coef.LarsPath	<i>Compute coefficients</i>
---------------	-----------------------------

---

### Description

Compute coefficients at a given level of penalty

### Usage

```
## S3 method for class 'LarsPath'
coef(object, index = NULL, mode = c("lambda",
  "step", "fraction", "norm"), ...)
```

### Arguments

object	a LarsParth object
index	If mode="norm", index represents the l1-norm of the coefficients with which we want to predict. If mode="fraction", index represents the ratio (l1-norm of the coefficients with which we want to predict)/(l1-norm maximal of the LarsPath object). If mode="lambda", index represents the value of the penalty parameter. If mode="step", index represents the number of the step at which we want coefficients.
mode	"fraction" or "norm" or "lambda" or "step".
...	other arguments. Not used

### Value

A vector containing the estimated coefficient for index

### Author(s)

Quentin Grimonprez

### See Also

[HDlars](#) [LarsPath](#)

### Examples

```
dataset <- simul(50, 10000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- HDlars(dataset$data[1:40,], dataset$response[1:40])
coeff <- coef(result, 0.3, "fraction")
```

---

coeff	<i>get coefficients at a given step.</i>
-------	--

---

**Description**

Get the vector of coefficients at a given step

**Usage**

```
coeff(x, step)
```

**Arguments**

x	A LarsPath object.
step	The step at which you want to get the coefficients.

**Value**

a vector of size p containing the value of coefficients at the desired step.

**See Also**

[HDLars](#) [Hdfusion](#) [LarsPath](#)

**Examples**

```
dataset <- simul(50, 1000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- Hdfusion(dataset$data, dataset$response)
coefficient <- coeff(result, result@nbStep) #get the coefficients
```

---

computeCoefficients	<i>Compute coefficients</i>
---------------------	-----------------------------

---

**Description**

Compute coefficients at a given level of penalty

**Usage**

```
computeCoefficients(x, lambda, mode = "fraction")
```

**Arguments**

x	a LarsParth object
lambda	If mode = "norm", lambda represents the l1-norm of the coefficients with which we want to predict. If mode="fraction", lambda represents the ratio (l1-norm of the coefficients with which we want to predict)/(l1-norm maximal of the LarsPath object).
mode	"fraction" or "norm" or "lambda".

**Value**

A list containing

**variable** Index of non-zeros coefficients.

**coefficient** non-zeros coefficients.

**Author(s)**

Quentin Grimonprez

**Examples**

```
dataset <- simul(50, 10000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- HDlars(dataset$data[1:40,], dataset$response[1:40])
coeff <- computeCoefficients(result, 0.3, "fraction")
```

---

EMcvfusedlasso

*cross validation for EM fused-lasso*

---

**Description**

cross validation function for [EMfusedlasso](#).

**Usage**

```
EMcvfusedlasso(X, y, lambda1, lambda2, nbFolds = 10, maxSteps = 1000,
  burn = 50, intercept = TRUE, model = c("linear", "logistic"),
  eps = 1e-05, eps0 = 1e-08, epsCG = 1e-08)
```

**Arguments**

X	the matrix (of size n*p) of the covariates.
y	a vector of length n with the response.
lambda1	Values of lambda1 at which prediction error should be computed. Can be a single value.

<code>lambda2</code>	Values of <code>lambda2</code> at which prediction error should be computed. Can be a single value.
<code>nbFolds</code>	the number of folds for the cross-validation.
<code>maxSteps</code>	Maximal number of steps for EM algorithm.
<code>burn</code>	Number of steps for the burn period.
<code>intercept</code>	If TRUE, there is an intercept in the model.
<code>model</code>	"linear" or "logistic".
<code>eps</code>	Tolerance of the algorithm.
<code>eps0</code>	Zero tolerance. Coefficients under this value are set to zero.
<code>epsCG</code>	Epsilon for the convergence of the conjugate gradient.

**Value**

A list containing

**cv** Mean prediction error for each value of index.

**cvError** Standard error of `cv`.

**minCv** Minimal cv criterion.

**lambda1** Values of `lambda1` at which prediction error should be computed.

**lambda2** Values of `lambda2` at which prediction error should be computed.

**lambda.optimal** Value of (`lambda1`,`lambda2`) for which the cv criterion is minimal.

**Author(s)**

Quentin Grimonprez, Serge Iovleff

**Examples**

```
dataset <- simul(50, 100, 0.4, 1, 10, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- EMcvfusedlasso(X = dataset$data, y = dataset$response, lambda1 = 3:1,
                        lambda2 = 3:1, nbFolds = 5, intercept = FALSE)
```

---

EMcvlasso

*cross validation for [EMlasso](#)*

---

**Description**

cross validation function for [EMlasso](#).

**Usage**

```
EMcvlasso(X, y, lambda = NULL, nbFolds = 10, maxSteps = 1000,
          intercept = TRUE, model = c("linear", "logistic"), burn = 30,
          threshold = 1e-08, eps = 1e-05, epsCG = 1e-08)
```

**Arguments**

X	the matrix (of size $n \times p$ ) of the covariates.
y	a vector of length $n$ with the response.
lambda	Values at which prediction error should be computed.
nbFolds	the number of folds for the cross-validation.
maxSteps	Maximal number of steps for EM algorithm.
intercept	If TRUE, there is an intercept in the model.
model	"linear" or "logistic".
burn	Number of steps for the burn period.
threshold	Zero tolerance. Coefficients under this value are set to zero.
eps	Tolerance of the EM algorithm.
epsCG	Epsilon for the convergence of the conjugate gradient.

**Value**

A list containing

**cv** Mean prediction error for each value of index.

**cvError** Standard error of lambda.

**minCv** Minimal lambda criterion.

**lambda** Values of lambda at which prediction error should be computed.

**lambda.optimal** Value of lambda for which the cv criterion is minimal.

**Author(s)**

Quentin Grimonprez, Serge Iovleff

**Examples**

```
dataset <- simul(50, 100, 0.4, 1, 10, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- EMcvlasso(X = dataset$data, y = dataset$response,
                    lambda = 5:1, nbFolds = 5, intercept = FALSE)
```

EMfusedlasso

*EM algorithm for fused-lasso penalty***Description**

EM algorithm for fused-lasso penalty

**Usage**

```
EMfusedlasso(X, y, lambda1, lambda2, maxSteps = 1000, burn = 50,
  intercept = TRUE, model = c("linear", "logistic"), eps = 1e-05,
  eps0 = 1e-08, epsCG = 1e-08)
```

**Arguments**

X	the matrix (of size n*p) of the covariates.
y	a vector of length n with the response.
lambda1	a positive real. Parameter associated with the lasso penalty.
lambda2	a positive real. Parameter associated with the fusion penalty.
maxSteps	Maximal number of steps for EM algorithm.
burn	Number of steps before regrouping some variables in segment.
intercept	If TRUE, there is an intercept in the model.
model	"linear" or "logistic"
eps	tolerance for convergence of the EM algorithm.
eps0	Zero tolerance. Coefficients under this value are set to zero.
epsCG	tolerance for convergence of the conjugate gradient.

**Value**

A list containing :

**step** Vector containing the number of steps of the algorithm for every lambda.

**variable** List of vector of size "step+1". The i+1-th item contains the index of non-zero coefficients at the i-th step.

**coefficient** List of vector of size "step+1". The i+1-th item contains the non-zero coefficients at the i-th step.

**lambda** Vector of length "step+1", containing the lambda at each step.

**mu** Intercept.

**Author(s)**

Quentin Grimonprez, Serge Iovleff



**See Also**[EMcvfusedlasso](#)**Examples**

```
dataset <- simul(50, 100, 0.4, 1, 10, matrix(c(0.1,0.9,0.02,0.02), nrow = 2))
result <- EMfusedlasso(dataset$data, dataset$response, 1, 1)
```

EMlasso

*EM algorithm for lasso penalty***Description**

EM algorithm for lasso penalty

**Usage**

```
EMlasso(X, y, lambda, maxSteps = 1000, intercept = TRUE,
        model = c("linear", "logistic"), burn = 50, threshold = 1e-08,
        eps = 1e-05, epsCG = 1e-08)
```

**Arguments**

X	the matrix (of size n*p) of the covariates.
y	a vector of length n with the response.
lambda	a sequence of l1 penalty regularization term. If no sequence is provided, the function computes his own sequence.
maxSteps	Maximal number of steps for EM algorithm.
intercept	If TRUE, there is an intercept in the model.
model	"linear" or "logistic"
burn	Number of steps before thresholding some variables to zero.
threshold	Zero tolerance. Coefficients under this value are set to zero.
eps	Epsilon for the convergence of the EM algorithm.
epsCG	Epsilon for the convergence of the conjugate gradient.

**Value**

A list containing :

**step** Vector containing the number of steps of the algorithm for every lambda.**variable** List of vector of the same length as lambda. The i-th item contains the index of non-zero coefficients for the i-th lambda value.**coefficient** List of vector of the same length as lambda. The i-th item contains the non-zero coefficients for the i-th lambda value.**lambda** Vector containing the lambda values.**mu** Intercept.

**Author(s)**

Quentin Grimonprez, Serge Iovleff

**See Also**

[EMcvlasso](#)

**Examples**

```
dataset <- simul(50, 100, 0.4, 1, 10, matrix(c(0.1,0.9,0.02,0.02), nrow = 2))
result <- EMLasso(dataset$data, dataset$response)
# Obtain estimated coefficient in matrix format
coefficient <- listToMatrix(result)
```

---

HDcvlars

*cross validation*

---

**Description**

cross validation function for lars algorithm

**Usage**

```
HDcvlars(X, y, nbFolds = 10, index = seq(0, 1, by = 0.01),
  mode = c("fraction", "lambda"), maxSteps = 3 * min(dim(X)),
  partition = NULL, intercept = TRUE, eps = .Machine$double.eps^0.5)
```

**Arguments**

X	the matrix (of size n*p) of the covariates.
y	a vector of length n with the response.
nbFolds	the number of folds for the cross-validation.
index	Values at which prediction error should be computed. When mode = "fraction", this is the fraction of the saturated lbetal. The default value is seq(0,1,by=0.01). When mode="lambda", this is values of lambda.
mode	Either "fraction" or "lambda". Type of values containing in partition.
maxSteps	Maximal number of steps for lars algorithm.
partition	partition in nbFolds folds of y. Must be a vector of same size than y containing the index of folds.
intercept	If TRUE, there is an intercept in the model.
eps	Tolerance of the algorithm.

**Value**

A list containing

**cv** Mean prediction error for each value of index.

**cvError** Standard error of cv.

**minCv** Minimal cv criterion.

**minIndex** Value of index for which the cv criterion is minimal.

**index** Values at which prediction error should be computed. This is the fraction of the saturated  $\beta$ . The default value is `seq(0,1,by=0.01)`.

**maxSteps** Maximum number of steps of the lars algorithm.

**Author(s)**

Quentin Grimonprez

**Examples**

```
dataset <- simul(50, 10000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow =2 ))
result <- HDcvlars(dataset$data, dataset$response, 5)
```

---

Hdfusion

*Fusion algorithm*


---

**Description**

It performs the lars algorithm for solving a special case of lasso problem. It is a linear regression problem with a l1-penalty on the difference of two successive coefficients.

**Usage**

```
Hdfusion(X, y, maxSteps = 3 * min(dim(X)), intercept = TRUE,
eps = .Machine$double.eps^0.5)
```

**Arguments**

X	the matrix (of size n*p) of the covariates.
y	a vector of length n with the response.
maxSteps	Maximal number of steps for lars algorithm.
intercept	If TRUE, there is an intercept in the model.
eps	Tolerance of the algorithm.

**Value**

An object of type `LarsPath`. [LarsPath-class](#).

**Author(s)**

Quentin Grimonprez

**References**

Efron, Hastie, Johnstone and Tibshirani (2003) "Least Angle Regression" (with discussion) *Annals of Statistics*

**See Also**

LarsPath HDlars

**Examples**

```
set.seed(10)
dataset <- simul(50, 10000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- HDFusion(dataset$data, dataset$response)
```

---

HDlars

*Lars algorithm*

---

**Description**

It performs the lars algorithm for solving lasso problem. It is a linear regression problem with a l1-penalty on the estimated coefficient.

**Usage**

```
HDlars(X, y, maxSteps = 3 * min(dim(X)), intercept = TRUE,
      eps = .Machine$double.eps^0.5)
```

**Arguments**

X	the matrix (of size n*p) of the covariates.
y	a vector of length n with the response.
maxSteps	Maximal number of steps for lars algorithm.
intercept	If TRUE, add an intercept to the model.
eps	Tolerance of the algorithm.

**Details**

The l1 penalty performs variable selection via shrinkage of the estimated coefficient. It depends on a penalty parameter called lambda controlling the amount of regularization. The objective function of lasso is :

$$\|y - X\beta\|_2 + \lambda\|\beta\|_1$$

**Value**

An object of type [LarsPath](#).

**Author(s)**

Quentin Grimonprez

**References**

Efron, Hastie, Johnstone and Tibshirani (2003) "Least Angle Regression" (with discussion) *Annals of Statistics*

**See Also**

[LarsPath](#) [HDCvlars](#) [listToMatrix](#)

**Examples**

```
dataset <- simul(50, 10000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- HDlars(dataset$data, dataset$response)
# Obtain estimated coefficient in matrix format
coefficient <- listToMatrix(result)
```

---

LarsPath-class

*Constructor of LarsPath class*

---

**Description**

This class stores the results of lars and fusion algorithms.

**Details**

**nbStep** Number of steps of the algorithm.

**variable** List of vector of size "step+1". The i+1-th item contains the index of non-zero coefficients at the i-th step.

**coefficient** List of vector of size "step+1". The i+1-th item contains the non-zero coefficients at the i-th step.

**l1norm** Vector of length "step+1", containing the L1-norm of the coefficients at each step.

**lambda** Vector of length "step+1", containing the lambda at each step.

**dropIndex** Vector of length "step" containing the index of the dropped variable at the i-th step, 0 means no variable has been dropped at this step.

**addIndex** Vector of length "step" containing the index of the added variable at the i-th step, 0 means no variable has been added at this step.

**mu** Intercept.

**meanX** Mean of columns of X.

**ignored** A vector containing index of ignored variables during the algorithm.

**p** Total number of covariates.

**fusion** If TRUE, results from HDfusion function.

**error** Error message from lars.

### See Also

[HDlars](#)

---

listToMatrix	<i>List to sparse matrix conversion</i>
--------------	---

---

### Description

create a matrix with all estimated coefficients from the output of [HDlars](#) or [EMlasso](#) functions.

### Usage

```
listToMatrix(x, row = c("covariates", "lambda"))
```

### Arguments

x	a <a href="#">LarsPath</a> or <a href="#">EMlasso</a> object
row	if covariates, covariates are in row

### Value

A sparse matrix containing the values of estimated coefficients for all penalty parameter and all covariates

### See Also

[HDlars](#) [EMlasso](#)

---

plot-methods	<i>plot methods for LarsPath object</i>
--------------	---

---

**Description**

plot the path of the lars algorithm.

**Usage**

```
## S4 method for signature 'LarsPath'
plot(x, sep.line = FALSE, abscissa = c("l1norm",
  "lambda"), log.scale = FALSE, ...)
```

**Arguments**

x	LarsPath object
sep.line	If TRUE, print vertical dashed line when a variable is added or dropped in the path
abscissa	either "l1norm" or "lambda". If "lambda", regularization parameter is used as abscissa, else l1 norm of the solution is used.
log.scale	If TRUE, use logarithm scale on abscissa
...	Other plot arguments

**See Also**

[HDlars LarsPath](#)

---

plot.HDcvlars	<i>plot cross validation mean square error</i>
---------------	--

---

**Description**

plot cross validation mean square error

**Usage**

```
## S3 method for class 'HDcvlars'
plot(x, ...)
```

**Arguments**

x	Output from HDcvlars function.
...	graphical parameters

**Author(s)**

Quentin Grimonprez

**Examples**

```
dataset <- simul(50, 10000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- HDcvlars(dataset$data, dataset$response, 5)
plot(result)
```

---

plotCoefficient	<i>Plot of coefficients</i>
-----------------	-----------------------------

---

**Description**

Plot of the coefficients of a step

**Usage**

```
plotCoefficient(x, step, ylab = "coefficients", xlab = "variables",
  ...)
```

**Arguments**

x	A LarsPath object.
step	The step at which you want to plot the coefficients.
ylab	Name of the y axis.
xlab	Name of the x axis.
...	Other plot arguments.

**See Also**

[HDlars LarsPath](#)

**Examples**

```
dataset <- simul(50, 1000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))
result <- HDfusion(dataset$data, dataset$response)
plotCoefficient(result, result@nbStep) #plot coefficients at the last step
```



---

predict.LarsPath      *Prediction of response*

---

### Description

Predict response of a new sample Xnew at a given level of penalty

### Usage

```
## S3 method for class 'LarsPath'  
predict(object, Xnew, lambda, mode = c("fraction",  
    "lambda", "norm"), ...)
```

### Arguments

object	a LarsParth object
Xnew	a matrix (of size n*object@p) of covariates.
lambda	If mode = "norm", lambda represents the l1-norm of the coefficients with which we want to predict. If mode="fraction", lambda represents the ratio (l1-norm of the coefficients with which we want to predict)/(l1-norm maximal of the LarsPath object).
mode	"fraction", "lambda" or "norm".
...	other arguments. Not used.

### Value

The predicted response

### Author(s)

Quentin Grimonprez

### Examples

```
dataset <- simul(50, 10000, 0.4, 10, 50, matrix(c(0.1,0.8,0.02,0.02), nrow = 2))  
result <- HDlars(dataset$data[1:40,], dataset$response[1:40])  
y <- predict(result, dataset$data[41:50,], 0.3, "fraction")
```

---

simul *Simulate copy number data for a case-control study.*

---

### Description

Simulate copy number data for a case-control study.

### Usage

```
simul(n, nbSNP, probCas, nbSeg, meanSegmentSize, prob, alpha = 15)
```

### Arguments

n	Number of individuals.
nbSNP	Size of the DNA sequence.
probCas	Probability to be a case individual.
nbSeg	Number of causal segments.
meanSegmentSize	The mean size of anormal segment.
prob	A 2*2 matrix containing probabilities: prob[1,1]=probability to have an anomaly to a SNP given the person does not have the disease and the SNP is causal. prob[1,2]=probability to have an anomaly to a SNP given the person does not have the disease and the SNP is not causal. prob[2,1]=probability to have an anomaly to a SNP given the person has the disease and the SNP is causal. prob[2,2]=probability to have an anomaly to a SNP given the person has the disease and the SNP is not causal.
alpha	Parameter of the beta(alpha,alpha).

### Value

a list containing:

data	A matrix of size n*nbSeg, containing values of the copy-number signal.
response	A vector of size n containing the cas/control status.
causalSNP	A vector of size nbSeg containing the center of causal segments.

### Author(s)

Quentin Grimonprez, Serge Iovleff

### Examples

```
data <- simul(50, 10000, 0.4, 10, 150, matrix(c(0.1,0.8,0.001,0.001), nrow = 2))
```

# Index

## \*Topic **package**

HDPenReg-package, 2

coef.LarsPath, 3

coeff, 4

computeCoefficients, 4

EMcvfusedlasso, 5, 9

EMcvlasso, 6, 10

EMfusedlasso, 5, 8

EMlasso, 6, 9, 14

HDcvlars, 2, 10, 13

HDfusion, 4, 11

HDlars, 2–4, 12, 14–16

HDPenReg (HDPenReg-package), 2

HDPenReg-package, 2

HDPenReg-package, (HDPenReg-package), 2

LarsPath, 3, 4, 11, 13–16

LarsPath (LarsPath-class), 13

LarsPath-class, 13

listToMatrix, 13, 14

plot, LarsPath-method (plot-methods), 15

plot-methods, 15

plot.HDcvlars, 15

plotCoefficient, 16

predict.LarsPath, 17

simul, 18