

# Package ‘CoGAPS’

October 12, 2023

**Version** 3.20.0

**Date** 2023-03-18

**Title** Coordinated Gene Activity in Pattern Sets

**Author** Jeanette Johnson, Ashley Tsang, Jacob Mitchell, Thomas Sherman, Wai-shing Lee, Conor Kelton, Ondrej Maxian, Jacob Carey, Genevieve Stein-O’Brien, Michael Considine, Maggie Wodicka, John Stansfield, Shawn Siyy, Carlo Colantuoni, Alexander Favorov, Mike Ochs, Elana Fertig

**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.

**Maintainer** Elana J. Fertig <ejfertig@jhmi.edu>,  
Thomas D. Sherman <tomsherman159@gmail.com>,  
Jeanette Johnson <jjohn450@jhmi.edu>

**Depends** R (>= 3.5.0)

**Imports** BiocParallel, cluster, methods, gplots, graphics, grDevices,  
RColorBrewer, Rcpp, S4Vectors, SingleCellExperiment, stats,  
SummarizedExperiment, tools, utils, rhdf5, biomaRt, dplyr,  
fgsea, forcats, ggplot2, msigdbr

**Suggests** testthat, knitr, rmarkdown, BiocStyle

**LinkingTo** Rcpp, BH

**VignetteBuilder** knitr

**LazyLoad** true

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**biocViews** GeneExpression, Transcription, GeneSetEnrichment,  
DifferentialExpression, Bayesian, Clustering, TimeCourse,  
RNASeq, Microarray, MultipleComparison, DimensionReduction,  
ImmunoOncology

**NeedsCompilation** yes

**RoxygenNote** 7.2.3

**Encoding** UTF-8

**Collate** 'class-CogapsParams.R' 'CoGAPS.R' 'DistributedCogaps.R'  
 'HelperFunctions.R' 'Package.R' 'RcppExports.R' 'SubsetData.R'  
 'class-CogapsResult.R' 'methods-CogapsParams.R'  
 'methods-CogapsResult.R'

**git\_url** <https://git.bioconductor.org/packages/CoGAPS>

**git\_branch** RELEASE\_3\_17

**git\_last\_commit** 6861e5b

**git\_last\_commit\_date** 2023-07-11

**Date/Publication** 2023-10-12

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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.

Package: CoGAPS  
 Type: Package  
 Version: 2.99.0  
 Date: 2018-01-24  
 License: LGPL

## Author(s)

Maintainer: Elana J. Fertig <ejfertig@jhmi.edu>, Michael F. Ochs <ochsm@tcnj.edu>

## 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	<i>Information About Package Compilation</i>
-------------	--

---

**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	<i>calculate statistic on sets of measurements (genes) or samples</i>
----------------	---

---

**Description**

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

**Usage**

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

*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*

---

## 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

<code>data</code>	data in a supported format
<code>allParams</code>	list of all parameters
<code>uncertainty</code>	uncertainty of data in the same format
<code>index</code>	index for which subset to run on
<code>sets</code>	list of all subsets
<code>geneNames</code>	names of all genes
<code>sampleNames</code>	names of all samples
<code>fixedMatrix</code>	matrix of matched patterns

## Value

CogapsResult object

---

`checkDataMatrix`

*check that provided data is valid*

---

## 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"),
  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>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

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, nIterations=25)

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

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

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*      *CogapsParams*

**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

---

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()`

---

<code>computeGeneGSProb</code>	<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

<code>object</code>	an object of type CogapsResult
<code>GStoGenes</code>	data.frame or list with gene sets
<code>numPerm</code>	number of permutations for null
<code>Pw</code>	weight on genes
<code>PwNull</code>	- 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     *convert any acceptable data input to a numeric matrix*

---

**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            *cluster patterns together*

---

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

---

**Description**

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

**Usage**

```
corrToMeanPattern(cluster)
```

**Value**

correlation of each pattern

---

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

---

**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`*partition genes/samples into subsets*

---

**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**

- |                        |                               |
|------------------------|-------------------------------|
| <code>data</code>      | either file name or matrix    |
| <code>allParams</code> | list of all CoGAPS parameters |

**Value**

list of sorted subsets of either genes or samples

---

`distributedCogaps`*CoGAPS Distributed Matrix Factorization Algorithm*

---

**Description**

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

**Usage**`distributedCogaps(data, allParams, uncertainty)`**Arguments**

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

**Details**

For file types CoGAPS supports csv, tsv, and mtx

**Value**`list`

---

findConsensusMatrix    *find the consensus pattern matrix across all subsets*

---

**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

---

gapsCat                  *wrapper around cat*

---

**Description**

cleans up message printing

**Usage**

`gapsCat(allParams, ...)`

**Arguments**

`allParams`    all cogaps parameters  
`...`    arguments forwarded to cat

**Value**

conditionally print message

---

```
getAmplitudeMatrix      return Amplitude matrix from CogapsResult object
```

---

**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")
getParam(params, "seed")
```

<code>getPatternMatrix</code>	<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**

<code>object</code>	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

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")
params <- setDistributedParams(params, nSets=2)
params <- setParam(params, "nIterations", 100)
params <- setParam(params, "nPatters", 3)
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, ...)
```

**Arguments**

.Object	CogapsParams object
distributed	either "genome-wide" or "single-cell" indicating which distributed algorithm should be used
...	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

<code>isRdsFile</code>	<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

---

ncolHelper

*get number of columns from supported file name or matrix*

---

### 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

*get number of rows from supported file name or matrix*

---

### 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`      *parse parameters passed through the ... variable*

### Description

parse parameters passed through the ... variable

### Usage

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

### Arguments

<code>allParams</code>	list of all parameters
<code>extraParams</code>	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`      *compute pattern markers statistic*

### Description

calculate the most associated pattern for each gene

### Usage

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

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

### Arguments

<code>object</code>	an object of type <code>CogapsResult</code>
<code>threshold</code>	the type of threshold to be used. The default "all" will distribute genes into pattern with the lowest ranking. The "cut" thresholds by the first gene to have a lower ranking, i.e. better fit to, a pattern.
<code>lp</code>	a vector of weights for each pattern to be used for finding markers. If NA markers for each pattern of the A matrix will be used.
<code>axis</code>	either 1 or 2, specifying if pattern markers should be calculated using the rows of the data (1) or the columns of the data (2)

**Value**

By default a non-overlapping list of genes associated with each 1p.

**Examples**

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

---

patternMatch

*Match Patterns Across Multiple Runs*

---

**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

---

plotPatternMarkers

*heatmap of original data clustered by pattern markers statistic*

---

**Description**

heatmap of original data clustered by pattern markers statistic

**Usage**

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

**Arguments**

<code>object</code>	an object of type CogapsResult
<code>data</code>	the original data as a matrix
<code>patternPalette</code>	a vector indicating what color should be used for each pattern
<code>sampleNames</code>	names of the samples to use for labeling
<code>samplePalette</code>	a vector indicating what color should be used for each sample
<code>heatmapCol</code>	pallelet giving color scheme for heatmap
<code>colDenogram</code>	logical indicating whether to display sample denogram
<code>scale</code>	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".
<code>...</code>	additional graphical parameters to be passed to <code>heatmap.2</code>

**Value**

heatmap of the data values for the `patternMarkers`

**See Also**

[heatmap.2](#)

`plotResiduals`

*plot of residuals*

**Description**

calculate residuals and produce heatmap

**Usage**

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

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

**Arguments**

<code>object</code>	an object of type CogapsResult
<code>data</code>	original data matrix run through GAPS
<code>uncertainty</code>	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

*reconstruct gene*

---

**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

*subset data by uniformly partitioning rows (cols)*

---

**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`

*subset rows (cols) proportional to the user provided weights*

---

**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` *use user provided subsets*

---

**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")
params <- setDistributedParams(params, nSets=2)
params <- setParam(params, "nIterations", 100)
params <- setParam(params, "nPatters", 3)
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")
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**

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

**Value**

the modified params object

## Examples

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

**setFixedPatterns**      *set the fixed patterns for either the A or the P matrix*

## 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")
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")
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

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