

Package ‘pdglasso’

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Maintainer The package maintainer <yourself@somewhere.net>
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R topics documented:

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admm.pdglasso	<i>Estimate a concentration matrix under the pdColG model using adaptive ADMM graphical lasso algorithm.</i>
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Description

Description here.

Usage

```
admm.pdglasso(
  S,
  lambda1 = 1,
  lambda2 = 1e-04,
  type = c("vertex", "inside.block.edge", "across.block.edge"),
  force.symm = NULL,
  X.init = NULL,
  rho1 = 1,
  rho2 = 1,
  varying.rho1 = TRUE,
  varying.rho2 = TRUE,
  max_iter = 1000,
  eps.abs = 1e-12,
  eps.rel = 1e-12,
  verbose = FALSE
)
```

Arguments

<code>S</code>	A $p \times p$ covariance (or correlation) matrix.
<code>lambda1</code>	A non-negative scalar (or vector) penalty that encourages sparsity in the concentration matrix. If a vector is provided, it should match the appropriate length, i.e.
<code>lambda2</code>	A non-negative scalar (or vector) penalty that encourages equality constraints in the concentration matrix. If a vector is provided, it should match the appropriate length, i.e.
<code>type</code>	A string or vector of strings for the type of equality constraints to be imposed; zero, one or more available options can be selected among: * "vertex", symmetries are imposed on the diagonal entries of the concentration matrix. * "inside.block.edge", symmetries are imposed between elements of the LL and RR block the concentration matrix. * "across.block.edge", symmetries are imposed between elements of the LR and RL block the concentration matrix. Shortened forms are accepted too, i.e. "V" or "vert" for "vertex".
<code>force.symm</code>	A string or vector of strings to impose forced symmetry on the corresponding block of the concentration matrix. Same options as "type".
<code>X.init</code>	(optional) A $p \times p$ initial guess for the concentration matrix and/or starting solution for the ADMM algorithm.
<code>rho1</code>	A scalar; tuning parameter of the ADMM algorithm to be used for the outer loop. It must be strictly positive.
<code>rho2</code>	A scalar; tuning parameter of the ADMM algorithm to be used for the inner loop. It must be strictly positive.
<code>varying.rho1</code>	A boolean value; if TRUE the parameter rho1 is updated iteratively to speed-up convergence.
<code>varying.rho2</code>	A boolean value; if TRUE the parameter rho2 is updated iteratively to speed-up convergence.
<code>max_iter</code>	An integer; maximum number of iterations to be run in case the algorithm does not converge.

eps.abs	A scalar; the absolute precision required for the computation of primal and dual residuals of the ADMM algorithm.
eps.rel	A scalar; the relative precision required for the computation of primal and dual residuals of the ADMM algorithm.
verbose	A boolean value; if TRUE the progress (and internal convergence of inner loop) is shown in the console while the algorithm is running.

Value

A list, whose element are: * X, the estimated concentration matrix under the pdglasso model; the model is identified by the values of lambda1 and lambda 2, together with the type of penalization imposed. * acronims, a vector of strings for the type of penalties and forced symmetries imposed when calling the function. * internal.par, a list of internal parameters passed to the function at the call, as well as convergence information.

Examples

```
!!! Create fake dataset
S <- cov(toy.data)
admm.pdglasso(S)
```

G.merge	<i>Conversion from the multiple matrix representation of the model to the single matrix representation.</i>
---------	---

Description

This is the inverse of the function [G.split](#), i.e. X is equal to G.split(G.merge(X)).

Usage

```
G.merge(X)
```

Arguments

X	list with three upper triangular matrices with entries 0 and 1: G, G.sym and G.across, any of G.sym and G.across may be NULL
---	--

Value

a pXp symmetric matrix with entries 0, 1, and 2

Examples

```
# random generation of a list(G=G, G.sym=G.sym, G.across=G.across)

q <- 5 # this can be any integer
p <- q*2

g.p <- matrix(sample(c(0,1), size=p^2, replace=TRUE), nrow=p, ncol=p)
g.p[lower.tri(g.p, diag=TRUE)] <- 0
```

```

g.q1 <- matrix(sample(c(0,1), size=q^2, replace=TRUE), nrow=q, ncol=q)
g.q1[lower.tri(g.q1, diag = FALSE)] <- 0

g.q2 <- matrix(sample(c(0,1), size=q^2, replace=TRUE), nrow=q, ncol=q)
g.q2[lower.tri(g.q2, diag = TRUE)] <- 0
g.q2sym <- g.q2+t(g.q2)

g.p[1:q, 1:q] <- g.p[1:q, 1:q] *(1-g.q1)
g.p[(q+1):p, (q+1):p] <- g.p[(q+1):p, (q+1):p] *(1-g.q1)
g.p[1:q, (q+1):p] <- g.p[1:q, (q+1):p] * (1-g.q2sym)

# list obtained

X <- list(G=g.p, G.sym=g.q1, G.across=g.q2)

g <- G.merge(X)
gs <- G.split(g)

identical(X, gs)

X <- list(G=g.p, G.sym=NULL, G.across=NULL)

g <- G.merge(X)
gs <- G.split(g)

identical(X, gs)

```

G.split

Conversion from the single matrix representation of the model to the multiple matrix representation.

Description

This is the inverse of the function `G.merge()`, i.e. `g` is equal to `G.merge(G.split(g))`.

Usage

```
G.split(g)
```

Arguments

`g` is a $p \times p$ symmetric matrix with entries 0, 1, and 2

Value

a list with three upper triangular matrices: `G`, `G.sym` and `G.across` with entries 0 and 1, any of `G.sym` and `G.across` may be `NULL`

Examples

```
# see example in function G.merge().
```

get.pdColG	<i>Build a graph from the output of a call to admm.pdglasso.</i>
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Description

Description here.

Usage

```
get.pdColG(admm.out, th1 = NULL, th2 = NULL, verbose = FALSE)
```

Arguments

admm.out	An object of list type, that is the output of a call to the admm-pdglasso function.
th1	(optional) A scalar, the threshold to identify edges in the graph; it must be non-negative.
th2	(optional) A scalar, the threshold to identify coloured edges in the graph; it must be non-negative.
verbose	(optional) if TRUE provides summary statistics of the graph.

Value

a list, containing:

- g, the graph in matrix form.
- dof, the degrees of freedom corresponding to the graph build under the pdglasso model provided.

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