# 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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# Contents

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`CoGAPS-package` *CoGAPS: Coordinated Gene Activity in Pattern Sets*

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## 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.

Package: CoGAPS

| | |
|---|---|
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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

```
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 calculatin p-value |
| ... | handles old arguments for backwards compatibility |

### Value

gene set statistics for each column of A

---

## calcGeneGSStat *probability gene belongs in gene set*

---

### 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` *compute z-score matrix*

## 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

| | |
|---|---|
| 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 *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 *check that all inputs are valid*

---

### 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` *Check if package was built with checkpoints enabled*

---

### Description

Check if package was built with checkpoints enabled

### Usage

```
checkpointsEnabled()
```

### Value

true/false if checkpoints are enabled

### Examples

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

---

`CoGAPS` *CoGAPS Matrix Factorization Algorithm*

---

### 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

| | |
|---|---|
| data | File name or R object (see details for supported types) |
| params | CogapsParams object |
| nPatterns | rank of the nmf decomposition |
| 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 neccesary 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 snapsjots 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` *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 matching

`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 *compute gene probability*

### 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 containined 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

| | |
|---|---|
| data | either file name or matrix |
| allParams | 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

| | |
|---|---|
| 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 *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

## fromCSV *read CoGAPS Result object from a directory with a set of csvs see toCSV*

### 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 *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` *extracts gene/sample names from the data*

---

## 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` *return featureLoadings matrix from CogapsResult object*

---

## 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 *extract gene names from data*

### Description

extract gene names from data

### Usage

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

### Value

vector of gene names

## getMeanChiSq *return chi-sq of final matrices*

### 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` *generate statistics associating patterns with gene sets*

---

## 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 *return pattern matrix from CogapsResult object*

### 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 *get specified number of retina subsets*

### 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 *return sampleFactors matrix from CogapsResult object*

### 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 *extract sample names from data*

### Description

extract sample names from data

### Usage

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

### Value

vector of sample names

## getSubsets *return the names of the genes (samples) in each subset*

### 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 *return version number used to generate this result*

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

### Description

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

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

### 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 neccesary 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 <- 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 *checks if file is rds format*

### 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 *Result of applying CoGAPS on the Toy example.*

### Description

Result of applying CoGAPS on the Toy example.

### Usage

```
data(modsimresult)
```

### Format

S4 class 'CogapsResult' [package "CoGAPS"] with 7 slots.

## 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

| | |
|---|---|
| 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 *compute pattern markers statistic*

## 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 *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

## plotPatternGeneSet *generate a barchart of most significant hallmark sets for a pattern*

### 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 *heatmap of original data clustered by pattern markers statistic*

## 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 | pallelet 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 *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

| | |
|---|---|
| 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 *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 partioning rows (cols)*

---

### Description

subset data by uniformly partioning rows (cols)

### Usage

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

### Arguments

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

## sampleWithExplictSets *use user provided subsets*

### Description

use user provided subsets

### Usage

```
sampleWithExplictSets(allParams)
```

### Arguments

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

### 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 neccesary 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 *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", nPatterns=3)
data(GIST)
params <- setFixedPatterns(params, getSampleFactors(GIST.result), 'P')
```

## setParam *set the value of a parameter*

### 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 *write start up message*

### Description

write start up message

### Usage

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

### Arguments

| | |
|---|---|
| data | data set |
| allParams | list of all parameters |

### Value

message displayed to screen

## stitchTogether *concatenate final results across subsets*

### 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 *checks if file is supported*

### Description

checks if file is supported

### Usage

```
supported(file)
```

### Arguments

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

### Value

TRUE if file is supported, FALSE if not

toCSV *save CoGAPS Result object as a set of csvs to directory see fromCSV*

## 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

# Index