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Title Family of Lasso Regression

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Imports methods

Description Provide the implementation of a family of Lasso variants including Dantzig Selector, LAD Lasso, SQRT Lasso, Lq Lasso for estimating high dimensional sparse linear model. We adopt the alternating direction method of multipliers and convert the original optimization problem into a sequential L1 penalized least square minimization problem, which can be efficiently solved by linearization algorithm. A multi-stage screening approach is adopted for further acceleration. Besides the sparse linear model estimation, we also provide the extension of these Lasso variants to sparse Gaussian graphical model estimation including TIGER and CLIME using either L1 or adaptive penalty. Missing values can be tolerated for Dantzig selector and CLIME. The computation is memory-optimized using the sparse matrix output. For more information, please refer to <<https://www.jmlr.org/papers/volume16/li15a/li15a.pdf>>.

License GPL-2

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flare-package	<i>flare: a new Family of Lasso Regression</i>
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Description

The package "flare" provides the implementation of a family of novel regression methods (Lasso, Dantzig Selector, LAD Lasso, SQRT Lasso, Lq Lasso) and their extensions to sparse precision matrix estimation (TIGER and CLIME using L1) in high dimensions. We adopt the alternating direction method of multipliers and convert the original optimization problem into a sequence of L1-penalized least square minimization problems with the linearization method and multi-stage screening of variables. Missing values can be tolerated for Dantzig selector in the design matrix and response vector, and CLIME in the data matrix. The computation is memory-optimized using the sparse matrix output. In addition, we also provide several convenient regularization parameter selection and visualization tools.

Details

Package:	flare
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Version:	1.7.0
Date:	2020-11-28

License: GPL-2

Author(s)

Xingguo Li, Tuo Zhao, Lie Wang , Xiaoming Yuan and Han Liu
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References

1. E. Candes and T. Tao. The Dantzig selector: Statistical estimation when p is much larger than n . *Annals of Statistics*, 2007.
2. A. Belloni, V. Chernozhukov and L. Wang. Pivotal recovery of sparse signals via conic programming. *Biometrika*, 2012.
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7. H. Liu and L. Wang. TIGER: A tuning-insensitive approach for optimally estimating large undirected graphs. *Technical Report*, 2012.
8. B. He and X. Yuan. On non-ergodic convergence rate of Douglas-Rachford alternating direction method of multipliers. *Technical Report*, 2012.

See Also

[sugm](#) and [slim](#).

 coef.slim

Extract Model Coefficients for an object with S3 class "slim"

Description

Extract estimated regression coefficient vectors from the solution path.

Usage

```
## S3 method for class 'slim'
coef(object, lambda.idx = c(1:3), beta.idx = c(1:3), ...)
```

Arguments

object	An object with S3 class "slim"
lambda.idx	The indices of the regularization parameters in the solution path to be displayed. The default values are <code>c(1:3)</code> .
beta.idx	The indices of the estimate regression coefficient vectors in the solution path to be displayed. The default values are <code>c(1:3)</code> .
...	Arguments to be passed to methods.

Author(s)

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See Also

[slim](#) and [flare-package](#).

eyedata	<i>The Bardet-Biedl syndrome Gene expression data from Scheetz et al. (2006)</i>
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Description

Gene expression data (20 genes for 120 samples) from the microarray experiments of mammalian-eye tissue samples of Scheetz et al. (2006).

Usage

```
data(eyedata)
```

Format

The format is a list containing contains a matrix and a vector. 1. `x` - an 120 by 200 matrix, which represents the data of 120 rats with 200 gene probes. 2. `y` - a 120-dimensional vector of, which represents the expression level of TRIM32 gene.

Details

This data set contains 120 samples with 200 predictors

Author(s)

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References

1. T. Scheetz, k. Kim, R. Swiderski, A. Philp, T. Braun, K. Knudtson, A. Dorrance, G. DiBona, J. Huang, T. Casavant, V. Sheffield, E. Stone .Regulation of gene expression in the mammalian eye and its relevance to eye disease. *Proceedings of the National Academy of Sciences of the United States of America*, 2006.

See Also

[flare-package](#).

Examples

```
data(eyedata)
image(x)
```

flare-internal	<i>Internal flare functions</i>
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Description

Internal flare functions

Usage

```
sugm.likelihood(Sigma, Omega)
sugm.tracel2(Sigma, Omega)
sugm.cv(obj, loss=c("likelihood", "tracel2"), fold=5)
part.cv(n, fold)
sugm.clime.ladm.scr(Sigma, lambda, nlambda, n, d, maxdf, rho, shrink, prec,
                    max.ite, verbose)
sugm.tiger.ladm.scr(data, n, d, maxdf, rho, lambda, shrink, prec,
                    max.ite, verbose)
slim.lad.ladm.scr.btr(Y, X, lambda, nlambda, n, d, maxdf, rho, max.ite, prec,
                     intercept, verbose)
slim.sqrt.ladm.scr(Y, X, lambda, nlambda, n, d, maxdf, rho, max.ite, prec,
                   intercept, verbose)
slim.dantzig.ladm.scr(Y, X, lambda, nlambda, n, d, maxdf, rho, max.ite, prec,
                     intercept, verbose)
slim.lq.ladm.scr.btr(Y, X, q, lambda, nlambda, n, d, maxdf, rho, max.ite, prec,
                    intercept, verbose)
slim.lasso.ladm.scr(Y, X, lambda, nlambda, n, d, maxdf, max.ite, prec,
                   intercept, verbose)
```

Arguments

<code>Sigma</code>	Covariance matrix.
<code>Omega</code>	Inverse covariance matrix.
<code>obj</code>	An object with S3 class returned from "sugm".
<code>loss</code>	Type of loss function for cross validation.
<code>fold</code>	The number of fold for cross validatio.
<code>n</code>	The number of observations (sample size).
<code>d</code>	Dimension of data.
<code>maxdf</code>	Maximal degree of freedom.
<code>lambda</code>	Grid of non-negative values for the regularization parameter lambda.
<code>nlambda</code>	The number of the regularization parameter lambda.
<code>shrink</code>	Shrinkage of regularization parameter based on precision of estimation.
<code>rho</code>	Value of augmented Lagrangian multipiler.
<code>prec</code>	Stopping criterion.
<code>max.ite</code>	Maximal value of iterations.
<code>data</code>	n by d data matrix.
<code>Y</code>	Dependent variables in linear regression.
<code>X</code>	Design matrix in linear regression.
<code>q</code>	The vector norm used for the loss term.
<code>intercept</code>	The indicator of whether including intercepts specifically.
<code>verbose</code>	Tracing information printing is disabled if <code>verbose = FALSE</code> . The default value is <code>TRUE</code> .

Details

These are not intended for use by users.

Author(s)

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See Also

[sugm](#), [slim](#) and [flare-package](#).

plot.roc *Plot Function for "roc"*

Description

Plot the ROC curve for an object with S3 class "roc"

Usage

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

Arguments

x An object with S3 class "roc"
... System reserved (No specific usage)

Author(s)

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See Also

[sugm.roc](#), [sugm](#) and [flare-package](#).

plot.select *Plot Function for "select"*

Description

Plot the optimal graph by model selection.

Usage

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

Arguments

x An object with S3 class "select"
... System reserved (No specific usage)

Author(s)

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See Also

[sugm](#) and [sugm.select](#)

plot.sim

Plot Function for "sim"

Description

Visualize the covariance matrix, the empirical covariance matrix, the adjacency matrix and the graph pattern of the true graph structure.

Usage

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

Arguments

x	An object with S3 class "sim"
...	Arguments to be passed to methods.

Author(s)

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See Also

[sugm.generator](#), [sugm](#) and [flare-package](#)

plot.slim	<i>Plot Function for "slim"</i>
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Description

Visualize the solution path of regression estimate corresponding to regularization parameters.

Usage

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

Arguments

x	An object with S3 class "slim".
...	Arguments to be passed to methods.

Author(s)

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See Also

[slim](#) and [flare-package](#).

plot.sugm	<i>Plot Function for "sugm"</i>
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Description

Plot sparsity level information and 3 typical sparse graphs from the graph path.

Usage

```
## S3 method for class 'sugm'  
plot(x, align = FALSE, ...)
```

Arguments

x	An object with S3 class "sugm"
align	If align = FALSE, 3 plotted graphs are aligned
...	Arguments to be passed to methods.

Author(s)

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See Also

[sugm](#) and [flare-package](#)

predict.slim

Prediction for an object with S3 class "slim"

Description

Predicting responses of the given design data.

Usage

```
## S3 method for class 'slim'
predict(object, newdata, lambda.idx = c(1:3), Y.pred.idx = c(1:5), ...)
```

Arguments

object	An object with S3 class "slim"
newdata	An optional data frame in which to look for variables with which to predict. If omitted, the training data of the are used.
lambda.idx	The indices of the regularization parameters in the solution path to be displayed. The default values are c(1:3).
Y.pred.idx	The indices of the predicted response vectors in the solution path to be displayed. The default values are c(1:5).
...	Arguments to be passed to methods.

Details

predict.slim produces predicted values of the responses of the newdata from the estimated beta values in the object, i.e.

$$\hat{Y} = \hat{\beta}_0 + X_{new}\hat{\beta}.$$

Value

Y.pred The predicted response vectors based on the estimated models.

Author(s)

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See Also

[slim](#) and [flare-package](#).

Examples

```
## load library
library(flare)
## generate data
set.seed(123)
n = 100
d = 200
d1 = 10
rho0 = 0.3
lambda = c(3:1)*sqrt(log(d)/n)
Sigma = matrix(0,nrow=d,ncol=d)
Sigma[1:d1,1:d1] = rho0
diag(Sigma) = 1
mu = rep(0,d)
X = mvrnorm(n=2*n,mu=mu,Sigma=Sigma)
X.fit = X[1:n,]
X.pred = X[(n+1):(2*n),]
eps = rt(n=n,df=n-1)
beta = c(rep(sqrt(1/3),3),rep(0,d-3))
Y.fit = X.fit%%beta+eps

## Regression with "dantzig".
out=slim(X=X.fit,Y=Y.fit,lambda=lambda,method = "lq",q=1)

## Display results
Y=predict(out,X.pred)
```

print.roc

Print Function for for an object with S3 class "roc"

Description

Print the information about true positive rates, false positive rates, the area under curve and maximum F1 score

Usage

```
## S3 method for class 'roc'
print(x, ...)
```

Arguments

x An object with S3 class "roc"
... Arguments to be passed to methods.

Author(s)

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See Also

[sugm.roc](#), [sugm](#) and [flare-package](#)

print.select

Print Function for for an object with S3 class "select"

Description

Print the information about the model usage, graph dimension, model selection criterion, sparsity level of the optimal graph

Usage

```
## S3 method for class 'select'  
print(x, ...)
```

Arguments

x	An object with S3 class "select"
...	Arguments to be passed to methods.

Author(s)

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See Also

[sugm.select](#), [sugm](#) and [flare-package](#)

print.sim	<i>Print Function for for an object with S3 class "sim"</i>
-----------	---

Description

Print the information about the sample size, the dimension, the pattern and sparsity of the true graph structure.

Usage

```
## S3 method for class 'sim'  
print(x, ...)
```

Arguments

x	An object with S3 class "sim".
...	Arguments to be passed to methods.

Author(s)

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See Also

[sugm](#) and [sugm.generator](#)

print.slim	<i>Print Function for an object with S3 class "slim"</i>
------------	--

Description

Print a summary of the information about an object with S3 class "slim".

Usage

```
## S3 method for class 'slim'  
print(x, ...)
```

Arguments

x	An object with S3 class "slim".
...	Arguments to be passed to methods.

Details

This call simply outlines the options used for computing a slim object.

Author(s)

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See Also

[slim](#) and [flare-package](#).

print.sugm

Print Function for an object with S3 class "sugm"

Description

Print a summary of the information about an object with S3 class "slim".

Usage

```
## S3 method for class 'sugm'  
print(x, ...)
```

Arguments

x	An object with S3 class "sugm".
...	Arguments to be passed to methods.

Details

This call simply outlines the options used for computing a sugm object.

Author(s)

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See Also

[sugm](#) and [flare-package](#).

slim *Sparse Linear Regression using Nonsmooth Loss Functions and L1 Regularization*

Description

The function "slim" implements a family of Lasso variants for estimating high dimensional sparse linear models including Dantzig Selector, LAD Lasso, SQRT Lasso, Lq Lasso for estimating high dimensional sparse linear model. We adopt the alternating direction method of multipliers (ADMM) and convert the original optimization problem into a sequential L1-penalized least square minimization problem, which can be efficiently solved by combining the linearization and multi-stage screening of variables. Missing values can be tolerated for Dantzig selector in the design matrix and response vector.

Usage

```
slim(X, Y, lambda = NULL, nlambda = NULL,
     lambda.min.value = NULL, lambda.min.ratio = NULL,
     rho = 1, method="lq", q = 2, res.sd = FALSE,
     prec = 1e-5, max.ite = 1e5, verbose = TRUE)
```

Arguments

Y	The n -dimensional response vector.
X	The n by d design matrix. $d \geq 2$ is required.
lambda	A sequence of decreasing positive numbers to control the regularization. Typical usage is to leave the input <code>lambda = NULL</code> and have the program compute its own <code>lambda</code> sequence based on <code>nlambda</code> and <code>lambda.min.ratio</code> . Users can also specify a sequence to override this. Default value is from <code>lambda.max</code> to <code>lambda.min.ratio*lambda.max</code> . For Lq regression, the default value of <code>lambda.max</code> is $\pi \sqrt{\log(d)/n}$. For Dantzig selector, the default value of <code>lambda.max</code> is the minimum regularization parameter, which yields an all-zero estimates.
nlambda	The number of values used in <code>lambda</code> . Default value is 5.
lambda.min.value	The smallest value for <code>lambda</code> , as a fraction of the upperbound (<code>lambda.max</code>) of the regularization parameter. The program can automatically generate <code>lambda</code> as a sequence of length = <code>nlambda</code> starting from <code>lambda.max</code> to <code>lambda.min.ratio*lambda.max</code> in log scale. The default value is $\log(d)/n$ for for Dantzig selector $0.3 * \lambda.max$ for Lq Lasso.
lambda.min.ratio	The smallest ratio of the value for <code>lambda</code> . The default value is 0.3 for Lq Lasso and 0.5 for Dantzig selector.
rho	The penalty parameter used in ADMM. The default value is \sqrt{d} .
method	Dantzig selector is applied if <code>method = "dantzig"</code> and L_q Lasso is applied if <code>method = "lq"</code> . Standard Lasso is provided if <code>method = "lasso"</code> . The default value is "lq".

q	The loss function used in Lq Lasso. It is only applicable when method = "lq" and must be in [1,2]. The default value is 2.
res.sd	Flag of whether the response variables are standardized. The default value is FALSE.
prec	Stopping criterion. The default value is 1e-5.
max.ite	The iteration limit. The default value is 1e5.
verbose	Tracing information printing is disabled if verbose = FALSE. The default value is TRUE.

Details

Standard Lasso

$$\min \frac{1}{2n} \|Y - X\beta\|_2^2 + \lambda \|\beta\|_1$$

Dantzig selector solves the following optimization problem

$$\min \|\beta\|_1, \quad \text{s.t. } \|X'(Y - X\beta)\|_\infty < \lambda$$

L_q loss Lasso solves the following optimization problem

$$\min n^{-\frac{1}{q}} \|Y - X\beta\|_q + \lambda \|\beta\|_1$$

where $1 \leq q \leq 2$. Lq Lasso is equivalent to LAD Lasso and SQR Lasso when $q = 1$ and $q = 2$ respectively.

Value

An object with S3 class "slim" is returned:

beta	A matrix of regression estimates whose columns correspond to regularization parameters.
intercept	The value of intercepts corresponding to regularization parameters.
Y	The value of Y used in the program.
X	The value of X used in the program.
lambda	The sequence of regularization parameters lambda used in the program.
nlambda	The number of values used in lambda.
method	The method from the input.
sparsity	The sparsity levels of the solution path.
ite	A list of vectors where ite[[1]] is the number of external iteration and ite[[2]] is the number of internal iteration with the i-th entry corresponding to the i-th regularization parameter.
verbose	The verbose from the input.

Author(s)

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References

1. E. Candes and T. Tao. The Dantzig selector: Statistical estimation when p is much larger than n . *Annals of Statistics*, 2007.
2. A. Belloni, V. Chernozhukov and L. Wang. Pivotal recovery of sparse signals via conic programming. *Biometrika*, 2012.
3. L. Wang. L1 penalized LAD estimator for high dimensional linear regression. *Journal of Multivariate Analysis*, 2012.
4. J. Liu and J. Ye. Efficient L1/Lq Norm Regularization. *Technical Report*, 2010.
5. S. Boyd, N. Parikh, E. Chu, B. Peleato, and J. Eckstein, Distributed Optimization and Statistical Learning via the Alternating Direction Method of Multipliers. *Foundations and Trends in Machine Learning*, 2011.
6. B. He and X. Yuan. On non-ergodic convergence rate of Douglas-Rachford alternating direction method of multipliers. *Technical Report*, 2012.

See Also

[flare-package](#), [print.slim](#), [plot.slim](#), [coef.slim](#) and [predict.slim](#).

Examples

```
## load library
library(flare)
## generate data
n = 50
d = 100
X = matrix(rnorm(n*d), n, d)
beta = c(3,2,0,1.5,rep(0,d-4))
eps = rnorm(n)
Y = X%%beta + eps
nlamb = 5
ratio = 0.3

## Regression with "dantzig", general "lq" and "lasso" respectively
out1 = slim(X=X,Y=Y,nlamb=nlamb,lambda.min.ratio=ratio,method="dantzig")
out2 = slim(X=X,Y=Y,nlamb=nlamb,lambda.min.ratio=ratio,method="lq",q=1)
out3 = slim(X=X,Y=Y,nlamb=nlamb,lambda.min.ratio=ratio,method="lq",q=1.5)
out4 = slim(X=X,Y=Y,nlamb=nlamb,lambda.min.ratio=ratio,method="lq",q=2)
out5 = slim(X=X,Y=Y,nlamb=nlamb,lambda.min.ratio=ratio,method="lasso")

## Display results
print(out4)
plot(out4)
coef(out4)
```

Description

The function "sugm" estimates sparse undirected graphical models, i.e. Gaussian precision matrix, in high dimensions. We adopt two estimation procedures based on column by column regression scheme: (1) Tuning-Insensitive Graph Estimation and Regression based on square root Lasso (tiger); (2) The Constrained L1 Minimization for Sparse Precision Matrix Estimation using either L1 penalty (clime). The optimization algorithm for all three methods are implemented based on the alternating direction method of multipliers (ADMM) with the linearization method and multi-stage screening of variables. Missing values can be tolerated for CLIME in the data matrix. The computation is memory-optimized using the sparse matrix output.

Usage

```
sugm(data, lambda = NULL, nlambda = NULL, lambda.min.ratio = NULL,
      rho = NULL, method = "tiger", sym = "or", shrink=NULL,
      prec = 1e-4, max.ite = 1e4, standardize = FALSE,
      perturb = TRUE, verbose = TRUE)
```

Arguments

data	There are 2 options for "clime": (1) data is an n by d data matrix (2) a d by d sample covariance matrix. The program automatically identifies the input matrix by checking the symmetry. (n is the sample size and d is the dimension). For "tiger", covariance input is not supported and $d \geq 3$ is required. For "clime", $d \geq 2$ is required.
lambda	A sequence of decreasing positive numbers to control the regularization. Typical usage is to leave the input lambda = NULL and have the program compute its own lambda sequence based on nlambda and lambda.min.ratio. Users can also specify a sequence to override this. Default value is from lambda.max to lambda.min.ratio*lambda.max. For "tiger", the default value of lambda.max is $\pi \sqrt{\log(d)/n}$. For "clime", the default value of lambda.max is the minimum regularization parameter, which yields an all-zero off-diagonal estimates.
nlambda	The number of values used in lambda. Default value is 5.
lambda.min.ratio	The smallest value for lambda, as a fraction of the upperbound (lambda.max) of the regularization parameter. The program can automatically generate lambda as a sequence of length = nlambda starting from lambda.max to lambda.min.ratio*lambda.max in log scale. The default value is 0.25 for "tiger" and 0.5 for "clime".
rho	Penalty parameter used in the optimization algorithm for clime. The default value is \sqrt{d} .
method	"tiger" is applied if method = "tiger" and "clime" is applied if method="clime". Default value is "tiger".

sym	Symmetrization of output graphs. If sym = "and", the edge between node i and node j is selected ONLY when both node i and node j are selected as neighbors for each other. If sym = "or", the edge is selected when either node i or node j is selected as the neighbor for each other. The default value is "or".
shrink	Shrinkage of regularization parameter based on precision of estimation. The default value is 1.5 if method = "clime" and the default value is 0 if method="tiger".
prec	Stopping criterion. The default value is 1e-4.
max.ite	The iteration limit. The default value is 1e4.
standardize	Variables are standardized to have mean zero and unit standard deviation if standardize = TRUE. The default value is FALSE.
perturb	The diagonal of Sigma is added by a positive value to guarantee that Sigma is positive definite if perturb = TRUE. User can specify a numeric value for perturbe. The default value is perturb = TRUE.
verbose	Tracing information printing is disabled if verbose = FALSE. The default value is TRUE.

Details

CLIME solves the following minimization problem

$$\min \|\Omega\|_1 \quad \text{s.t.} \quad \|S\Omega - I\|_\infty \leq \lambda,$$

where $\|\cdot\|_1$ and $\|\cdot\|_\infty$ are element-wise 1-norm and ∞ -norm respectively.

"tiger" solves the following minimization problem

$$\min \|X - XB\|_{2,1} + \lambda\|B\|_1 \quad \text{s.t.} \quad B_{jj} = 0,$$

where $\|\cdot\|_1$ and $\|\cdot\|_{2,1}$ are element-wise 1-norm and $L_{2,1}$ -norm respectively.

Value

An object with S3 class "sugm" is returned:

data	The n by d data matrix or d by d sample covariance matrix from the input.
cov.input	An indicator of the sample covariance.
lambda	The sequence of regularization parameters lambda used in the program.
nlambda	The number of values used in lambda.
icov	A list of d by d precision matrices corresponding to regularization parameters.
sym	The sym from the input.
method	The method from the input.
path	A list of d by d adjacency matrices of estimated graphs as a graph path corresponding to lambda.
sparsity	The sparsity levels of the graph path.

ite	If method = "clime", it is a list of two matrices where ite[[1]] is the number of external iterations and ite[[2]] is the number of internal iterations with the entry of (i,j) as the number of iteration of i-th column and j-th lambda. If method="tiger", it is a matrix of iteration with the entry of (i,j) as the number of iteration of i-th column and j-th lambda.
df	It is a d by nlambda matrix. Each row contains the number of nonzero coefficients along the lasso solution path.
standardize	The standardize from the input.
perturb	The perturb from the input.
verbose	The verbose from the input.

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References

1. T. Cai, W. Liu and X. Luo. A constrained L1 minimization approach to sparse precision matrix estimation. *Journal of the American Statistical Association*, 2011.
2. H. Liu, L. Wang. TIGER: A tuning-insensitive approach for optimally estimating large undirected graphs. *Technical Report*, 2012.
3. B. He and X. Yuan. On non-ergodic convergence rate of Douglas-Rachford alternating direction method of multipliers. *Technical Report*, 2012.

See Also

[flare-package](#), [sugm.generator](#), [sugm.select](#), [sugm.plot](#), [sugm.roc](#), [plot.sugm](#), [plot.select](#), [plot.roc](#), [plot.sim](#), [print.sugm](#), [print.select](#), [print.roc](#) and [print.sim](#).

Examples

```
## load package required
library(flare)

## generating data
n = 50
d = 50
D = sugm.generator(n=n,d=d,graph="band",g=1)
plot(D)

## sparse precision matrix estimation with method "clime"
out1 = sugm(D$data, method = "clime")
plot(out1)
sugm.plot(out1$path[[4]])

## sparse precision matrix estimation with method "tiger"
out2 = sugm(D$data, method = "tiger")
plot(out2)
sugm.plot(out2$path[[5]])
```

sugm.generator *Data generator for sparse undirected graph estimation.*

Description

Implements the data generation from multivariate normal distributions with different graph structures, including "random", "hub", "cluster", "band", and "scale-free".

Usage

```
sugm.generator(n = 200, d = 50, graph = "random", v = NULL, u = NULL,
              g = NULL, prob = NULL, seed = NULL, vis = FALSE, verbose = TRUE)
```

Arguments

n	The number of observations (sample size). The default value is 200.
d	The number of variables (dimension). For "hub" and "cluster", $d \geq 4$ is required. For "random", "band" and "scale-free", $d \geq 3$ is required. The default value is 50.
graph	The graph structure with 5 options: "random", "hub", "cluster", "band", and "scale-free".
v	The off-diagonal elements of the precision matrix, controlling the magnitude of partial correlations with u. The default value is 0.3.
u	A positive number being added to the diagonal elements of the precision matrix, to control the magnitude of partial correlations. The default value is 0.1.
g	For "cluster" or "hub" graph, g is the number of hubs or clusters in the graph. The default value is about $d/20$ if $d \geq 40$ and 2 if $d < 40$. For "band" graph, g is the bandwidth and the default value is 1. NOT applicable to "random" graph.
prob	For "random" graph, it is the probability that a pair of nodes has an edge. The default value is $3/d$. For "cluster" graph, it is the probability that a pair of nodes has an edge in each cluster. The default value is $6*g/d$ if $d/g \leq 30$ and 0.3 if $d/g > 30$. NOT applicable to "hub", "band", and "scale-free" graphs.
seed	Set seed for data generation. The default value is 1.
vis	Visualize the adjacency matrix of the true graph structure, the graph pattern, the covariance matrix and the empirical covariance matrix. The default value is FALSE.
verbose	If verbose = FALSE, tracing information printing is disabled. The default value is TRUE.

Details

Given the adjacency matrix θ , the graph patterns are generated as below:

(I) "random": Each pair of off-diagonal elements are randomly set $\theta[i, j] = \theta[j, i] = 1$

for $i \neq j$ with probability prob , and 0 otherwise. It results in about $d \cdot (d-1) \cdot \text{prob} / 2$ edges in the graph.

(II) "hub": The row/columns are evenly partitioned into g disjoint groups. Each group is associated with a "center" row i in that group. Each pair of off-diagonal elements are set $\text{theta}[i, j] = \text{theta}[j, i] = 1$ for $i \neq j$ if j also belongs to the same group as i and 0 otherwise. It results in $d - g$ edges in the graph.

(III) "cluster": The row/columns are evenly partitioned into g disjoint groups. Each pair of off-diagonal elements are set $\text{theta}[i, j] = \text{theta}[j, i] = 1$ for $i \neq j$ with the probability prob both i and j belong to the same group, and 0 otherwise. It results in about $g \cdot (d/g) \cdot (d/g - 1) \cdot \text{prob} / 2$ edges in the graph.

(IV) "band": The off-diagonal elements are set to be $\text{theta}[i, j] = 1$ if $1 \leq |i - j| \leq g$ and 0 otherwise. It results in $(2d - 1 - g) \cdot g / 2$ edges in the graph.

(V) "scale-free": The graph is generated using B-A algorithm. The initial graph has two connected nodes and each new node is connected to only one node in the existing graph with the probability proportional to the degree of the each node in the existing graph. It results in d edges in the graph.

The adjacency matrix theta has all diagonal elements equal to 0 . To obtain a positive definite covariance matrix, the smallest eigenvalue of $\text{theta} \cdot v$ (denoted by e) is computed. Then we set the covariance matrix equal to $\text{cov2cor}(\text{solve}(\text{theta} \cdot v + (|e| + 0.1 + u) \cdot I))$ to generate multivariate normal data.

Value

An object with S3 class "sim" is returned:

<code>data</code>	The n by d matrix for the generated data
<code>sigma</code>	The covariance matrix for the generated data
<code>omega</code>	The precision matrix for the generated data
<code>sigmahat</code>	The empirical covariance matrix for the generated data
<code>theta</code>	The adjacency matrix of true graph structure (in sparse matrix representation) for the generated data

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See Also

[flare](#) and [flare-package](#)

Examples

```

## load package required
library(flare)

## band graph with bandwidth 3
L = sugm.generator(graph = "band", g = 3)
plot(L)

## random sparse graph
L = sugm.generator(vis = TRUE)

## hub graph with 6 hubs
L = sugm.generator(graph = "hub", g = 6, vis = TRUE)

## cluster graph with 8 clusters
L = sugm.generator(graph = "cluster", g = 8, vis = TRUE)

## scale-free graphs
L = sugm.generator(graph="scale-free", vis = TRUE)

```

sugm.plot

*Graph visualization for an object with S3 class "sugm"***Description**

Implements the graph visualization using adjacency matrix. It can automatic organize 2D embedding layout.

Usage

```
sugm.plot(G, epsflag = FALSE, graph.name = "default", cur.num = 1,
          location)
```

Arguments

G	The adjacency matrix corresponding to the graph.
epsflag	If epsflag = TRUE, save the plot as an eps file in the target directory. The default value is FALSE.
graph.name	The name of the output eps files. The default value is "default".
cur.num	The number of plots saved as eps files. Only applicable when epsflag = TRUE. The default value is 1.
location	Target directory. The default value is the current working directory.

Details

The user can change cur.num to plot several figures and select the best one. The implementation is based on the popular package "igraph".

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See Also

[flare](#) and [flare-package](#)

Examples

```
## load package required
library(flare)

## visualize the hub graph
L = sugm.generator(graph = "hub")
sugm.plot(L$theta)

## visualize the band graph
L = sugm.generator(graph = "band",g=5)
sugm.plot(L$theta)

## visualize the cluster graph
L = sugm.generator(graph = "cluster")
sugm.plot(L$theta)

## Not run:
#show working directory
getwd()
#plot 5 graphs and save the plots as eps files in the working directory
sugm.plot(L$theta, epsflag = TRUE, cur.num = 5)

## End(Not run)
```

sugm.roc

Draw ROC Curve for an object with S3 class "sugm"

Description

Draws ROC curve for a graph path according to the true graph structure.

Usage

```
sugm.roc(path, theta, verbose = TRUE)
```

Arguments

path	A graph path.
theta	The true graph structure.
verbose	If verbose = FALSE, tracing information printing is disabled. The default value is TRUE.

Details

To avoid the horizontal oscillation, false positive rates is automatically sorted in the ascent order and true positive rates also follow the same order.

Value

An object with S3 class "roc" is returned:

F1	The F1 scores along the graph path.
tp	The true positive rates along the graph path
fp	The false positive rates along the graph paths
AUC	Area under the ROC curve

Note

For a lasso regression, the number of nonzero coefficients is at most $n-1$. If $d \gg n$, even when regularization parameter is very small, the estimated graph may still be sparse. In this case, the AUC may not be a good choice to evaluate the performance.

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See Also

[sugm](#) and [flare-package](#)

Examples

```
## load package required
library(flare)

#generate data
L = sugm.generator(d = 30, graph = "random", prob = 0.1)
out1 = sugm(L$data, lambda=10^(seq(log10(.4), log10(0.03), length.out=20)))

#draw ROC curve
Z1 = sugm.roc(out1$path,L$theta)

#Maximum F1 score
max(Z1$F1)
```

sugm.select

*Model selection for high-dimensional undirected graphical models***Description**

Implements the regularization parameter selection for high dimensional undirected graphical models. The optional approaches are stability approach to regularization selection (stars) and cross validation selection (cv).

Usage

```
sugm.select(est, criterion = "stars", stars.subsample.ratio = NULL,
            stars.thresh = 0.1, rep.num = 20, fold = 5,
            loss="likelihood", verbose = TRUE)
```

Arguments

est	An object with S3 class "sugm"
criterion	Model selection criterion. "stars" and "cv" are available for both graph estimation methods. The default value is "stars".
stars.subsample.ratio	The subsampling ratio. The default value is $10 \times \sqrt{n}/n$ when $n > 144$ and 0.8 when $n \leq 144$, where n is the sample size. Only applicable when criterion = "stars".
stars.thresh	The variability threshold in stars. The default value is 0.1. Only applicable when criterion = "stars".
rep.num	The number of subsamplings. The default value is 20.
fold	The number of folds used in cross validation. The default value is 5. Only applicable when criterion = "cv".
loss	Loss to be used in cross validation. Two losses are available: "likelihood" and "tracel2". Default "likelihood". Only applicable when criterion = "cv".
verbose	If verbose = FALSE, tracing information printing is disabled. The default value is TRUE.

Details

Stability approach to regularization selection (stars) is a natural way to select optimal regularization parameter for all three estimation methods. It selects the optimal graph by variability of subsamplings and tends to over-select edges in Gaussian graphical models. Besides selecting the regularization parameters, stars can also provide an additional estimated graph by merging the corresponding subsampled graphs using the frequency counts. The K-fold cross validation is also provided for selecting the parameter λ , and two loss functions are adopted as follow

$$likelihood : Tr(\Sigma\Omega) - \log |\Omega|$$

$$tracel2 : Tr(diag(\Sigma\Omega - I)^2).$$

Value

An object with S3 class "select" is returned:

refit	The optimal graph selected from the graph path
opt.icov	The optimal precision matrix selected.
merge	The graph path estimated by merging the subsampling paths. Only applicable when the input criterion = "stars".
variability	The variability along the subsampling paths. Only applicable when the input criterion = "stars".
opt.index	The index of the selected regularization parameter.
opt.lambda	The selected regularization/thresholding parameter.
opt.sparsity	The sparsity level of "refit".

and anything else included in the input est

Note

The model selection is NOT available when the data input is the sample covariance matrix.

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References

1. T. Cai, W. Liu and X. Luo. A constrained ℓ_1 minimization approach to sparse precision matrix estimation. *Journal of the American Statistical Association*, 2011.
2. B. He and X. Yuan. On non-ergodic convergence rate of Douglas-Rachford alternating direction method of multipliers. *Technical Report*, 2012.

See Also

[sugm](#) and [flare-package](#).

Examples

```
## load package required
library(flare)

#generate data
L = sugm.generator(d = 10, graph="hub")
out1 = sugm(L$data)

#model selection using stars
#out1.select1 = sugm.select(out1, criterion = "stars", stars.thresh = 0.1)
#plot(out1.select1)

#model selection using cross validation
```

```
out1.select2 = sugm.select(out1, criterion = "cv")  
plot(out1.select2)
```

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