

Package ‘causalDisco’

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Title Tools for Causal Discovery on Observational Data

Version 0.9.5

Description Various tools for inferring causal models from observational data. The package includes an implementation of the temporal Peter-Clark (TPC) algorithm. Petersen, Osler and Ekstrøm (2021) <[doi:10.1093/aje/kwab087](https://doi.org/10.1093/aje/kwab087)>. It also includes general tools for evaluating differences in adjacency matrices, which can be used for evaluating performance of causal discovery procedures.

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'gausCorScore.R' 'maketkz.R' 'metrics.R' 'misc.R' 'nDAGs.R'
'pc.R' 'plot.R' 'plotTempoMech.R' 'probmata2amat.R' 'regTest.R'
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Author Anne Helby Petersen [aut],
Tobias Ellegaard Larsen [ctb],
Claus Thorn Ekstrøm [ctb],
Bjarke Hautop Kristensen [ctb, cre]

Maintainer Bjarke Hautop Kristensen <bjarke.kristensen@sund.ku.dk>

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<i>adj_confusion</i>	<i>Compute confusion matrix for comparing two adjacency matrices</i>
----------------------	----------------------------------------------------------------------

Description

Two adjacency matrices are compared either in terms of adjacencies (`type = "adj"`) or orientations (`type = "dir"`).

Usage

```
adj_confusion(est_amat, true_amat)
```

Arguments

- `est_amat` The estimated adjacency matrix, or tpdag/cpdag object as obtained from [tpc](#) or [pc](#)
- `true_amat` The true adjacency matrix, or tpdag/cpdag object as obtained from [tpc](#) or [pc](#)

Details

Adjacency comparison: The confusion matrix is a cross-tabulation of adjacencies. Hence, a true positive means that the two inputs agree on the presence of an adjacency. A true negative means that the two inputs agree on no adjacency. A false positive means that `est_amat` places an adjacency where there should be none. A false negative means that `est_amat` does not place an adjacency where there should have been one.

Orientation comparison: The orientation confusion matrix is conditional on agreement on adjacency. This means that only adjacencies that are shared in both input matrices are considered, and agreement wrt. orientation is then computed only among these edges that occur in both matrices. A true positive is a correctly placed arrowhead (1), a false positive marks placement of arrowhead (1) where there should have been a tail (0), a false negative marks placement of tail (0) where there should have been an arrowhead (1), and a true negative marks correct placement of a tail (0).

Value

A list with entries `$tp` (number of true positives), `$tn` (number of true negatives), `$fp` (number of false positives), and `$fn` (number of false negatives).

Examples

```
#####
# Compare two adjacency matrices #####
#####
x1 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               1, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
x2 <- matrix(c(0, 0, 1, 0,
               1, 0, 0, 0,
               0, 0, 0, 0,
               1, 0, 1, 0), 4, 4, byrow = TRUE)

# confusion matrix for adjacencies
confusion(x2, x1)

# confusion matrix for conditional orientations
confusion(x2, x1, type = "dir")

#####
# Compare estimated cpdag with true adjacency matrix #####
#####
# simulate DAG adjacency matrix and Gaussian data
set.seed(123)
x3 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               0, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
ex_data <- simGausFromDAG(x3, n = 50)
pcres <- pc(ex_data, sparsity = 0.1, test = corTest)

# compare adjacencies with true amat (x1)
confusion(pcres, x3)

# compare conditional orientations with true amat
confusion(pcres, x1, type = "dir")
```

amat

Extract adjacency matrix from tpdag, cpdag, tpag or pag object

Description

If the input is a tpdag or cpdag, the resulting adjacency matrix A is "from-to" matrix encoded as follows: - $A(i,j) = 1$ and $A(j,i) = 0$ means there is an edge $j \rightarrow i$. - $A(j,i) = 1$ and $A(i,j) = 0$ means there is an edge $i \rightarrow j$. - $A(i,j) = 1$ and $A(j,i) = 1$ means there is an undirected edge between i and j , $i - j$. - $A(i,j) = 0$ and $A(j,i) = 0$ means there is no edge between i and j .

Usage

```
amat(x)
```

Arguments

x tpdag, cpdag, tpag, or pag object as obtained from [tpc](#), [pc](#), [tfci](#), or [fci](#), respectively.

Details

If the inout is a tpag or pag, there are four possible entry values: 0 (no edge), 1 (circle), 2 (arrow-head), 3 (tail). It is still encoded as a "to-from" adjacency matrix, which means that e.g. $A(i,j) = 1$ places a circle in the directed from j to i. For example, if $A(i,j) = 1$ and $A(j,i) = 2$, we have that $i \rightarrow j$. Similarly, $A(i,j) = 2$ and $A(j,i) = 3$ mean that $i \leftarrow j$.

as.graphNEL

Convert adjacency matrix to graphNEL object

Description

Convert adjacency matrix to graphNEL object

Usage

```
as.graphNEL(amat)
```

Arguments

amat An adjacency matrix

Value

A graphNEL object, see [graphNEL-class](#).

average_degree

Compute average degree for adjacency matrix

Description

Computes the average degree, i.e. the number of edges divided by the number of nodes.

Usage

```
average_degree(amat)
```

Arguments

`amat` An adjacency matrix

Value

A numeric.

<code>compare</code>	<i>Compare two tpdag or tskeleton objects</i>
----------------------	-----------------------------------------------

Description

Compare edges in two tpdag objects or two tskeleton objects. Note that they should be based on the same variables. Only edge absence/presence is compared, not edge orientation.

Usage

```
compare(x, y = NULL)
```

Arguments

`x` First object
`y` Second object (optional)

Value

A list with entries: `$nedges1` (the number of edges in the first object), `$nedges2` (the number of edges in the second object), `$psi1` (the test significance level of the first object), `$psi2` (the test significance level of the second object), `$nadded` (the number of additional edges in object 2, relative to object 1), and `nremoved` (the number of absent edges in object 2, relative to object 1).

<code>confusion</code>	<i>Compute confusion matrix for comparing two adjacency matrices</i>
------------------------	----------------------------------------------------------------------

Description

Two adjacency matrices are compared either in terms of adjacencies (`type = "adj"`) or orientations (`type = "dir"`).

Usage

```
confusion(est_amat, true_amat, type = "adj")
```

Arguments

<code>est_amat</code>	The estimated adjacency matrix, or <code>tpdag/cpdag</code> object as obtained from tpc or pc
<code>true_amat</code>	The true adjacency matrix, or <code>tpdag/cpdag</code> object as obtained from tpc or pc
<code>type</code>	String indicating whether the confusion matrix should be computed for adjacencies (<code>"adj"</code> , the default) or for (conditional) orientations (<code>dir</code>).

Details

Adjacency comparison: The confusion matrix is a cross-tabulation of adjacencies. Hence, a true positive means that the two inputs agree on the presence of an adjacency. A true negative means that the two inputs agree on no adjacency. A false positive means that `est_amat` places an adjacency where there should be none. A false negative means that `est_amat` does not place an adjacency where there should have been one.

Orientation comparison: The orientation confusion matrix is conditional on agreement on adjacency. This means that only adjacencies that are shared in both input matrices are considered, and agreement wrt. orientation is then computed only among these edges that occur in both matrices. A true positive is a correctly placed arrowhead (1), a false positive marks placement of arrowhead (1) where there should have been a tail (0), a false negative marks placement of tail (0) where there should have been an arrowhead (1), and a true negative marks correct placement of a tail (0).

Value

A list with entries `$tp` (number of true positives), `$tn` (number of true negatives), `$fp` (number of false positives), and `$fn` (number of false negatives).

Examples

```
#####
# Compare two adjacency matrices #####
#####
x1 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               1, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
x2 <- matrix(c(0, 0, 1, 0,
               1, 0, 0, 0,
               0, 0, 0, 0,
               1, 0, 1, 0), 4, 4, byrow = TRUE)

# confusion matrix for adjacencies
confusion(x2, x1)

# confusion matrix for conditional orientations
confusion(x2, x1, type = "dir")

#####
# Compare estimated cpdag with true adjacency matrix #####
#####
```

```
# simulate DAG adjacency matrix and Gaussian data
set.seed(123)
x3 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               0, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
ex_data <- simGausFromDAG(x3, n = 50)
pcres <- pc(ex_data, sparsity = 0.1, test = corTest)

# compare adjacencies with true amat (x1)
confusion(pcres, x3)

# compare conditional orientations with true amat
confusion(pcres, x1, type = "dir")
```

corTest	<i>Test for vanishing partial correlations</i>
---------	------------------------------------------------

Description

This function simply calls the [gaussCItest](#) function from the `pcalg` package.

Usage

```
corTest(x, y, S, suffStat)
```

Arguments

x	Index of x variable
y	Index of y variable
S	Index of S variable(s), possibly NULL
suffStat	Sufficient statistic; list with data, binary variables and order.

Value

A numeric, which is the p-value of the test.

dir_confusion	<i>Compute confusion matrix for comparing two adjacency matrices</i>
---------------	----------------------------------------------------------------------

Description

Two adjacency matrices are compared either in terms of adjacencies (type = "adj") or orientations (type = "dir").

Usage

```
dir_confusion(est_amat, true_amat)
```

Arguments

est_amat	The estimated adjacency matrix, or tpdag/cpdag object as obtained from tpc or pc
true_amat	The true adjacency matrix, or tpdag/cpdag object as obtained from tpc or pc

Details

Adjacency comparison: The confusion matrix is a cross-tabulation of adjacencies. Hence, a true positive means that the two inputs agree on the presence of an adjacency. A true negative means that the two inputs agree on no adjacency. A false positive means that est_amat places an adjacency where there should be none. A false negative means that est_amat does not place an adjacency where there should have been one.

Orientation comparison: The orientation confusion matrix is conditional on agreement on adjacency. This means that only adjacencies that are shared in both input matrices are considered, and agreement wrt. orientation is then computed only among these edges that occur in both matrices. A true positive is a correctly placed arrowhead (1), a false positive marks placement of arrowhead (1) where there should have been a tail (0), a false negative marks placement of tail (0) where there should have been an arrowhead (1), and a true negative marks correct placement of a tail (0).

Value

A list with entries \$tp (number of true positives), \$tn (number of true negatives), \$fp (number of false positives), and \$tp (number of false negatives).

Examples

```
#####
# Compare two adjacency matrices #####
#####
x1 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               1, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
x2 <- matrix(c(0, 0, 1, 0,
               1, 0, 0, 0,
```

```

      0, 0, 0, 0,
      1, 0, 1, 0), 4, 4, byrow = TRUE)

# confusion matrix for adjacencies
confusion(x2, x1)

# confusion matrix for conditional orientations
confusion(x2, x1, type = "dir")

#####
# Compare estimated cpdag with true adjacency matrix #####
#####
# simulate DAG adjacency matrix and Gaussian data
set.seed(123)
x3 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               0, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
ex_data <- simGausFromDAG(x3, n = 50)
pcres <- pc(ex_data, sparsity = 0.1, test = corTest)

# compare adjacencies with true amat (x1)
confusion(pcres, x3)

# compare conditional orientations with true amat
confusion(pcres, x1, type = "dir")

```

dir_confusion_original

Compute confusion matrix for comparing two adjacency matrices

Description

Two adjacency matrices are compared either in terms of adjacencies (type = "adj") or orientations (type = "dir").

Usage

```
dir_confusion_original(est_amat, true_amat)
```

Arguments

est_amat	The estimated adjacency matrix, or tpdag/cpdag object as obtained from tpc or pc
true_amat	The true adjacency matrix, or tpdag/cpdag object as obtained from tpc or pc

Details

This is an old version of the function, included for possible backwards compatibility. Edges are scored as follows: A correctly unoriented edge counts as a true negative (TN). An undirected edge that should have been directed counts as a false negative (FN). A directed edge that should have been undirected counts as a false positive (FP). A directed edge oriented in the correct direction counts as a true positive (TP). A directed edge oriented in the incorrect direction counts as both a false positive (FP) and a false negative (FN).

Value

A list with entries \$tp (number of true positives), \$tn (number of true negatives), \$fp (number of false positives), and \$fn (number of false negatives).

Examples

```
#####
# Compare two adjacency matrices #####
#####
x1 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               1, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
x2 <- matrix(c(0, 0, 1, 0,
               1, 0, 0, 0,
               0, 0, 0, 0,
               1, 0, 1, 0), 4, 4, byrow = TRUE)

# confusion matrix for adjacencies
confusion(x2, x1)

# confusion matrix for conditional orientations
confusion(x2, x1, type = "dir")

#####
# Compare estimated cpdag with true adjacency matrix #####
#####
# simulate DAG adjacency matrix and Gaussian data
set.seed(123)
x3 <- matrix(c(0, 0, 0, 0,
               1, 0, 1, 0,
               0, 0, 0, 0,
               0, 0, 1, 0), 4, 4, byrow = TRUE)
ex_data <- simGausFromDAG(x3, n = 50)
pcres <- pc(ex_data, sparsity = 0.1, test = corTest)

# compare adjacencies with true amat (x1)
confusion(pcres, x3)

# compare conditional orientations with true amat
confusion(pcres, x1, type = "dir")
```

edges	<i>List of edges in adjacency matrix</i>
-------	------------------------------------------

Description

Produces a list of edges from an adjacency matrix.

Usage

```
edges(amat)
```

Arguments

amat	An adjacency matrix.
------	----------------------

Value

A list consisting of two lists: One for oriented edges (`$dir`), and one for unoriented edges (`$undir`).

essgraph2amat	<i>Convert essential graph to adjacency matrix</i>
---------------	----------------------------------------------------

Description

Extracts the adjacency matrix from an [EssGraph-class](#) object. This object is returned by score-based causal discovery algorithms in the `pcalg` package.

Usage

```
essgraph2amat(essgraph, p = length(essgraph$field(".nodes")))
```

Arguments

essgraph	An EssGraph object
p	The number of nodes in the graph

Value

An adjacency matrix (square matrix with 0/1 entries).

evaluate

Evaluate adjacency matrix estimation

Description

Applies several different metrics to evaluate difference between estimated and true adjacency matrices. Intended to be used to evaluate performance of causal discovery algorithms.

Usage

```
evaluate(est, true, metrics, ...)
```

Arguments

<code>est</code>	Estimated adjacency matrix/matrices.
<code>true</code>	True adjacency matrix/matrices.
<code>metrics</code>	List of metrics, see details.
<code>...</code>	Further arguments that depend on input type. Currently only <code>list.out</code> is allowed, and only if the first argument is a matrix (see details under Value).

Details

Two options for input are available: Either `est` and `true` can be two adjacency matrices, or they can be two arrays of adjacency matrices. The arrays should have shape $n * p * p$ where n is the number of of matrices, and p is the number of nodes/variables.

The metrics should be given as a list with slots `$adj`, `$dir` and `$other`. Metrics under `$adj` are applied to the adjacency confusion matrix, while metrics under `$dir` are applied to the conditional orientation confusion matrix (see [confusion](#)). Metrics under `$other` are applied without computing confusion matrices first.

Available metrics to be used with confusion matrices are [precision](#), [recall](#), [specificity](#), [FOR](#), [FDR](#), [NPV](#), [F1](#) and [G1](#). The user can supply custom metrics as well: They need to have the confusion matrix as their first argument and should return a numeric.

Available metrics to be used as "other" is: [shd](#). The user can supply custom metrics as well: They need to have arguments `est_amat` and `true_amat`, where the former is the estimated adjacency matrix and the latter is the true adjacency matrix. The metrics should return a numeric.

Value

A `data.frame` with one column for each computed metric and one row per evaluated matrix pair. Adjacency metrics are prefixed with "adj_", orientation metrics are prefixed with "dir_", other metrics do not get a prefix. If the first argument is a matrix, `list.out = TRUE` can be used to change the return object to a list instead. This list will contain three lists, where adjacency, orientation and other metrics are reported, respectively.

evaluate.array

Evaluate adjacency matrix estimation

Description

Applies several different metrics to evaluate difference between estimated and true adjacency matrices. Intended to be used to evaluate performance of causal discovery algorithms.

Usage

```
## S3 method for class 'array'
evaluate(est, true, metrics, ...)
```

Arguments

est	Estimated adjacency matrix/matrices.
true	True adjacency matrix/matrices.
metrics	List of metrics, see details.
...	Further arguments that depend on input type. Currently only <code>list.out</code> is allowed, and only if the first argument is a matrix (see details under Value).

Details

Two options for input are available: Either `est` and `true` can be two adjacency matrices, or they can be two arrays of adjacency matrices. The arrays should have shape $n * p * p$ where n is the number of of matrices, and p is the number of nodes/variables.

The metrics should be given as a list with slots `$adj`, `$dir` and `$other`. Metrics under `$adj` are applied to the adjacency confusion matrix, while metrics under `$dir` are applied to the conditional orientation confusion matrix (see [confusion](#)). Metrics under `$other` are applied without computing confusion matrices first.

Available metrics to be used with confusion matrices are [precision](#), [recall](#), [specificity](#), [FOR](#), [FDR](#), [NPV](#), [F1](#) and [G1](#). The user can supply custom metrics as well: They need to have the confusion matrix as their first argument and should return a numeric.

Available metrics to be used as "other" is: [shd](#). The user can supply custom metrics as well: They need to have arguments `est_amat` and `true_amat`, where the former is the estimated adjacency matrix and the latter is the true adjacency matrix. The metrics should return a numeric.

Value

A `data.frame` with one column for each computed metric and one row per evaluated matrix pair. Adjacency metrics are prefixed with "adj_", orientation metrics are prefixed with "dir_", other metrics do not get a prefix. If the first argument is a matrix, `list.out = TRUE` can be used to change the return object to a list instead. This list will contain three lists, where adjacency, orientation and other metrics are reported, respectively.

evaluate.matrix	<i>Evaluate adjacency matrix estimation</i>
-----------------	---------------------------------------------

Description

Applies several different metrics to evaluate difference between estimated and true adjacency matrices. Intended to be used to evaluate performance of causal discovery algorithms.

Usage

```
## S3 method for class 'matrix'
evaluate(est, true, metrics, list.out = FALSE, ...)
```

Arguments

<code>est</code>	Estimated adjacency matrix/matrices.
<code>true</code>	True adjacency matrix/matrices.
<code>metrics</code>	List of metrics, see details.
<code>list.out</code>	If FALSE (default), output is returned as a data.frame, otherwise it will be a list.
<code>...</code>	Further arguments that depend on input type. Currently only <code>list.out</code> is allowed, and only if the first argument is a matrix (see details under Value).

Details

Two options for input are available: Either `est` and `true` can be two adjacency matrices, or they can be two arrays of adjacency matrices. The arrays should have shape $n * p * p$ where n is the number of of matrices, and p is the number of nodes/variables.

The metrics should be given as a list with slots `$adj`, `$dir` and `$other`. Metrics under `$adj` are applied to the adjacency confusion matrix, while metrics under `$dir` are applied to the conditional orientation confusion matrix (see [confusion](#)). Metrics under `$other` are applied without computing confusion matrices first.

Available metrics to be used with confusion matrices are [precision](#), [recall](#), [specificity](#), [FOR](#), [FDR](#), [NPV](#), [F1](#) and [G1](#). The user can supply custom metrics as well: They need to have the confusion matrix as their first argument and should return a numeric.

Available metrics to be used as "other" is: [shd](#). The user can supply custom metrics as well: They need to have arguments `est_amat` and `true_amat`, where the former is the estimated adjacency matrix and the latter is the true adjacency matrix. The metrics should return a numeric.

Value

A data.frame with one column for each computed metric and one row per evaluated matrix pair. Adjacency metrics are prefixed with "adj_", orientation metrics are prefixed with "dir_", other metrics do not get a prefix. If the first argument is a matrix, `list.out = TRUE` can be used to change the return object to a list instead. This list will contain three lists, where adjacency, orientation and other metrics are reported, respectively.

evaluate.tamat

Evaluate adjacency matrix estimation

Description

Applies several different metrics to evaluate difference between estimated and true adjacency matrices. Intended to be used to evaluate performance of causal discovery algorithms.

Usage

```
## S3 method for class 'tamat'
evaluate(est, true, metrics, ...)
```

Arguments

est	Estimated adjacency matrix/matrices.
true	True adjacency matrix/matrices.
metrics	List of metrics, see details.
...	Further arguments that depend on input type. Currently only list.out is allowed, and only if the first argument is a matrix (see details under Value).

Details

Two options for input are available: Either est and true can be two adjacency matrices, or they can be two arrays of adjacency matrices. The arrays should have shape $n * p * p$ where n is the number of of matrices, and p is the number of nodes/variables.

The metrics should be given as a list with slots \$adj, \$dir and \$other. Metrics under \$adj are applied to the adjacency confusion matrix, while metrics under \$dir are applied to the conditional orientation confusion matrix (see [confusion](#)). Metrics under \$other are applied without computing confusion matrices first.

Available metrics to be used with confusion matrices are [precision](#), [recall](#), [specificity](#), [FOR](#), [FDR](#), [NPV](#), [F1](#) and [G1](#). The user can supply custom metrics as well: They need to have the confusion matrix as their first argument and should return a numeric.

Available metrics to be used as "other" is: [shd](#). The user can supply custom metrics as well: They need to have arguments est_amat and true_amat, where the former is the estimated adjacency matrix and the latter is the true adjacency matrix. The metrics should return a numeric.

Value

A data.frame with one column for each computed metric and one row per evaluated matrix pair. Adjacency metrics are prefixed with "adj_", orientation metrics are prefixed with "dir_", other metrics do not get a prefix. If the first argument is a matrix, list.out = TRUE can be used to change the return object to a list instead. This list will contain three lists, where adjacency, orientation and other metrics are reported, respectively.

F1	<i>F1 score</i>
----	-----------------

Description

Computes F1 score from a confusion matrix, see [confusion](#). The F1 score is defined as $2 * TP / (2 * TP + FP + FN)$, where TP are true positives, FP are false positives, and FN are false negatives. If $TP + FP + FN = 0$, 1 is returned.

Usage

```
F1(confusion)
```

Arguments

confusion	Confusion matrix as obtained from confusion
-----------	-------------------------------------------------------------

Value

A numeric in [0,1].

fci	<i>Perform causal discovery using the FCI algorithm</i>
-----	---------------------------------------------------------

Description

This is a wrapper function for the [fci](#) function as implemented in the pcalg package. All computations are carried out by the pcalg package. The default output object however matches that of [tfci](#), see this function for details about how the adjacency matrix is encoded.

Usage

```
fci(
  data = NULL,
  sparsity = 10^(-1),
  test = regTest,
  suffStat = NULL,
  method = "stable.fast",
  methodNA = "none",
  methodOri = "conservative",
  output = "pag",
  varnames = NULL,
  ...
)
```

Arguments

<code>data</code>	A <code>data.frame</code> with data. All variables should be assigned to exactly one period by prefixing them with the period name (see example below).
<code>sparsity</code>	The sparsity level to be used for independence testing (i.e. significance level threshold to use for each test).
<code>test</code>	A procedure for testing conditional independence. The default, <code>regTest</code> uses a regression-based information loss test. Another available option is <code>corTest</code> which tests for vanishing partial correlations. User supplied functions may also be used, see details below about the required syntax.
<code>suffStat</code>	Sufficient statistic. If this argument is supplied, the sufficient statistic is not computed from the inputted data. The format and contents of the sufficient statistic depends on which test is being used.
<code>method</code>	Which method to use for skeleton construction, must be <code>"stable"</code> , <code>"original"</code> , or <code>"stable.fast"</code> (the default). See skeleton for details.
<code>methodNA</code>	Method for handling missing information (NA values). Must be one of <code>"none"</code> (default, an error is thrown if NAs are present), <code>"cc"</code> (complete case analysis, deletes all observations that have any NA values), or <code>"twd"</code> (test wise deletion, omits observations with missing information test-by-test) (further details below).
<code>methodOri</code>	Method for handling conflicting separating sets when orienting edges, must be one of <code>"standard"</code> , <code>"conservative"</code> (the default) or <code>"maj.rule"</code> . See pc for further details.
<code>output</code>	One of <code>"pag"</code> or <code>"fciAlgo"</code> . If <code>"pag"</code> a partial ancestral graph (PAG) object is outputted. If <code>"fciAlgo"</code> the PAG is outputted as the object class fciAlgo-class from the <code>pcalg</code> package. This is intended for compatability with tools from that package.
<code>varnames</code>	A character vector of variable names. It only needs to be supplied if the <code>data</code> argument is not used, and data are hence passed exclusively through the <code>suffStat</code> argument.
<code>...</code>	Further optional arguments which are passed to skeleton for the skeleton constructing phase.

Examples

```
# simulate linear Gaussian data w unobserved variable L1
n <- 100
L1 <- rnorm(n)
X1 <- rnorm(n)
X2 <- L1 + X1 + rnorm(n)
X3 <- X1 + rnorm(n)
X4 <- X3 + L1 + rnorm(n)
d <- data.frame(p1_X1 = X1,
                p1_X2 = X2,
                p2_X3 = X3,
                p2_X4 = X4)
```

```
# use FCI algorithm to recover PAG
fci(d, test = corTest)
```

FDR	<i>False Discovery Rate</i>
-----	-----------------------------

Description

Computes false discovery rate from a confusion matrix, see [confusion](#). False discovery rate is defined as $FP/(FP + TP)$, where FP are false positives and TP are true positives. If $FP + TP = 0$, 0 is returned.

Usage

```
FDR(confusion)
```

Arguments

confusion Confusion matrix as obtained from [confusion](#)

Value

A numeric in $[0,1]$.

FOR	<i>False Omission Rate</i>
-----	----------------------------

Description

Computes false omission rate from a confusion matrix, see [confusion](#). False omission rate is defined as $FN/(FN + TN)$, where FN are false negatives and TN are true negatives. If $FN + TN = 0$, 0 is returned.

Usage

```
FOR(confusion)
```

Arguments

confusion Confusion matrix as obtained from [confusion](#)

Value

A numeric in $[0,1]$.

G1	<i>G1 score</i>
----	-----------------

Description

Computes G1 score from a confusion matrix, see [confusion](#). G1 score is F1 score with reversed roles of 0/1 classifications, see Petersen et al. 2022. The G1 score is defined as $2 * TN / (2 * TN + FN + FP)$, where TN are true negatives, FP are false positives, and FN are false negatives. If $TN + FN + FP = 0$, 1 is returned.

Usage

```
G1(confusion)
```

Arguments

confusion	Confusion matrix as obtained from confusion
-----------	-------------------------------------------------------------

Value

A numeric in [0,1].

References

Petersen, Anne Helby, et al. "Causal discovery for observational sciences using supervised machine learning." arXiv preprint arXiv:2202.12813 (2022).

gausCorScore	<i>Gaussian L0 score computed on correlation matrix</i>
--------------	---------------------------------------------------------

Description

The score is intended to be used with score-based causal discovery algorithms from the pcalg package. It is identical to the [pcalg::GaussL0OpenObsScore](#), except that it takes in a correlation matrix instead of the full data set.

Usage

```
gausCorScore(cormat, n, p = NULL, lambda = NULL, ...)
```

Arguments

cormat	A correlation matrix. Needs to be symmetric.
n	The number of observations in the dataset that the correlation matrix was computed from.
p	The number of variables. This is inferred from the cormat if not supplied.
lambda	Penalty to use for the score. If NULL (default), the BIC score penalty is used. See pcalg::GaussLOpenObsScore for further details.
...	Other arguments passed along to pcalg::GaussLOpenObsScore .

Value

A Score object (S4), see [pcalg::Score](#).

Examples

```
# Simulate data and compute correlation matrix
x1 <- rnorm(100)
x2 <- rnorm(100)
x3 <- x1 + x2 + rnorm(100)
d <- data.frame(x1, x2, x3)
cmat <- cor(d)

# Use gausCorScore with pcalg::ges()
pcalg::ges(gausCorScore(cmat, n = 100))
```

getvar.character	<i>Get variables with a specific prefix (character method)</i>
------------------	----------------------------------------------------------------

Description

Get variables with a specific prefix (character method)

Usage

```
## S3 method for class 'character'
getvar(x, prefix, sep = "_")
```

Arguments

x	Character vector
prefix	Prefix to match
sep	Separator, default "_"

Value

Subset of x that matches the prefix

<code>getvar.data.frame</code>	<i>Get variables with a specific prefix (data.frame method)</i>
--------------------------------	-----------------------------------------------------------------

Description

Get variables with a specific prefix (data.frame method)

Usage

```
## S3 method for class 'data.frame'
getvar(x, prefix, sep = "_")
```

Arguments

<code>x</code>	A data.frame
<code>prefix</code>	Prefix to match
<code>sep</code>	Separator, default "_"

Value

Names of columns in x that match the prefix

<code>graph2amat</code>	<i>Convert graphNEL object to adjacency matrix</i>
-------------------------	----------------------------------------------------

Description

Convert graphNEL object to adjacency matrix

Usage

```
graph2amat(graph, toFrom = TRUE, type = "pdag")
```

Arguments

<code>graph</code>	A graphNEL object.
<code>toFrom</code>	Logical indicating whether the resulting adjacency matrix is "to-from" (default), or "from-to", see details.
<code>type</code>	The type of adjacency matrix, must be one of "pdag" or "ag". "pdag" should be used for directed graphs, namely DAG, CPDAG, MPDAG, TPDAG and PDAG adjacency matrices, i.e. adjacency matrices where $A(i,j) = A(j,i) = 1$ is interpreted as an undirected edge. "ag" may be used for ADMGs, MAGs, PAGs and TPAGs, where further possible arrowhead options are available (see amat)

Details

A "to-from" pdag adjacency matrix is encoded as follows: $A(i,j) = 1$ and $A(j,i) = 0$ means there is an edge $i \rightarrow j$. $A(j,i) = 1$ and $A(i,j) = 0$ means there is an edge $j \rightarrow i$. $A(i,j) = 1$ and $A(j,i) = 1$ means there is an undirected edge between i and j , $i - j$. $A(i,j) = 0$ and $A(j,i) = 0$ means there is no edge between i and j .

A "from-to" adjacency matrix is the transpose of a "to-from" adjacency matrix. A "from-to" pdag adjacency matrix is hence encoded as follows: $A(i,j) = 1$ and $A(j,i) = 0$ means there is an edge $j \rightarrow i$. $A(j,i) = 1$ and $A(i,j) = 0$ means there is an edge $i \rightarrow j$. $A(i,j) = 1$ and $A(j,i) = 1$ means there is an undirected edge between i and j , $i - j$. $A(i,j) = 0$ and $A(j,i) = 0$ means there is no edge between i and j .

See [amat](#) for details about how an ag adjacency matrix is encoded.

is_cpdag

Check for CPDAG

Description

Check for CPDAG

Usage

```
is_cpdag(amat)
```

Arguments

amat An adjacency matrix

Details

Check: Is adjacency matrix proper CPDAG? See [isValidGraph](#) for definition.

Value

A logical.

is_pdag	<i>Check for PDAG</i>
---------	-----------------------

Description

Check for PDAG

Usage

```
is_pdag(amat)
```

Arguments

amat	An adjacency matrix
------	---------------------

Details

Check: Is adjacency matrix proper PDAG? See [isValidGraph](#) for definition.

Value

A logical.

maketikz	<i>Generate Latex tikz code for plotting a temporal DAG, PDAG or PAG.</i>
----------	---------------------------------------------------------------------------

Description

Generate Latex tikz code for plotting a temporal DAG, PDAG or PAG.

Usage

```
maketikz(
  model,
  xjit = 2,
  yjit = 2,
  markperiods = TRUE,
  xpgap = 4,
  annotateEdges = NULL,
  addAxis = TRUE,
  varLabels = NULL,
  periodLabels = NULL,
  annotationLabels = NULL,
  clipboard = TRUE,
  rawout = FALSE,
  colorAnnotate = NULL,
  bandedges = FALSE
)
```


Arguments

<code>model</code>	tpdag, tskeleton, tpag, or tamat object to plot.
<code>xjit</code>	How much should nodes within a period be jittered horizontally.
<code>yjit</code>	Vertical distance between nodes within a period.
<code>markperiods</code>	If TRUE, gray boxes are drawn behind each period.
<code>xpgap</code>	Horizontal gap between different periods.
<code>annotateEdges</code>	If TRUE, add a text annotation to edges. If <code>annotationLabels</code> are supplied, these labels will be used. Otherwise, the value in the inputted adjacency matrix corresponding to the edge will be used. Cannot be used for tpag input objects (or ag amat types).
<code>addAxis</code>	If TRUE, a horizontal axis with period labels are added.
<code>varLabels</code>	Optional labels for nodes (variables). Should be given as a named list, where the name is the variable name, and the entry is the label, e.g. <code>list(vname = "Label for vname")</code> .
<code>periodLabels</code>	Optional labels for periods. Should be given as a named list, where the name is the period name (as stored in the tamat), and the entry is the label, e.g. <code>list(periodname = "Label for period")</code> .
<code>annotationLabels</code>	Optional labels for edge annotations. Only used if <code>annotateEdges = TRUE</code> . Should be given as a named list, where the name is the edge annotation (as stored in the tamat), and the entry is the label, e.g. <code>list(h = "High")</code> .
<code>clipboard</code>	If TRUE, the tikz code is not printed, but instead copied to the clipboard, so it can easily be pasted into a Latex document.
<code>rawout</code>	If TRUE, the tikz code is only returned as a character vector.
<code>colorAnnotate</code>	Named list of colors to use to mark edge annotations instead of labels. This overrides <code>annotateEdges</code> and both are not available at the same time. The list should be given with annotations as names and colors as entries, e.g. <code>list(h = "blue")</code> . Cannot be used for tpag input objects (or ag amat types).
<code>bendedges</code>	If TRUE, all edges are bend 10 degrees to the right, thereby avoiding having edges exactly on top of each other.

Details

Note that it is necessary to read in relevant tikz libraries in the Latex preamble. The relevant lines of code are (depending a bit on parameter settings):

```
\usepackage{tikz}
\usetikzlibrary{arrows.meta,arrows,shapes,decorations,automata,backgrounds,petri}
\usepackage{pgfplots}
```

Value

Silently returns a character vector with lines of tikz code. The function furthermore has a side-effect. If `clipboard = TRUE`, the side-effect is that the tikz code is also copied to the clipboard. If `clipboard = FALSE`, the tikz code is instead printed in the console.

Examples

```
# Make tikz figure code from tpdag, print code to screen
data(tpcExample)
tpdag1 <- tpc(tpcExample, order = c("child", "youth", "oldage"), sparsity = 0.01,
             test = corTest)
maketikz(tpdag1, clipboard = FALSE)

# Make tikz figure code from tamat, copy code to clipboard
thisdag <- simDAG(5)
rownames(thisdag) <- colnames(thisdag) <- c("child_x", "child_y",
                                             "child_z", "adult_x",
                                             "adult_y")
thistamat <- tamat(thisdag, order = c("child", "adult"))
## Not run:
maketikz(thistamat)

## End(Not run)
```

maxnedges

Compute maximal number of edges for graph

Description

Computes the number of edges a graph with p nodes will have if its fully connected.

Usage

```
maxnedges(p)
```

Arguments

p The number of nodes in the graph

Value

A numeric.

nDAGs	<i>Number of different DAGs</i>
-------	---------------------------------

Description

Computes the number of different possible DAGs that can be constructed over a given number of nodes.

Usage

nDAGs(p)

Arguments

p The number of nodes.

Value

A numeric.

nedges	<i>Number of edges in adjacency matrix</i>
--------	--------------------------------------------

Description

Counts the number of edges in an adjacency matrix.

Usage

nedges(amat)

Arguments

amat An adjacency matrix

Value

A numeric (non-negative integer).

NPV	<i>Negative predictive value</i>
-----	----------------------------------

Description

Computes negative predictive value recall from a confusion matrix, see [confusion](#). Negative predictive value is defined as $TN/(TN + FN)$, where TN are true negatives and FN are false negatives. If $TN + FN = 0$, 0 is returned.

Usage

```
NPV(confusion)
```

Arguments

confusion	Confusion matrix as obtained from confusion
-----------	-------------------------------------------------------------

Value

A numeric in [0,1].

pc	<i>Perform causal discovery using the PC algorithm</i>
----	--------------------------------------------------------

Description

This is a wrapper function for the [pc](#) function as implemented in the pcalg package. All computations are carried out by the pcalg package.

Usage

```
pc(
  data = NULL,
  sparsity = 10^(-1),
  test = regTest,
  suffStat = NULL,
  method = "stable.fast",
  methodNA = "none",
  methodOri = "conservative",
  output = "cpdag",
  varnames = NULL,
  conservative = TRUE,
  ...
)
```

Arguments

data	A data.frame with data. All variables should be assigned to exactly one period by prefixing them with the period name (see example below).
sparsity	The sparsity level to be used for independence testing (i.e. significance level threshold to use for each test).
test	A procedure for testing conditional independence. The default, <code>regTest</code> uses a regression-based information loss test. Another available option is <code>corTest</code> which tests for vanishing partial correlations. User supplied functions may also be used, see details below about the required syntax.
suffStat	Sufficient statistic. If this argument is supplied, the sufficient statistic is not computed from the inputted data. The format and contents of the sufficient statistic depends on which test is being used.
method	Which method to use for skeleton construction, must be "stable", "original", or "stable.fast" (the default). See skeleton for details.
methodNA	Method for handling missing information (NA values). Must be one of "none" (default, an error is thrown if NAs are present), "cc" (complete case analysis, deletes all observations that have any NA values), or "twd" (test wise deletion, omits observations with missing information test-by-test) (further details below).
methodOri	Method for handling conflicting separating sets when orienting edges, must be one of "standard", "conservative" (the default) or "maj.rule". See pc for further details.
output	One of "cpdag", "skeleton" or "pcAlgo". If "skeleton", a skeleton is constructed and outputted, but the edges are not directed. If "cpdag" (the default), the edges are directed, resulting in a completed partially directed acyclic graph (CPDAG). If "pcAlgo" the CPDAG is outputted as the object class pcAlgo-class from the <code>pcalg</code> package. This is intended for compatability with tools from that package.
varnames	A character vector of variable names. It only needs to be supplied if the data argument is not used, and data are hence passed exclusively through the <code>suffStat</code> argument.
conservative	Logical, if TRUE the conservative version of PC is used (see pc for details).
...	Further optional arguments which are passed to skeleton if <code>output = "skeleton"</code> or to pc otherwise.

Details

Note that all independence test procedures implemented in the `pcalg` package may be used, see [pc](#).

The methods for handling missing information require that the data, rather than the `suffStat` argument is used for inputting data; the latter assumes no missing information and hence always sets `methodNA = "none"`. If the test is `corTest`, test-wise deletion is performed when computing the sufficient statistic (correlation matrix) (so for each pair of variables, only complete cases are used). If the test is `regTest`, test-wise deletion is performed for each conditional independence test instead.

Value

A tpdag or tskeleton object. Both return types are S3 objects, i.e., lists with entries: \$amat (the estimated adjacency matrix), \$order (character vector with the order, as inputted to this function), \$psi (the significance level used for testing), and \$ntests (the number of tests conducted).

Examples

```
# PC on included example data, use sparsity psi = 0.01, default test (regression-based
#information loss):
data(tpcExample)
pc(tpcExample, sparsity = 0.01)

# PC on included example data, use sparsity psi = 0.01, use test for vanishing partial
# correlations:
data(tpcExample)
pc(tpcExample, sparsity = 0.01, test = corTest)
```

plot.pag

Plot partial ancestral graph (PAG)

Description

Plot partial ancestral graph (PAG)

Usage

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

Arguments

x	pag object to be plotted (as outputted from fci).
...	Currently not in use.

Value

No return value, the function is called for its side-effects (plotting).

Author(s)

This code is a modification of the fciAlgo plotting method implemented in the pcalg package.

Examples

```
# simulate linear Gaussian data w unobserved variable L1
n <- 100
L1 <- rnorm(n)
X1 <- rnorm(n)
X2 <- L1 + X1 + rnorm(n)
X3 <- X1 + rnorm(n)
X4 <- X3 + L1 + rnorm(n)
d <- data.frame(p1_X1 = X1,
                p1_X2 = X2,
                p2_X3 = X3,
                p2_X4 = X4)

# use FCI algorithm to recover PAG
res <- fci(d, test = corTest)

# plot
plot(res)
```

plot.tamat

Plot adjacency matrix with order information

Description

Plot adjacency matrix with order information

Usage

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

Arguments

x tamam (temporal adjacency matrix) object to be plotted (as outputted from [tamam](#)).

... Further plotting arguments passed along to [plotTempoMech](#).

Value

No return value, the function is called for its side-effects (plotting).

plot.tpag

Plot temporal partial ancestral graph (TPAG)

Description

Plot temporal partial ancestral graph (TPAG)

Usage

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

Arguments

x tpag object to be plotted (as outputted from [tfci](#)).
... Currently not in use.

Value

No return value, the function is called for its side-effects (plotting).

Author(s)

This code is a modification of the fciAlgo plotting method implemented in the pcalg package.

Examples

```
# simulate linear Gaussian data w unobserved variable L1
n <- 100
L1 <- rnorm(n)
X1 <- rnorm(n)
X2 <- L1 + X1 + rnorm(n)
X3 <- X1 + rnorm(n)
X4 <- X3 + L1 + rnorm(n)
d <- data.frame(p1_X1 = X1,
                p1_X2 = X2,
                p2_X3 = X3,
                p2_X4 = X4)

# use FCI algorithm to recover PAG
res <- tfci(d, order = c("p1", "p2"), test = corTest)

# plot
plot(res)
```

plot.tpdag	<i>Plot temporal partially directed acyclic graph (TPDAG)</i>
------------	---------------------------------------------------------------

Description

Plot temporal partially directed acyclic graph (TPDAG)

Usage

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

Arguments

x	tpdag (temporal partially directed acyclic graph) object to be plotted (as outputted from tpc).
...	Further plotting arguments passed along to plotTempoMech .

Value

No return value, the function is called for its side-effects (plotting).

plot.tskeleton	<i>Plot temporal skeleton</i>
----------------	-------------------------------

Description

Plot temporal skeleton

Usage

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

Arguments

x	tskeleton (temporal skeleton) object to be plotted (as outputted from tpc).
...	Further plotting arguments passed along to plotTempoMech .

Value

No return value, the function is called for its side-effects (plotting).

plotTempoMech	<i>Plot temporal data generating mechanism</i>
---------------	------------------------------------------------

Description

Plots tpdag, tskeleton and tamat objects.

Usage

```
plotTempoMech(
  x,
  addTimeAxis = TRUE,
  addPsi = TRUE,
  varLabels = NULL,
  periodLabels = NULL,
  colors = NULL,
  ...
)
```

Arguments

x	The tpdag/tskeleton or tamat to plot.
addTimeAxis	Logical indicating whether a time axis should be added to the plot.
addPsi	Logical indicating whether the sparsity level should be added to the plot.
varLabels	A named list of variable labels.
periodLabels	A character vector with labels for periods.
colors	A character vector with colors to use for marking periods. Should have at least as many elements as the numbers of periods.
...	Additional arguments passed to plot.igraph .

Value

No return value, the function is called for its side-effects (plotting).

precision	<i>Precision</i>
-----------	------------------

Description

Computes precision (aka positive predictive value) from a confusion matrix, see [confusion](#). Precision is defined as $TP/(TP + FP)$, where TP are true positives and FP are false positives. If $TP + FP = 0$, 0 is returned.

Usage

```
precision(confusion)
```

Arguments

confusion Confusion matrix as obtained from [confusion](#)

Value

A numeric in [0,1].

probmata2amat	<i>Convert a matrix of probabilities into an adjacency matrix</i>
---------------	-------------------------------------------------------------------

Description

Convert a matrix of probabilities into an adjacency matrix

Usage

```
probmata2amat(
  probmat,
  threshold,
  method = "cutoff",
  keep_vnames = TRUE,
  graph_criterion = "pdag",
  deletesym = FALSE
)
```

Arguments

probmat	Square matrix of probabilities.
threshold	Value between 0 and 1. Any probabilities lower than this value will be set to 0 (no arrowhead).
method	Either "cutoff" or "bpco", see details.
keep_vnames	If TRUE, variable names (provided as rownames in the input probmat) will be preserved in the output.
graph_criterion	Which criterion to check if the output graph fulfills for the bpco method. Should be one of "dag", "pdag" or "cpdag" or NULL. Choosing NULL (the default) puts no further restrictions on the output. See isValidGraph for definitions.
deletesym	If TRUE, edges are deleted symmetrically in the bpco method. This means that instead of removing arrowheads (setting singular elements to 0), the procedure removes full edges (setting both potential arrowheads for the given edge to zero). This only makes a difference if the graph may include undirected edges, which should be encoded as bidirected edges.

Details

Two methods for converting the probability matrix into an adjacency matrix are implemented. First, the cutoff-method (`method = "cutoff"`) simply uses a threshold value and sets all values below that to zero in the outputted adjacency matrix. No checks are performed to ensure that the resulting matrix is a proper dag/pdag/cpdag adjacency matrix. Second, the backwards PC orientation method (`method = "bpc"`) first uses a cutoff, and then sets further elements to zero until the resulting matrix can be converted into a proper adjacency matrix (using the graph criterion specified in the `graph_criterion` argument) by applying the PC algorithm orientation rules. See Petersen et al. 2022 for further details.

Value

A square matrix of probabilities (all entries in $[0,1]$).

References

Petersen, Anne Helby, et al. "Causal discovery for observational sciences using supervised machine learning." arXiv preprint arXiv:2202.12813 (2022).

Examples

```
#Make random probability matrix that can be
#converted into adjacency matrix
pmat <- matrix(runif(25, 0, 1), 5, 5)
diag(pmat) <- 0

#Convert to adjacency matrix using cutoff-method (threshold = 0.5)
probm2amat(pmat, threshold = 0.5)

#Convert to adjacency matrix using BPCO-method (threshold = 0.5)
probm2amat(pmat, threshold = 0.5, method = "bpc")
```

recall

Recall

Description

Computes recall from a confusion matrix, see [confusion](#). Recall is defined as $TP/(TP + FN)$, where TP are true positives and FN are false negatives. If $TP + FN = 0$, 0 is returned.

Usage

```
recall(confusion)
```

Arguments

confusion Confusion matrix as obtained from [confusion](#)

Value

A numeric in [0,1].

regTest

Regression-based information loss test

Description

We test whether x and y are associated, given S using a generalized linear model.

Usage

```
regTest(x, y, S, suffStat)
```

Arguments

x	Index of x variable
y	Index of y variable
S	Index of S variable(s), possibly NULL
suffStat	Sufficient statistic; list with data, binary variables and order.

Details

All included variables should be either numeric or binary. If y is binary, a logistic regression model is fitted. If y is numeric, a linear regression model is fitted. x and S are included as explanatory variables. Any numeric variables among x and S are modeled with spline expansions (natural splines, 3 df). This model is tested against a numeric where x (including a possible spline expansion) has been left out using a likelihood ratio test. The model is fitted in both directions (interchanging the roles of x and y). The final p-value is the maximum of the two obtained p-values.

Value

A numeric, which is the p-value of the test.

shd	<i>Structural hamming distance between adjacency matrices</i>
-----	---------------------------------------------------------------

Description

Computes the structural hamming distance between two adjacency matrices. This implementation is a modification of the [shd](#) function from the pcalg package, but here we avoid working on the heavy graphNEL objects for representing graphs that are used in the pcalg package.

Usage

```
shd(est_amat, true_amat)
```

Arguments

est_amat	Estimated adjacency matrix
true_amat	True adjacency matrix

Details

Note that the function is symmetric in the two inputted adjacency matrices.

Value

A numeric (a non-negative integer).

simDAG	<i>Simulate a random DAG</i>
--------	------------------------------

Description

Simulates a random directed acyclic graph adjacency (DAG) matrix with the provided edge sparsity. The edge sparsity is the percentage of edges that are absent, relative to a fully connected DAG.

Usage

```
simDAG(p, sparsity = NULL, sparsityLim = c(0, 0.8), permute = TRUE)
```

Arguments

p	The number of nodes.
sparsity	If NULL (the default), a random edge sparsity is sampled from the interval provided in sparsityLim. Otherwise, the sparsity should be provided as a numeric in [0,1].
sparsityLim	A vector of two numerics, both must be in [0,1].
permute	If FALSE, the adjacency matrix will include nodes in their causal ordering. This is avoided by setting permute = TRUE, in which case the node order is permuted randomly.

Value

An adjacency matrix.

Examples

```
# Simulate a DAG adjacency matrix with 5 nodes
simDAG(5)
```

simGausFromDAG	<i>Simulate Gaussian data according to DAG</i>
----------------	------------------------------------------------

Description

Simulates a jointly Gaussian dataset given a DAG adjacency matrix ("from-to" encoding, see [amat](#) for details). The data is simulated using linear structural equations and the parameters (residual standard deviations and regression coefficients) are sampled from chosen intervals.

Usage

```
simGausFromDAG(
  amat,
  n,
  regparLim = c(0.5, 2),
  resSDLim = c(0.1, 1),
  pnegRegpar = 0.4,
  standardize = FALSE
)
```

Arguments

amat	An adjacency matrix.
n	The number of observations that should be simulated.
regparLim	The interval from which regression parameters are sampled.
resSDLim	The interval from which residual standard deviations are sampled.
pnegRegpar	The probability of sampling a negative regression parameter.
standardize	If FALSE (the default), the raw data are returned. If TRUE, the data are first standardized, i.e., each variable will have its mean subtracted and be divided by its standard deviation.

Details

A variable X_i is simulated as

$$X_i := \sum_{Z \in pa(X_i)} \beta_Z * Z + e_i$$

where $pa(X_i)$ are the parents of X_i in the DAG. The residual, e_i , is drawn from a normal distribution.

Value

A data.frame of identically distributed simulated observations.

Examples

```
# Simulate DAG adjacency matrix with 6 nodes
ex_dag <- simDAG(6)

# Simulate Gaussian data (100 iid observations)
ex_data <- simGausFromDAG(ex_dag, n = 100)
```

specificity	<i>Specificity</i>
-------------	--------------------

Description

Computes specificity from a confusion matrix, see [confusion](#). Specificity is defined as $TN/(TN + FP)$, where TN are true negatives and FP are false positives. If $TN + FP = 0$, 0 is returned.

Usage

```
specificity(confusion)
```

Arguments

confusion Confusion matrix as obtained from [confusion](#)

Value

A numeric in [0,1].

tamat	<i>Make a temporal adjacency matrix</i>
-------	-----------------------------------------

Description

Make a temporal adjacency matrix

Usage

```
tamat(amat, order, type = NULL)
```


Arguments

amat	Adjacency matrix. Row names and column names should be identical and be the names of the variables/nodes. Variable names should be prefixed with their period, e.g. "child_x" for variable "x" at period "child"
order	A character vector with the periods in their order.
type	The type of adjacency matrix, must be one of "pdag" or "ag". If NULL (default), the function first checks for a tamat_type attribute in the input object and makes sure the output matches that, and if no the input does not have this attribute, it is set to "tpdag". Otherwise, the user can specify a type manually as follows: "pdag" should be used for directed graphs, namely DAG, CPDAG, MPDAG, TPDAG and PDAG adjacency matrices, i.e. adjacency matrices where $A(i,j) = A(j,i) = 1$ is interpreted as an undirected edge. "ag" may be used for ADMGs, MAGs, PAGs and TPAGs, where further possible arrowhead options are available (see amat)

Value

A tamat object, which is a matrix with a "order" attribute(a character vector listing the temporal order of the variables in the adjacency matrix).

TemporalBDeu-class	<i>Temporal Bayesian Dirichlet equivalent uniform (Score criterion)</i>
--------------------	-------------------------------------------------------------------------

Description

A reference class for categorical observational data Scoring with Tiered Background Knowledge. This class represents a score for causal discovery using tiered background knowledge from observational categorical data; it is used in the causal discovery function [tges](#).

Arguments

data	A numeric matrix with n rows and p columns. Each row corresponds to one observational realization.
order	A vector specifying the order each variable. Can be either a vector of integers or an vector of prefixes. If integers, such that the i th entry will detail the order of the i th variable in the dataset. Must start at 1 and increase with increments of 1. If prefixes, must be in order.
iss	Imaginary Sample Size (ISS), also referred to as Equivalent Sample Size (ESS), determines how much weight is assigned to the prior in terms of the size of an imaginary sample supporting it. Increasing the ISS will increase the density of the estimated graph.

Details

The class implements a score which scores all edges contradicting the ordering (edge going from a later tier to an earlier) to minus ∞ . If the edges does not contradict, the score is equal to that of the standard BDeu.

Extends

Class `pcalg::Score` directly.

All reference classes extend and inherit methods from `envRefClass`.

Constructor

```
new("TemporalBdeu",
    data = matrix(1, 1, 1),
    order = rep(1, ncol(data)),
    iss = 1
    ...)
```

Alternative implementation of TemporalBDeu score

We provide here a faster alternative to the implemented version of TemporalBDeu. However, this version relies on a non-exported function from **bnlearn**. We provide the code for it here:

```
setRefClass("TemporalBDeu",
    contains = "DataScore",

    fields = list(
        .order = "vector",
        .iss = "numeric"),

    methods = list(
        initialize = function(data = matrix(1, 1, 1),
                                nodes = colnames(data),
                                iss = 1,
                                order = rep(1, ncol(data)),
                                ...) {
            .order <- order
            .iss <- iss
            callSuper(data = data,
                        nodes = nodes,
                        iss = iss,
                        ...)},

        local.score = function(vertex, parents,...) {
            ## Check validity of arguments
            validate.vertex(vertex)
            validate.parents(parents)
            order <- .order
            iss <- .iss
            if (order[vertex] >= max(c(order[parents], -Inf))){
                #Checks if the tier of parents are before or same as node

                # Create local dataset
```

```

D <- pp.dat$data
pa_nam <- colnames(D)[parents]
ve_nam <- colnames(D)[vertex]
res_nam <- colnames(D)[-c(parents,vertex)]

# Create local bn object
if (length(parents) > 0){
  mod_string <- paste(c("[",
                        paste(c(pa_nam,res_nam),collapse = "]["),
                        "][",
                        ve_nam,
                        "|",
                        paste(pa_nam,collapse = ":"),
                        "]"
                        ),collapse = "")
} else {
  mod_string <- paste(c("[",
                        paste(c(pa_nam,res_nam),collapse = "]["),
                        "][",
                        ve_nam,
                        "]"
                        ),collapse = "")
}

bn_ob <- as.bn(mod_string)
BdeuScore <- bnlearn::per.node.score(network = bn_ob,
                                     data = D,
                                     score = "bde",
                                     targets = ve_nam,
                                     extra.args = list(iss = iss,
                                                       prior = "uniform"))

  return(BdeuScore)
}
else { skip <- -Inf
return(skip)}#set score to minus infinity if vertex earlier than parents
}),
inheritPackage = TRUE
)

```

Author(s)

Tobias Ellegaard Larsen

See Also[tges](#)

Examples

```

# For reproducibility
set.seed(123)

# Number of samples
n <- 1000

# Define probabilities for A
p_A <- c(0.4, 0.35, 0.25) # Probabilities for A = {1, 2, 3}

# Simulate A from a categorical distribution
A <- sample(1:3, n, replace = TRUE, prob = p_A)

# Define conditional probabilities for B given A
p_B_given_A <- list(
  c(0.7, 0.3), # P(B | A=1)
  c(0.4, 0.6), # P(B | A=2)
  c(0.2, 0.8)  # P(B | A=3)
)

# Sample B based on A
B <- sapply(A, function(a) sample(1:2, 1, prob = p_B_given_A[[a]]))

# Define conditional probabilities for C given A and B
p_C_given_A_B <- list(
  "1_1" = c(0.6, 0.4), # P(C | A=1, B=1)
  "1_2" = c(0.3, 0.7), # P(C | A=1, B=2)
  "2_1" = c(0.5, 0.5), # P(C | A=2, B=1)
  "2_2" = c(0.2, 0.8), # P(C | A=2, B=2)
  "3_1" = c(0.7, 0.3), # P(C | A=3, B=1)
  "3_2" = c(0.4, 0.6)  # P(C | A=3, B=2)
)

# Sample C based on A and B
C <- mapply(function(a, b) sample(1:2, 1, prob = p_C_given_A_B[[paste(a, b, sep = "_")]]), A, B)

# Create dataset
simdata <- data.frame(as.factor(A), as.factor(B), as.factor(C))

# Define order in prefix way
colnames(simdata) <- c("child_A", "child_B", "adult_C")
prefix_order <- c("child", "adult")

# Define TemporalBDeu score
t_score <- new("TemporalBDeu", order = prefix_order
              , data = simdata)

# Run tges
tges_pre <- tges(t_score)

# Plot MPDAG
plot(tges_pre)

```

```
# Define order in integer way
colnames(simdata) <- c("A","B","C")
integer_order <- c(1,1,2)

# Define TemporalBDeu score
t_score <- new("TemporalBDeu", order = integer_order
              , data = simdata)

# Run tges
tges_integer <- tges(t_score)

# Plot MPDAG
plot(tges_integer)
```

TemporalBIC-class	<i>Temporal Bayesian Information Criterion (Score criterion)</i>
-------------------	------------------------------------------------------------------

Description

A Reference Class for Gaussian Observational Data Scoring with Tiered Background Knowledge. This class represents a score for causal discovery using tiered background knowledge from observational Gaussian data; it is used in the causal discovery function [tges](#).

Arguments

data	A numeric matrix with n rows and p columns. Each row corresponds to one observational realization.
order	A vector specifying the order each variable. Can be either a vector of integers or an vector of prefixes. If integers, such that the i th entry will detail the order of the i th variable in the dataset. Must start at 1 and increase with increments of 1. If prefixes, must be in order.
lambda	Penalization constant (see details).
intercept	Logical; indicates whether an intercept is allowed in the linear structural equations (i.e., whether a nonzero mean is allowed).

Details

The class implements a score which scores all edges contradicting the ordering (edge going from a later tier to an earlier) to minus ∞ . If the edges does not contradict, the score is equal to that of [pcalg::GaussL0penObsScore](#): The class implements an ℓ_0 -penalized Gaussian maximum likelihood estimator. The penalization is a constant (specified by the argument `lambda` in the constructor) times the number of parameters of the DAG model. By default, the constant λ is chosen as $\log(n)/2$, which corresponds to the BIC score.

Extends

Class [pcalg::GaussL0penObsScore](#) directly.

All reference classes extend and inherit methods from [envRefClass](#).

Constructor

```
new("TemporalBIC",
    data = matrix(1, 1, 1),
    order = rep(1, ncol(data)),
    lambda = 0.5 * log(nrow(data)),
    intercept = TRUE,
    ...)
```

Author(s)

Tobias Ellegaard Larsen

See Also

[tges](#)

Examples

```
#Simulate Gaussian data
set.seed(123)
n <- 500
child_x <- rnorm(n)
child_y <- 0.5*child_x + rnorm(n)
child_z <- 2*child_x + child_y + rnorm(n)

adult_x <- child_x + rnorm(n)
adult_z <- child_z + rnorm(n)
adult_w <- 2*adult_z + rnorm(n)
adult_y <- 2*child_x + adult_w + rnorm(n)

simdata <- data.frame(child_x, child_y, child_z,
                      adult_x, adult_z, adult_w,
                      adult_y)

# Define order in prefix way
prefix_order <- c("child", "adult")

# Define TBIC score
t_score <- new("TemporalBIC", order = prefix_order
              , data = simdata)

# Run tges
tges_pre <- tges(t_score)

# Plot MPDAG
plot(tges_pre)

# Define order in integer way
integer_order <- c(1,1,1,2,2,2,2)

# Define TBIC score
t_score <- new("TemporalBIC", order = integer_order
```

```

                                , data = simdata)
# Run tges
tges_int <- tges(t_score)

# Plot MPDAG
plot(tges_int)

```

tfci

Perform causal discovery using the temporal FCI algorithm (TFCI)

Description

Use a modification of the FCI algorithm that makes use of background knowledge in the format of a partial ordering. This may for instance come about when variables can be assigned to distinct tiers or periods (i.e., a temporal ordering).

Usage

```

tfci(
  data = NULL,
  order,
  sparsity = 10^(-1),
  test = regTest,
  suffStat = NULL,
  method = "stable.fast",
  methodNA = "none",
  methodOri = "conservative",
  varnames = NULL,
  ...
)

```

Arguments

data	A data.frame with data. All variables should be assigned to exactly one period by prefixing them with the period name (see example below).
order	A character vector with period-prefixes in their temporal order (see example below).
sparsity	The sparsity level to be used for independence testing (i.e. significance level threshold to use for each test).
test	A procedure for testing conditional independence. The default, regTest uses a regression-based information loss test. Another available option is corTest which tests for vanishing partial correlations. User supplied functions may also be used, see details below about the required syntax.
suffStat	Sufficient statistic. If this argument is supplied, the sufficient statistic is not computed from the inputted data. The format and contents of the sufficient statistic depends on which test is being used.

method	Which method to use for skeleton construction, must be "stable", "original", or "stable.fast" (the default). See skeleton for details.
methodNA	Method for handling missing information (NA values). Must be one of "none" (default, an error is thrown if NAs are present), "cc" (complete case analysis, deletes all observations that have any NA values), or "twd" (test wise deletion, omits observations with missing information test-by-test) (further details below).
methodOri	Method for handling conflicting separating sets when orienting edges, must be one of "standard", "conservative" (the default) or "maj.rule". See pc for further details.
varnames	A character vector of variable names. It only needs to be supplied if the data argument is not used, and data are hence passed exclusively through the suffStat argument.
...	Further optional arguments which are passed to skeleton for the skeleton constructing phase.

Details

The temporal/tiered background information enters several places in the TFCI algorithm: 1) In the skeleton construction phase, when looking for separating sets Z between two variables X and Y , Z is not allowed to contain variables that are strictly after both X and Y in the temporal order. 2) This also applies to the subsequent phase where the algorithm searches for possible D-SEP sets. 3) Prior to other orientation steps, any cross-tier edges get an arrowhead placed at their latest node.

After this, the usual FCI orientation rules are applied, see [udag2pag](#) for details.

Value

The default output is a tpag object. This is an S3 object, i.e., a list, with entries: \$amat (the estimated adjacency matrix), \$order (character vector with the order, as inputted to this function), \$psi (the significance level used for testing), and \$ntests (the number of tests conducted).

The adjacency matrix A has four possible entry values: 0 (no edge), 1 (circle), 2 (arrowhead), 3 (tail). It is encoded as a "to-from" adjacency matrix, which means that e.g. $A(i,j) = 1$ places a circle in the directed from j to i . For example, if $A(i,j) = 1$ and $A(j,i) = 2$, we have that $i \rightarrow j$. Similarly, $A(i,j) = 2$ and $A(j,i) = 3$ mean that $i <- j$. Note that this is a transposed version of the adjacency matrix outputted for fciAlgo objects from the pcalg package, which is "to-from". But it is consistent with the "from-to" adjacency matrices used for pcAlgo objects from the same package.

Author(s)

Anne Helby Petersen, Qixiang Chen, and Daniel Malinsky.

Examples

```
# simulate linear Gaussian data w unobserved variable L1
set.seed(123)
n <- 100
L1 <- rnorm(n)
X1 <- rnorm(n)
```



```

X2 <- L1 + X1 + rnorm(n)
X3 <- X1 + rnorm(n)
X4 <- X3 + L1 + rnorm(n)
d <- data.frame(p1_X1 = X1,
                p1_X2 = X2,
                p2_X3 = X3,
                p2_X4 = X4)

# use tfci algorithm to recover tpag (conservative edge orientation)
tfci(d, test = corTest, order = c("p1", "p2"))

# use tfci with standard (non-conservative) method for edge orientation
tfci(d, test = corTest, order = c("p1", "p2"), methodOri = "standard")

```

tges	<i>Estimate the restricted Markov equivalence class using Temporal Greedy Equivalence Search</i>
------	--------------------------------------------------------------------------------------------------

Description

Perform causal discovery using the temporal greedy equivalence search algorithm.

Usage

```
tges(score, verbose = FALSE)
```

Arguments

score	tiered scoring object to be used. At the moment only scores supported re TemporalBIC and TemporalBDeu .
verbose	indicates whether debug output should be printed.

Value

tges returns a tamat object which is a matrix with a "order" attribute (a character vector listing the temporal order of the variables in the adjacency matrix).

Author(s)

Tobias Ellegaard Larsen

See Also

[TemporalBIC](#)

Examples

```
#Simulate Gaussian data
set.seed(123)
n <- 500
child_x <- rnorm(n)
child_y <- 0.5*child_x + rnorm(n)
child_z <- 2*child_x + child_y + rnorm(n)

adult_x <- child_x + rnorm(n)
adult_z <- child_z + rnorm(n)
adult_w <- 2*adult_z + rnorm(n)
adult_y <- 2*child_x + adult_w + rnorm(n)

simdata <- data.frame(child_x, child_y, child_z,
                      adult_x, adult_z, adult_w,
                      adult_y)

# Define order in prefix way
prefix_order <- c("child", "adult")

# Define TBIC score
t_score <- new("TemporalBIC", order = prefix_order
              , data = simdata)

# Run tges
tges_pre <- tges(t_score)

# Plot MPDAG
plot(tges_pre)

# Define order in integer way
integer_order <- c(1,1,1,2,2,2,2)

# Define TBIC score
t_score <- new("TemporalBIC", order = integer_order
              , data = simdata)

# Run tges
tges_int <- tges(t_score)

# Plot MPDAG
plot(tges_int)
```

tpc

Perform causal discovery using the temporal PC algorithm (TPC)

Description

Perform causal discovery using the temporal PC algorithm (TPC)

Usage

```

tpc(
  data = NULL,
  order,
  sparsity = 10^(-1),
  test = regTest,
  suffStat = NULL,
  method = "stable.fast",
  methodNA = "none",
  methodOri = "conservative",
  output = "tpdag",
  varnames = NULL,
  ...
)

```

Arguments

data	A data.frame with data. All variables should be assigned to exactly one period by prefixing them with the period name (see example below).
order	A character vector with period-prefixes in their temporal order (see example below).
sparsity	The sparsity level to be used for independence testing (i.e. significance level threshold to use for each test).
test	A procedure for testing conditional independence. The default, <code>regTest</code> uses a regression-based information loss test. Another available option is <code>corTest</code> which tests for vanishing partial correlations. User supplied functions may also be used, see details below about the required syntax.
suffStat	Sufficient statistic. If this argument is supplied, the sufficient statistic is not computed from the inputted data. The format and contents of the sufficient statistic depends on which test is being used.
method	Which method to use for skeleton construction, must be "stable", "original", or "stable.fast" (the default). See skeleton for details.
methodNA	Method for handling missing information (NA values). Must be one of "none" (default, an error is thrown if NAs are present), "cc" (complete case analysis, deletes all observations that have any NA values), or "twd" (test wise deletion, omits observations with missing information test-by-test) (further details below).
methodOri	Method for handling conflicting separating sets when orienting edges. Currently only the conservative method is available.
output	One of "tpdag", "tskeleton" or "pcAlgo". If "tskeleton", a temporal skeleton is constructed and outputted, but the edges are not directed. If "tpdag" (the default), a the edges are directed, resulting in a temporal partially directed acyclic graph (TPDAG). If "pcAlgo" the TPDAG is outputted as the object class pcAlgo-class from the <code>pcalg</code> package. This is intended for compatability with tools from that package.

varnames	A character vector of variable names. It only needs to be supplied if the data argument is not used, and data are hence passed exclusively through the suffStat argument.
...	Further optional arguments which are passed to skeleton for the skeleton constructing phase.

Details

Note that all independence test procedures implemented in the `pca1g` package may be used, see [pc](#).

The methods for handling missing information require that the data, rather than the suffStat argument is used for inputting data; the latter assumes no missing information and hence always sets `methodNA = "none"`. If the test is `corTest`, test-wise deletion is performed when computing the sufficient statistic (correlation matrix) (so for each pair of variables, only complete cases are used). If the test is `regTest`, test-wise deletion is performed for each conditional independence test instead.

Value

A `tpdag` or `tskeleton` object. Both return types are S3 objects, i.e., lists with entries: `$amat` (the estimated adjacency matrix), `$order` (character vector with the order, as inputted to this function), `$psi` (the significance level used for testing), and `$ntests` (the number of tests conducted).

Examples

```
#TPC on included example data, use sparsity psi = 0.01, default test (regression-based
#information loss):
data(tpcExample)
tpc(tpcExample, order = c("child", "youth", "oldage"), sparsity = 0.01)
```

```
#TPC on included example data, use sparsity psi = 0.01, use test for vanishing partial
# correlations:
data(tpcExample)
tpc(tpcExample, order = c("child", "youth", "oldage"), sparsity = 0.01,
test = corTest)
```

```
#TPC on another simulated data set
```

```
#Simulate data
set.seed(123)
n <- 500
child_x <- rnorm(n)^2
child_y <- 0.5*child_x + rnorm(n)
child_z <- sample(c(0,1), n, replace = TRUE,
                 prob = c(0.3, 0.7))

adult_x <- child_x + rnorm(n)
adult_z <- as.numeric(child_z + rnorm(n) > 0)
adult_w <- 2*adult_z + rnorm(n)
adult_y <- 2*sqrt(child_x) + adult_w^2 + rnorm(n)
```

```

simdata <- data.frame(child_x, child_y, child_z,
                     adult_x, adult_z, adult_w,
                     adult_y)

#Define order
simorder <- c("child", "adult")

#Perform TPC with sparsity psi = 0.001
results <- tpc(simdata, order = simorder, sparsity = 10^(-3))

```

tpcExample

Simulated data example

Description

A small simulated data example intended to showcase the TPC algorithm. Note that the variable name prefixes defines with period they are related to ("child", "youth" or "oldage").

Usage

```
tpcExample
```

Format

A data.frame with 200 rows and 6 variables.

child_x1 Structural equation: $X1 := \epsilon1$ with $\epsilon1 \text{ Unif}(0, 1)$

child_x2 Structural equation: $X2 := 2 * X1 + \epsilon2$ with $\epsilon2 \text{ N}(0, 1)$

youth_x3 Structural equation: $X3 := \epsilon3$ with $\epsilon3 \text{ Unif}(0, 1)$

youth_x4 Structural equation: $X4 := X2 + \epsilon4$ with $\epsilon4 \text{ N}(0, 1)$

oldage_x5 Structural equation: $X5 := X3^2 + X3 - 3 * X2 + \epsilon5$ with $\epsilon5 \text{ N}(0, 1)$

oldage_x6 Structural equation: $X6 := X4^3 + X4^2 + 2 * X5 + \epsilon6$ with $\epsilon6 \text{ N}(0, 1)$

References

Petersen, AH; Osler, M and Ekstrøm, CT (2021): Data-Driven Model Building for Life-Course Epidemiology, American Journal of Epidemiology.

Examples

```
data(tpcExample)
```

tplot	<i>Plot temporal graph via Latex</i>
-------	--------------------------------------

Description

Generates a plot of a `tamat`, `tpdag` or `tpag` object by use of Latex `tikz`. Note that a working Latex installation is required. Note also that this function is slower than typical R graphics options and may take some time to terminate.

Usage

```
tplot(
  x,
  filename = "causaldisco_tplot_temp",
  keepfiles = FALSE,
  bendedges = TRUE,
  ...
)
```

Arguments

<code>x</code>	A <code>tamat</code> , <code>tpdag</code> , or <code>tpag</code> object as obtained from tamat , tpc , or tfci , respectively.
<code>filename</code>	Name of files that will be used internally during the function's runtime. This filename will be appended with both <code>.rmd</code> and <code>.pdf</code> . Note that unless <code>keepfiles = TRUE</code> , these files will automatically be deleted again.
<code>keepfiles</code>	If <code>FALSE</code> (default), temporary files used for making the plot are deleted, otherwise they are kept and will be placed in the working directory.
<code>bendedges</code>	If <code>TRUE</code> (default), all edges are bent 10 degrees to the right, thereby avoiding edges placed exactly on top of eachother.
<code>...</code>	Additional argument passed to maketikz .

Details

The function renders Latex code using `rmarkdown`, which relies on a working installation of Latex. Afterwards, the resulting pdf graphic is loaded into R and displayed in a browser. If working in Rstudio it may be opened in the built-in viewer, depending on Rstudio global settings.

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