

Package ‘CoGAPS’

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Title Coordinated Gene Activity in Pattern Sets

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Description Coordinated Gene Activity in Pattern Sets (CoGAPS) implements a Bayesian MCMC matrix factorization algorithm, GAPS, and links it to gene set statistic methods to infer biological process activity. It can be used to perform sparse matrix factorization on any data, and when this data represents biomolecules, to do gene set analysis.

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CoGAPS-package

*CoGAPS: Coordinated Gene Activity in Pattern Sets***Description**

CoGAPS implements a Bayesian MCMC matrix factorization algorithm, GAPS, and links it to gene set statistic methods to infer biological process activity. It can be used to perform sparse matrix factorization on any data, and when this data represents biomolecules, to do gene set analysis.

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References

Fertig EJ, Ding J, Favorov AV, Parmigiani G, Ochs MF. CoGAPS: an R/C++ package to identify patterns and biological process activity in transcriptomic data. *Bioinformatics*. 2010 Nov 1;26(21):2792-3

binaryA

*binary heatmap for standardized feature matrix***Description**

creates a binarized heatmap of the A matrix in which the value is 1 if the value in Amean is greater than threshold * Asd and 0 otherwise

Usage

```
binaryA(object, threshold = 3)

## S4 method for signature 'CogapsResult'
binaryA(object, threshold = 3)
```

Arguments

object an object of type CogapsResult
 threshold the number of standard deviations above zero that an element of Amean must be to get a value of 1

Value

plots a heatmap of the A Matrix

Examples

```
data(GIST)
# to expensive to call since it plots
# binaryA(GIST.result, threshold=3)
```

buildReport

Information About Package Compilation

Description

Information About Package Compilation

Usage

```
buildReport()
```

Details

returns information about how the package was compiled, i.e. which compiler/version was used, which compile time options were enabled, etc...

Value

string containing build report

Examples

```
CoGAPS::buildReport()
```

calcCoGAPSStat

calculate statistic on sets of measurements (genes) or samples

Description

calculates a statistic to determine if a pattern is enriched in a a particular set of measurements or samples.

Usage

```
calcCoGAPStat(  
  object,  
  sets = NULL,  
  whichMatrix = "featureLoadings",  
  numPerm = 1000,  
  ...  
)  
  
## S4 method for signature 'CogapsResult'  
calcCoGAPStat(  
  object,  
  sets = NULL,  
  whichMatrix = "featureLoadings",  
  numPerm = 1000,  
  ...  
)
```

Arguments

object	an object of type CogapsResult
sets	list of sets of measurements/samples
whichMatrix	either "featureLoadings" or "sampleFactors" indicating which matrix to calculate the statistics for
numPerm	number of permutations to use when calculatin p-value
...	handles old arguments for backwards compatibility

Value

gene set statistics for each column of A

calcGeneGSStat	<i>probability gene belongs in gene set</i>
----------------	---

Description

calculates the probability that a gene listed in a gene set behaves like other genes in the set within the given data set

Usage

```
calcGeneGSStat(  
  object,  
  GStoGenes,  
  numPerm,  
  Pw = rep(1, ncol(object@featureLoadings)),
```

```

    nullGenes = FALSE
  )

  ## S4 method for signature 'CogapsResult'
  calcGeneGSStat(
    object,
    GStoGenes,
    numPerm,
    Pw = rep(1, ncol(object@featureLoadings)),
    nullGenes = FALSE
  )

```

Arguments

object	an object of type CogapsResult
GStoGenes	data.frame or list with gene sets
numPerm	number of permutations for null
Pw	weight on genes
nullGenes	logical indicating gene adjustment

Value

gene similiarity statistic

calcZ	<i>compute z-score matrix</i>
-------	-------------------------------

Description

calculates the Z-score for each element based on input mean and standard deviation matrices

Usage

```

calcZ(object, whichMatrix)

## S4 method for signature 'CogapsResult'
calcZ(object, whichMatrix)

```

Arguments

object	an object of type CogapsResult
whichMatrix	either "featureLoadings" or "sampleFactors" indicating which matrix to calculate the z-score for

Value

matrix of z-scores

Examples

```
data(GIST)
featureZScore <- calcZ(GIST.result, "featureLoadings")
```

callInternalCoGAPS	<i>make correct call to internal CoGAPS dispatch function, CoGAPS could be called directly, but to avoid any re-entrant behavior this function is called instead. It is a light wrapper around cogaps_cpp that handles setting the distributed parameters</i>
--------------------	---

Description

make correct call to internal CoGAPS dispatch function, CoGAPS could be called directly, but to avoid any re-entrant behavior this function is called instead. It is a light wrapper around cogaps_cpp that handles setting the distributed parameters

Usage

```
callInternalCoGAPS(data, allParams, uncertainty, subsetIndices, workerID)
```

Arguments

data	data in a supported format
allParams	list of all parameters
uncertainty	uncertainty of data in the same format as data
subsetIndices	indices of the subset of data to run on
workerID	worker ID for parallelization

Value

CogapsResult object

checkDataMatrix	<i>check that provided data is valid</i>
-----------------	--

Description

check that provided data is valid

Usage

```
checkDataMatrix(data, uncertainty, params)
```

Arguments

data	data matrix
uncertainty	uncertainty matrix, can be null
params	CogapsParams object

Value

throws an error if data has problems

checkInputs	<i>check that all inputs are valid</i>
-------------	--

Description

check that all inputs are valid

Usage

```
checkInputs(data, uncertainty, allParams)
```

Arguments

data	data matrix
uncertainty	uncertainty matrix, can be null
allParams	list of all parameters

Value

throws an error if inputs are invalid

checkpointsEnabled	<i>Check if package was built with checkpoints enabled</i>
--------------------	--

Description

Check if package was built with checkpoints enabled

Usage

```
checkpointsEnabled()
```

Value

true/false if checkpoints are enabled

Examples

```
CoGAPS::checkpointsEnabled()
```

Description

calls the C++ MCMC code and performs Bayesian matrix factorization returning the two matrices that reconstruct the data matrix

Usage

```
CoGAPS(
  data,
  params = new("CogapsParams", nPatterns = nPatterns),
  nPatterns,
  nThreads = 1,
  messages = TRUE,
  outputFrequency = 1000,
  uncertainty = NULL,
  checkpointOutFile = "gaps_checkpoint.out",
  checkpointInterval = 0,
  checkpointInFile = NULL,
  transposeData = FALSE,
  BPPARAM = NULL,
  workerID = 1,
  asynchronousUpdates = TRUE,
  nSnapshots = 0,
  snapshotPhase = "sampling",
  ...
)
```

Arguments

<code>data</code>	File name or R object (see details for supported types)
<code>params</code>	CogapsParams object
<code>nPatterns</code>	rank of the nmf decomposition
<code>nThreads</code>	maximum number of threads to run on
<code>messages</code>	T/F for displaying output
<code>outputFrequency</code>	number of iterations between each output (set to 0 to disable status updates, other output is controlled by @code messages)
<code>uncertainty</code>	uncertainty matrix - either a matrix or a supported file type
<code>checkpointOutFile</code>	name of the checkpoint file to create
<code>checkpointInterval</code>	number of iterations between each checkpoint (set to 0 to disable checkpoints)

checkpointInFile	if this is provided, CoGAPS runs from the checkpoint contained in this file
transposeData	T/F for transposing data while reading it in - useful for data that is stored as samples x genes since CoGAPS requires data to be genes x samples
BPPARAM	BiocParallel backend
workerID	if calling CoGAPS in parallel the worker ID can be specified, only worker 1 prints output and each worker outputs when it finishes, this is not necessary when using the default parallel methods (i.e. distributed CoGAPS) but only when the user is manually calling CoGAPS in parallel
asynchronousUpdates	enable asynchronous updating which allows for multi-threaded runs
nSnapshots	how many snapshots to take in each phase, setting this to 0 disables snapshots
snapshotPhase	which phase to take snapshots in e.g. "equilibration", "sampling", "all"
...	allows for overwriting parameters in params

Details

The supported R types are: matrix, data.frame, SummarizedExperiment, SingleCellExperiment. The supported file types are csv, tsv, and mtx.

Value

CogapsResult object

Examples

```
# Running from R object
data(GIST)
resultA <- CoGAPS(GIST.data_frame, nPatterns=3, nIterations=25)

# Running from file name
gist_path <- system.file("extdata/GIST.mtx", package="CoGAPS")
resultB <- CoGAPS(gist_path, nPatterns=3, nIterations=25)

# Setting Parameters
params <- new("CogapsParams", nPatterns=3)
params <- setParam(params, "nIterations", 25)
resultC <- CoGAPS(GIST.data_frame, params)
```

CogapsParams

CogapsParams constructor

Description

create a CogapsParams object

Usage

```
CogapsParams(...)
```

Arguments

... parameters for the initialization method

Value

CogapsParams object

Examples

```
params <- CogapsParams(nPatterns=10)
params
```

CogapsParams-class	<i>CogapsParams</i>
--------------------	---------------------

Description

Encapsulates all parameters for the CoGAPS algorithm

Slots

nPatterns number of patterns CoGAPS will learn
 nIterations number of iterations for each phase of the algorithm
 alphaA sparsity parameter for feature matrix
 alphaP sparsity parameter for sample matrix
 maxGibbsMassA atomic mass restriction for feature matrix
 maxGibbsMassP atomic mass restriction for sample matrix
 seed random number generator seed
 sparseOptimization speeds up performance with sparse data (roughly >80 default uncertainty
 distributed either "genome-wide" or "single-cell" indicating which distributed algorithm should
 be used
 nSets [distributed parameter] number of sets to break data into
 cut [distributed parameter] number of branches at which to cut dendrogram used in pattern match-
 ing
 minNS [distributed parameter] minimum of individual set contributions a cluster must contain
 maxNS [distributed parameter] maximum of individual set contributions a cluster can contain
 explicitSets [distributed parameter] specify subsets by index or name
 samplingAnnotation [distributed parameter] specify categories along the rows (cols) to use for
 weighted sampling

`samplingWeight` [distributed parameter] weights associated with `samplingAnnotation`
`subsetIndices` set of indices to use from the data
`subsetDim` which dimension (1=rows, 2=cols) to subset
`geneNames` vector of names of genes in data
`sampleNames` vector of names of samples in data
`fixedPatterns` fix either 'A' or 'P' matrix to these values, in the context of distributed CoGAPS (GWCoGAPS/scCoGAPS), the first phase is skipped and `fixedPatterns` is used for all sets - allowing manual pattern matching, as well as fixed runs of standard CoGAPS
`whichMatrixFixed` either 'A' or 'P', indicating which matrix is fixed
`takePumpSamples` whether or not to take PUMP samples
`checkpointInterval` how many iterations between each checkpoint (set to 0 to disable)
`checkpointInFile` file path to load checkpoint from
`checkpointOutFile` file path where checkpoint should be written to

CogapsResult-class *CogapsResult*

Description

Contains all output from Cogaps run

Slots

`factorStdDev` std dev of the sampled P matrices
`loadingStdDev` std dev of the sampled A matrices

`compiledWithOpenMPSupport`
Check if compiler supported OpenMP

Description

Check if compiler supported OpenMP

Usage

`compiledWithOpenMPSupport()`

Value

true/false if OpenMP was supported

Examples

`CoGAPS::compiledWithOpenMPSupport()`

computeGeneGSProb	<i>compute gene probability</i>
-------------------	---------------------------------

Description

Computes the p-value for gene set membership using the CoGAPS-based statistics developed in Fertig et al. (2012). This statistic refines set membership for each candidate gene in a set specified in GSGenes by comparing the inferred activity of that gene to the average activity of the set.

Usage

```
computeGeneGSProb(
  object,
  GStoGenes,
  numPerm = 500,
  Pw = rep(1, ncol(object@featureLoadings)),
  PwNull = FALSE
)

## S4 method for signature 'CogapsResult'
computeGeneGSProb(
  object,
  GStoGenes,
  numPerm = 500,
  Pw = rep(1, ncol(object@featureLoadings)),
  PwNull = FALSE
)
```

Arguments

object	an object of type CogapsResult
GStoGenes	data.frame or list with gene sets
numPerm	number of permutations for null
Pw	weight on genes
PwNull	- logical indicating gene adjustment

Value

A vector of length GSGenes containing the p-values of set membership for each gene contained in the set specified in GSGenes.

convertDataToMatrix	<i>convert any acceptable data input to a numeric matrix</i>
---------------------	--

Description

convert supported R objects containing the data to a numeric matrix, if data is a file name do nothing. Exits with an error if data is not a supported type.

Usage

convertDataToMatrix(data)

Arguments

data data input

Value

data matrix

corcut	<i>cluster patterns together</i>
--------	----------------------------------

Description

cluster patterns together

Usage

corcut(allPatterns, cut, minNS)

Arguments

allPatterns matrix of all patterns across subsets
cut number of branches at which to cut dendrogram
minNS minimum of individual set contributions a cluster must contain

Value

patterns listed by which cluster they belong to

corrToMeanPattern	<i>calculate correlation of each pattern in a cluster to the cluster mean</i>
-------------------	---

Description

calculate correlation of each pattern in a cluster to the cluster mean

Usage

```
corrToMeanPattern(cluster)
```

Value

correlation of each pattern

createCogapsResult	<i>convert list output from c++ code to a CogapsResult object</i>
--------------------	---

Description

convert list output from c++ code to a CogapsResult object

Usage

```
createCogapsResult(returnList, allParams)
```

Arguments

returnList	list from cogaps_cpp
allParams	list of all parameters

Value

CogapsResult object

createSets	<i>partition genes/samples into subsets</i>
------------	---

Description

either genes or samples or partitioned depending on the type of distributed CoGAPS (i.e. genome-wide or single-cell)

Usage

```
createSets(data, allParams)
```

Arguments

data	either file name or matrix
allParams	list of all CoGAPS parameters

Value

list of sorted subsets of either genes or samples

distributedCogaps	<i>CoGAPS Distributed Matrix Factorization Algorithm</i>
-------------------	--

Description

runs CoGAPS over subsets of the data and stitches the results back together

Usage

```
distributedCogaps(data, allParams, uncertainty)
```

Arguments

data	File name or R object (see details for supported types)
allParams	list of all parameters used in computation
uncertainty	uncertainty matrix (same supported types as data)

Details

For file types CoGAPS supports csv, tsv, and mtx

Value

list

findConsensusMatrix	<i>find the consensus pattern matrix across all subsets</i>
---------------------	---

Description

find the consensus pattern matrix across all subsets

Usage

```
findConsensusMatrix(unmatchedPatterns, gapsParams)
```

Arguments

unmatchedPatterns	list of all unmatched pattern matrices from initial run of CoGAPS
gapsParams	list of all CoGAPS parameters

Value

matrix of consensus patterns

fromCSV	<i>read CoGAPS Result object from a directory with a set of csvs see toCSV</i>
---------	--

Description

save as csv

Usage

```
fromCSV(save_location = ".")

## S4 method for signature 'character'
fromCSV(save_location = ".")
```

Arguments

save_location	directory to read from
---------------	------------------------

Value

CogapsResult object

gapsCat	<i>wrapper around cat</i>
---------	---------------------------

Description

cleans up message printing

Usage

```
gapsCat(allParams, ...)
```

Arguments

allParams	all cogaps parameters
...	arguments forwarded to cat

Value

conditionally print message

getAmplitudeMatrix	<i>return Amplitude matrix from CogapsResult object</i>
--------------------	---

Description

return Amplitude matrix from CogapsResult object

Usage

```
getAmplitudeMatrix(object)

## S4 method for signature 'CogapsResult'
getAmplitudeMatrix(object)
```

Arguments

object	an object of type CogapsResult
--------	--------------------------------

Value

amplitude matrix

Examples

```
data(GIST)
amplitudeMatrix <- getAmplitudeMatrix(GIST.result)
```

getClusteredPatterns *return clustered patterns from set of all patterns across all subsets*

Description

return clustered patterns from set of all patterns across all subsets

Usage

```
getClusteredPatterns(object)

## S4 method for signature 'CogapsResult'
getClusteredPatterns(object)
```

Arguments

object an object of type CogapsResult

Value

CogapsParams object

Examples

```
data(GIST)
clusteredPatterns <- getClusteredPatterns(GIST.result)
```

getCorrelationToMeanPattern
 return correlation between each pattern and the cluster mean

Description

return correlation between each pattern and the cluster mean

Usage

```
getCorrelationToMeanPattern(object)

## S4 method for signature 'CogapsResult'
getCorrelationToMeanPattern(object)
```

Arguments

object an object of type CogapsResult

Value

CogapsParams object

Examples

```
data(GIST)
corrToMeanPattern <- getCorrelationToMeanPattern(GIST.result)
```

getDimNames	<i>extracts gene/sample names from the data</i>
-------------	---

Description

extracts gene/sample names from the data

Usage

```
getDimNames(data, allParams)
```

Arguments

data	data matrix
allParams	list of all parameters

Value

list of all parameters with added gene names

getFeatureLoadings	<i>return featureLoadings matrix from CogapsResult object</i>
--------------------	---

Description

return featureLoadings matrix from CogapsResult object

Usage

```
getFeatureLoadings(object)

## S4 method for signature 'CogapsResult'
getFeatureLoadings(object)
```

Arguments

object	an object of type CogapsResult
--------	--------------------------------

Value

featureLoadings matrix

Examples

```
data(GIST)
fLoadings <- getFeatureLoadings(GIST.result)
```

getGeneNames	<i>extract gene names from data</i>
--------------	-------------------------------------

Description

extract gene names from data

Usage

```
getGeneNames(data, transpose)
```

Value

vector of gene names

getMeanChiSq	<i>return chi-sq of final matrices</i>
--------------	--

Description

return chi-sq of final matrices

Usage

```
getMeanChiSq(object)

## S4 method for signature 'CogapsResult'
getMeanChiSq(object)
```

Arguments

object	an object of type CogapsResult
--------	--------------------------------

Value

chi-sq error

Examples

```
data(GIST)
getMeanChiSq(GIST.result)
```

getOriginalParameters *return original parameters used to generate this result*

Description

return original parameters used to generate this result

Usage

```
getOriginalParameters(object)

## S4 method for signature 'CogapsResult'
getOriginalParameters(object)
```

Arguments

object an object of type CogapsResult

Value

CogapsParams object

Examples

```
data(GIST)
params <- getOriginalParameters(GIST.result)
```

getParam *get the value of a parameter*

Description

get the value of a parameter

Usage

```
getParam(object, whichParam)

## S4 method for signature 'CogapsParams'
getParam(object, whichParam)
```

Arguments

object an object of type CogapsParams
whichParam a string with the name of the requested parameter

Value

the value of the parameter

Examples

```
params <- new("CogapsParams", nPatterns=3)
getParam(params, "seed")
```

getPatternGeneSet	<i>generate statistics associating patterns with gene sets</i>
-------------------	--

Description

generate statistics associating patterns with gene sets

Usage

```
getPatternGeneSet(
  object,
  gene.sets,
  method = c("enrichment", "overrepresentation"),
  ...
)

## S4 method for signature 'CogapsResult,list,character'
getPatternGeneSet(
  object,
  gene.sets,
  method = c("enrichment", "overrepresentation"),
  ...
)
```

Arguments

object	an object of type CogapsResult
gene.sets	a list of gene sets to test. List names should be the names of the gene sets
method	enrichment or overrepresentation. Conducts a test for gene set enrichment using fgsea::gsea ranking features by pattern amplitude or a test for gene set overrepresentation in pattern markers using fgsea::fora, respectively.
...	additional parameters passed to patternMarkers if using overrepresentation method

Value

list of dataframes containing gene set enrichment or gene set overrepresentation statistics

Examples

```
data(GIST)
gs.test <- list(
  "gs1" = c("Hs.2", "Hs.4", "Hs.36", "Hs.96", "Hs.202"),
  "gs2" = c("Hs.699463", "Hs.699288", "Hs.699280", "Hs.699154", "Hs.697294")
)
getPatternGeneSet(object = GIST.result, gene.sets = gs.test, method = "enrichment")
getPatternGeneSet(object = GIST.result, gene.sets = gs.test, method = "overrepresentation")
```

getPatternMatrix	<i>return pattern matrix from CogapsResult object</i>
------------------	---

Description

return pattern matrix from CogapsResult object

Usage

```
getPatternMatrix(object)

## S4 method for signature 'CogapsResult'
getPatternMatrix(object)
```

Arguments

object an object of type CogapsResult

Value

pattern matrix

Examples

```
data(GIST)
patternMatrix <- getPatternMatrix(GIST.result)
```

getRetinaSubset	<i>get specified number of retina subsets</i>
-----------------	---

Description

combines retina subsets from extdata directory

Usage

```
getRetinaSubset(n = 1)
```

Arguments

n number of subsets to use

Value

matrix of RNA counts

Examples

```
retSubset <- getRetinaSubset()
dim(retSubset)
```

getSampleFactors	<i>return sampleFactors matrix from CogapsResult object</i>
------------------	---

Description

return sampleFactors matrix from CogapsResult object

Usage

```
getSampleFactors(object)

## S4 method for signature 'CogapsResult'
getSampleFactors(object)
```

Arguments

object an object of type CogapsResult

Value

sampleFactors matrix

Examples

```
data(GIST)
sFactors <- getSampleFactors(GIST.result)
```

getSampleNames	<i>extract sample names from data</i>
----------------	---------------------------------------

Description

extract sample names from data

Usage

```
getSampleNames(data, transpose)
```

Value

vector of sample names

getSubsets	<i>return the names of the genes (samples) in each subset</i>
------------	---

Description

return the names of the genes (samples) in each subset

Usage

```
getSubsets(object)  
  
## S4 method for signature 'CogapsResult'  
getSubsets(object)
```

Arguments

object	an object of type CogapsResult
--------	--------------------------------

Value

CogapsParams object

Examples

```
data(GIST)  
subsets <- getSubsets(GIST.result)
```

getUnmatchedPatterns *return unmatched patterns from each subset*

Description

return unmatched patterns from each subset

Usage

```
getUnmatchedPatterns(object)

## S4 method for signature 'CogapsResult'
getUnmatchedPatterns(object)
```

Arguments

object an object of type CogapsResult

Value

CogapsParams object

Examples

```
data(GIST)
unmatchedPatterns <- getUnmatchedPatterns(GIST.result)
```

getValueOrRds *get input that might be an RDS file*

Description

get input that might be an RDS file

Usage

```
getValueOrRds(input)
```

Arguments

input some user input

Value

if input is an RDS file, read it - otherwise return input

getVersion	<i>return version number used to generate this result</i>
------------	---

Description

return version number used to generate this result

Usage

```
getVersion(object)

## S4 method for signature 'CogapsResult'
getVersion(object)
```

Arguments

object an object of type CogapsResult

Value

version number

Examples

```
data(GIST)
getVersion(GIST.result)
```

GIST.data_frame	<i>GIST gene expression data from Ochs et al. (2009)</i>
-----------------	--

Description

GIST gene expression data from Ochs et al. (2009)

GIST.matrix	<i>GIST gene expression data from Ochs et al. (2009)</i>
-------------	--

Description

GIST gene expression data from Ochs et al. (2009)

GIST.result

CoGAPS result from running on GIST dataset

Description

CoGAPS result from running on GIST dataset

GIST.uncertainty

GIST gene expression uncertainty matrix from Ochs et al. (2009)

Description

GIST gene expression uncertainty matrix from Ochs et al. (2009)

GWCoGAPS

Genome Wide CoGAPS

Description

wrapper around genome-wide distributed algorithm for CoGAPS

Usage

```
GWCoGAPS(
  data,
  params = new("CogapsParams"),
  nThreads = 1,
  messages = TRUE,
  outputFrequency = 500,
  uncertainty = NULL,
  checkpointOutFile = "gaps_checkpoint.out",
  checkpointInterval = 1000,
  checkpointInFile = NULL,
  transposeData = FALSE,
  BPPARAM = NULL,
  workerID = 1,
  asynchronousUpdates = FALSE,
  ...
)
```

Arguments

<code>data</code>	File name or R object (see details for supported types)
<code>params</code>	CogapsParams object
<code>nThreads</code>	maximum number of threads to run on
<code>messages</code>	T/F for displaying output
<code>outputFrequency</code>	number of iterations between each output (set to 0 to disable status updates, other output is controlled by @code messages)
<code>uncertainty</code>	uncertainty matrix - either a matrix or a supported file type
<code>checkpointOutFile</code>	name of the checkpoint file to create
<code>checkpointInterval</code>	number of iterations between each checkpoint (set to 0 to disable checkpoints)
<code>checkpointInFile</code>	if this is provided, CoGAPS runs from the checkpoint contained in this file
<code>transposeData</code>	T/F for transposing data while reading it in - useful for data that is stored as samples x genes since CoGAPS requires data to be genes x samples
<code>BPPARAM</code>	BiocParallel backend
<code>workerID</code>	if calling CoGAPS in parallel the worker ID can be specified, only worker 1 prints output and each worker outputs when it finishes, this is not necessary when using the default parallel methods (i.e. distributed CoGAPS) but only when the user is manually calling CoGAPS in parallel
<code>asynchronousUpdates</code>	enable asynchronous updating which allows for multi-threaded runs
<code>...</code>	allows for overwriting parameters in params

Value

CogapsResult object

Examples

```
## Not run:
data(GIST)
params <- new("CogapsParams", nPatterns=3)
params <- setDistributedParams(params, nSets=2)
params <- setParam(params, "nIterations", 100)
result <- GWCoGAPS(GIST.matrix, params, BPPARAM=BiocParallel::SerialParam())

## End(Not run)
```

```
initialize,CogapsParams-method
      constructor for CogapsParams
```

Description

constructor for CogapsParams

Usage

```
## S4 method for signature 'CogapsParams'
initialize(.Object, distributed = NULL, nPatterns, ...)
```

Arguments

.Object	CogapsParams object
distributed	either "genome-wide" or "single-cell" indicating which distributed algorithm should be used
nPatterns	number of patterns
...	initial values for slots

Value

initialized CogapsParams object

```
initialize,CogapsResult-method
      Constructor for CogapsResult
```

Description

Constructor for CogapsResult

Usage

```
## S4 method for signature 'CogapsResult'
initialize(
  .Object,
  Amean,
  Pmean,
  Asd,
  Psd,
  meanChiSq,
  geneNames,
```

```

    sampleNames,
    diagnostics = NULL,
    ...
)

```

Arguments

.Object	CogapsResult object
Amean	mean of sampled A matrices
Pmean	mean of sampled P matrices
Asd	std dev of sampled A matrices
Psd	std dev of sampled P matrices
meanChiSq	mean value of ChiSq statistic
geneNames	names of genes in data
sampleNames	names of samples in data
diagnostics	assorted diagnostic reports from the run
...	initial values for slots

Value

initialized CogapsResult object

isRdsFile	<i>checks if file is rds format</i>
-----------	-------------------------------------

Description

checks if file is rds format

Usage

```
isRdsFile(file)
```

Arguments

file	path to file
------	--------------

Value

TRUE if file is .rds, FALSE if not

MANOVA

MANOVA statistical test for patterns between sample groups

Description

MANOVA statistical test—wraps base R manova

Usage

```
MANOVA(interestedVariables, object)
```

```
## S4 method for signature 'matrix,CogapsResult'
MANOVA(interestedVariables, object)
```

Arguments

interestedVariables
study design for manova

object
CogapsResult object

Value

list of manova fit results

modsimdata

Toy example to run CoGAPS on.

Description

- V1..V20. some variables, for example levels of gene expression

Usage

```
data(modsimdata)
```

Format

'data.frame': 25 obs. of 20 variables.

modsimresult	<i>Result of applying CoGAPS on the Toy example.</i>
--------------	--

Description

Result of applying CoGAPS on the Toy example.

Usage

```
data(modsimresult)
```

Format

S4 class ‘CogapsResult’ [package “CoGAPS”] with 7 slots.

ncolHelper	<i>get number of columns from supported file name or matrix</i>
------------	---

Description

get number of columns from supported file name or matrix

Usage

```
ncolHelper(data)
```

Arguments

data	either a file name or a matrix
------	--------------------------------

Value

number of columns

nrowHelper	<i>get number of rows from supported file name or matrix</i>
------------	--

Description

get number of rows from supported file name or matrix

Usage

```
nrowHelper(data)
```

Arguments

data	either a file name or a matrix
------	--------------------------------

Value

number of rows

parseExtraParams	<i>parse parameters passed through the ... variable</i>
------------------	---

Description

parse parameters passed through the ... variable

Usage

```
parseExtraParams(allParams, extraParams)
```

Arguments

allParams	list of all parameters
extraParams	list of parameters in ...

Value

allParams with any valid parameters in extraParams added

Note

will halt with an error if any parameters in extraParams are invalid

patternMarkers	<i>compute pattern markers statistic</i>
----------------	--

Description

estimate the most associated pattern for each feature

Usage

```
patternMarkers(object, threshold = "all", lp = NULL, axis = 1)
```

```
## S4 method for signature 'CogapsResult'
patternMarkers(object, threshold = "all", lp = NULL, axis = 1)
```

Arguments

object	an object of type CogapsResult
threshold	the type of threshold to be used. The default "all" will distribute features into patterns with the highest ranking as ranked by the increasing Euclidian distance between feature loadings and lp. The alternative "cut" will only keep the features that are ranked higher than the first feature having greater intra-pattern compared to inter-pattern rank. This is useful to limit the number of markers ranked similarly everywhere. Features may be present in multiple patterns for "cut".
lp	list of vectors of weights for each pattern to be used for finding markers. If NULL, list of synthetic one-hot markers for each pattern will be generated and matched against.
axis	controls the matrix to use for ranking. 1 for featureLoadings, 2 for sampleFactors.

Value

List of: list of marker features for each pattern (best rank first), a matrix of ranks of each feature in each pattern, a matrix of scores for each feature in each pattern.

List of: list of marker features for each pattern (best rank first), and a matrix of ranks of each feature in each pattern.

Examples

```
data(GIST)
pm <- patternMarkers(GIST.result)
```

patternMatch	<i>Match Patterns Across Multiple Runs</i>
--------------	--

Description

Match Patterns Across Multiple Runs

Usage

```
patternMatch(allPatterns, gapsParams)
```

Arguments

allPatterns	matrix of patterns stored in the columns
gapsParams	CoGAPS parameters object

Value

a matrix of consensus patterns

plotPatternGeneSet	<i>generate a barchart of most significant hallmark sets for a pattern</i>
--------------------	--

Description

generate a barchart of most significant hallmark sets for a pattern

Usage

```
plotPatternGeneSet(patterngeneset, whichpattern = 1, padj_threshold = 0.05)

## S4 method for signature 'list,numeric,numeric'
plotPatternGeneSet(patterngeneset, whichpattern = 1, padj_threshold = 0.05)
```

Arguments

patterngeneset	output from getPatternGeneSet
whichpattern	which pattern to generate bar chart for
padj_threshold	maximum adjusted p-value of gene sets rendered on the resulting plot

Value

image object of barchart

plotPatternMarkers	<i>heatmap of original data clustered by pattern markers statistic</i>
--------------------	--

Description

heatmap of original data clustered by pattern markers statistic

Usage

```
plotPatternMarkers(
  object,
  data,
  patternMarkers,
  patternPalette,
  sampleNames,
  samplePalette = NULL,
  heatmapCol = bluered,
  colDendrogram = TRUE,
  scale = "row",
  ...
)
```

Arguments

object	an object of type CogapsResult
data	the original data as a matrix
patternMarkers	pattern markers to be plotted, as generated by the patternMarkers function
patternPalette	a vector indicating what color should be used for each pattern
sampleNames	names of the samples to use for labeling
samplePalette	a vector indicating what color should be used for each sample
heatmapCol	pallet giving color scheme for heatmap
colDendrogram	logical indicating whether to display sample dendrogram
scale	character indicating if the values should be centered and scaled in either the row direction or the column direction, or none. The default is "row".
...	additional graphical parameters to be passed to heatmap.2

Value

heatmap of the data values for the patternMarkers

See Also

[heatmap.2](#)

plotResiduals	<i>plot of residuals</i>
---------------	--------------------------

Description

calculate residuals and produce heatmap

Usage

```
plotResiduals(object, data, uncertainty = NULL)

## S4 method for signature 'CogapsResult'
plotResiduals(object, data, uncertainty = NULL)
```

Arguments

object	an object of type CogapsResult
data	original data matrix run through GAPS
uncertainty	original standard deviation matrix run through GAPS

Value

creates a residual plot

Examples

```
data(GIST)
# to expensive to call since it plots
# plotResiduals(GIST.result, GIST.matrix)
```

reconstructGene	<i>reconstruct gene</i>
-----------------	-------------------------

Description

reconstruct gene

Usage

```
reconstructGene(object, genes = NULL)

## S4 method for signature 'CogapsResult'
reconstructGene(object, genes = NULL)
```

Arguments

object	an object of type CogapsResult
genes	an index of the gene or genes of interest

Value

the D' estimate of a gene or set of genes

Examples

```
data(GIST)
estimatedD <- reconstructGene(GIST.result)
```

sampleUniformly	<i>subset data by uniformly partitioning rows (cols)</i>
-----------------	--

Description

subset data by uniformly partitioning rows (cols)

Usage

```
sampleUniformly(allParams, total, setSize)
```

Arguments

allParams	list of all CoGAPS parameters
total	total number of rows (cols) that are being partitioned
setSize	the size of each subset of the total

Value

list of subsets

sampleWithAnnotationWeights
<i>subset rows (cols) proportional to the user provided weights</i>

Description

subset rows (cols) proportional to the user provided weights

Usage

sampleWithAnnotationWeights(allParams, setSize)

Arguments

- | | |
|-----------|--------------------------------------|
| allParams | list of all CoGAPS parameters |
| setSize | the size of each subset of the total |

Value

list of subsets

sampleWithExplicitSets	<i>use user provided subsets</i>
------------------------	----------------------------------

Description

use user provided subsets

Usage

sampleWithExplicitSets(allParams)

Arguments

- | | |
|-----------|--|
| allParams | list of all CoGAPS parameters |
| total | total number of rows (cols) that are being partitioned |

Value

list of subsets

scCoGAPS

*Single Cell CoGAPS***Description**

wrapper around single-cell distributed algorithm for CoGAPS

Usage

```
scCoGAPS(
  data,
  params = new("CogapsParams"),
  nThreads = 1,
  messages = TRUE,
  outputFrequency = 500,
  uncertainty = NULL,
  checkpointOutFile = "gaps_checkpoint.out",
  checkpointInterval = 1000,
  checkpointInFile = NULL,
  transposeData = FALSE,
  BPPARAM = NULL,
  workerID = 1,
  asynchronousUpdates = FALSE,
  ...
)
```

Arguments

data	File name or R object (see details for supported types)
params	CogapsParams object
nThreads	maximum number of threads to run on
messages	T/F for displaying output
outputFrequency	number of iterations between each output (set to 0 to disable status updates, other output is controlled by @code messages)
uncertainty	uncertainty matrix - either a matrix or a supported file type
checkpointOutFile	name of the checkpoint file to create
checkpointInterval	number of iterations between each checkpoint (set to 0 to disable checkpoints)
checkpointInFile	if this is provided, CoGAPS runs from the checkpoint contained in this file
transposeData	T/F for transposing data while reading it in - useful for data that is stored as samples x genes since CoGAPS requires data to be genes x samples

BPPARAM	BiocParallel backend
workerID	if calling CoGAPS in parallel the worker ID can be specified, only worker 1 prints output and each worker outputs when it finishes, this is not necessary when using the default parallel methods (i.e. distributed CoGAPS) but only when the user is manually calling CoGAPS in parallel
asynchronousUpdates	enable asynchronous updating which allows for multi-threaded runs
...	allows for overwriting parameters in params

Value

CogapsResult object

Examples

```
## Not run:
data(GIST)
params <- new("CogapsParams", nPatterns=3)
params <- setDistributedParams(params, nSets=2)
params <- setParam(params, "nIterations", 100)
result <- scCoGAPS(t(GIST.matrix), params, BPPARAM=BiocParallel::SerialParam())

## End(Not run)
```

setAnnotationWeights *set the annotation labels and weights for subsetting the data*

Description

these parameters are interrelated so they must be set together

Usage

```
setAnnotationWeights(object, annotation, weights)
```

```
## S4 method for signature 'CogapsParams'
setAnnotationWeights(object, annotation, weights)
```

Arguments

object	an object of type CogapsParams
annotation	vector of labels
weights	vector of weights

Value

the modified params object

Examples

```
params <- new("CogapsParams", nPatterns=3)
params <- setAnnotationWeights(params, c('a', 'b', 'c'), c(1,2,1))
```

setDistributedParams *set the value of parameters for distributed CoGAPS*

Description

these parameters are interrelated so they must be set together

Usage

```
setDistributedParams(
  object,
  nSets = NULL,
  cut = NULL,
  minNS = NULL,
  maxNS = NULL
)

## S4 method for signature 'CogapsParams'
setDistributedParams(
  object,
  nSets = NULL,
  cut = NULL,
  minNS = NULL,
  maxNS = NULL
)
```

Arguments

object	an object of type CogapsParams
nSets	number of sets to break data into
cut	number of branches at which to cut dendrogram used in pattern matching
minNS	minimum of individual set contributions a cluster must contain
maxNS	maximum of individual set contributions a cluster can contain

Value

the modified params object

Examples

```
params <- new("CogapsParams", nPatterns=3)
params <- setDistributedParams(params, 5)
```

setFixedPatterns	<i>set the fixed patterns for either the A or the P matrix</i>
------------------	--

Description

these parameters are interrelated so they must be set together

Usage

```
setFixedPatterns(object, fixedPatterns, whichMatrixFixed)

## S4 method for signature 'CogapsParams'
setFixedPatterns(object, fixedPatterns, whichMatrixFixed)
```

Arguments

object	an object of type CogapsParams
fixedPatterns	values for either the A or P matrix
whichMatrixFixed	either 'A' or 'P' indicating which matrix is fixed

Value

the modified params object

Examples

```
params <- new("CogapsParams", nPatterns=3)
data(GIST)
params <- setFixedPatterns(params, getSampleFactors(GIST.result), 'P')
```

setParam	<i>set the value of a parameter</i>
----------	-------------------------------------

Description

set the value of a parameter

Usage

```
setParam(object, whichParam, value)

## S4 method for signature 'CogapsParams'
setParam(object, whichParam, value)
```

Arguments

object	an object of type CogapsParams
whichParam	a string with the name of the parameter to be changed
value	the value to set the parameter to

Value

the modified params object

Examples

```
params <- new("CogapsParams", nPatterns=3)
params <- setParam(params, "seed", 123)
```

startupMessage	<i>write start up message</i>
----------------	-------------------------------

Description

write start up message

Usage

```
startupMessage(data, allParams)
```

Arguments

data	data set
allParams	list of all parameters

Value

message displayed to screen

stitchTogether	<i>concatenate final results across subsets</i>
----------------	---

Description

concatenate final results across subsets

Usage

```
stitchTogether(result, allParams, sets)
```

Arguments

result	list of CogapsResult object from all runs across subsets
allParams	list of all CoGAPS parameters
sets	indices of sets used to break apart data

Value

list with all CoGAPS output

supported	<i>checks if file is supported</i>
-----------	------------------------------------

Description

checks if file is supported

Usage

```
supported(file)
```

Arguments

file	path to file
------	--------------

Value

TRUE if file is supported, FALSE if not

toCSV	<i>save CoGAPS Result object as a set of csvs to directory see fromCSV</i>
-------	--

Description

save as csv

Usage

```
toCSV(object, save_location = ".")  
  
## S4 method for signature 'CogapsResult,character'  
toCSV(object, save_location = ".")
```

Arguments

object CogapsResult object
save_location directory to write to

Value

none

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