

# Package: rOCEAN (via r-universe)

November 13, 2024

**Type** Package

**Title** Two-Way Feature Set Testing for Multi-Omics

**Version** 1.0

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**Description** For any two way feature-set from a pair of pre-processed omics data, pairwise-TDP, Column-TDP and row-TDP are calculated. Due to embeded closed testing procedure, the choice of feature-sets can be changed infinite times and even after seeing the data withuot any change in type I error rate. For more details refer to the refrence article.

**License** GPL (>= 2)

**Date** 2024-10-29

**Encoding** UTF-8

**RoxygenNote** 7.3.2

**Imports** ff

**Repository** <https://mitra-ep.r-universe.dev>

**RemoteUrl** <https://github.com/mitra-ep/rocean>

**RemoteRef** HEAD

**RemoteSha** 4c44201690d1773e0a3f10a738fca4ba4241e933

## Contents

corPs . . . . .	2
getCat . . . . .	2
ocean . . . . .	3
pairTDP . . . . .	4
runbab . . . . .	5
simesCT . . . . .	5
singleStep . . . . .	6

<b>Index</b>	<b>7</b>
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corPs *Calculate pairwise p-value*

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### Description

Calculates pairwise matrix of p-values based on Pearson's correlation test for two matrices. To gain speed and manage RAM usage, the matrices are split into several smaller chunks.

### Usage

```
corPs(pm1, pm2, type = c("Mat", "Vec"), pthresh = 0.05)
```

### Arguments

pm1, pm2	Subsets of two omics data sets where rows are the features and columns are samples. The rows of the two matrices would define the two-way feature set of interest.
type	Two options are available. Mat: Calculate the correlation of subsets and return a matrix; Vec: calculate the correlation matrix, subset by the given threshold and return a vector of p-values.
pthresh	Only relevant for type="Vec". The threshold by which the p-values are filtered (p>pthresh is removed). Default value is 0.05.

### Value

Either a matrix or vector of pairwise p-values, as indicated by type parameter.

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getCat *Calculate p-categories*

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### Description

Calculates cumulative p-categories for a given matrix of p-values.

### Usage

```
getCat(mps, gCT, scale = c("col", "row"))
```

### Arguments

mps	Matrix of p-values, representing pairwise associations between two feature sets.
gCT	Parameters of the global closed testing, which is the output of simesCT function.
scale	Scale of the quantification, a character string. Possible choices are "col" and "row".

**Value**

Matrix of p-categories.

**See Also**

[simesCT](#)

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ocean

*OCEAN algorithm*

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**Description**

Calculates heuristic and lower bound for the true discovery proportion (TDP) in 3 scales for a specified two-way feature set (Algorithm 1 in the reference). The input is either two omics data sub-matrices or the pre-calculated matrix of p-values for pairwise associations. In case the result is not exact, the function adopts branch and bound (Algorithm 2 in the reference), if nMax allows.

**Usage**

```
ocean(
  pm1,
  pm2,
  gCT,
  scale = c("pair", "row", "col"),
  mps,
  nMax = 100,
  verbose = TRUE
)
```

**Arguments**

pm1, pm2	Matrix; Subsets of two omics data sets where rows are the features and columns are samples. The rows of the two matrices would define the two-way feature set of interest.
gCT	Vector; Parameters of the global closed testing, output of simesCT function.
scale	Optional character vector; It specifies the scale of TDP quantification. Possible choices are "pair" (pair-TDP), "col" (col-TDP ) and "row" (for row-TDP'). If not specified, all three scales are returned.
mps	Optional matrix of p-values; A sub-matrix of pairwise associations, representing the two-way feature set of interest. If provided, pm1 and pm2 are not required. If not provided, matrix of pairwise associations will be derived from pm1 and pm2 based on Pearson's correlation.
nMax	Maximum number of steps for branch and bound algorithm, if set to 1 branch and bound is skipped even if the result is not exact. The default value is a 100. The algorithm may stop before the nMax is reached if it converges sooner.
verbose	Logical; if TRUE, progress messages will be displayed during the function's execution. Default is TRUE.

**Value**

TDP is returned for the specified scales, along with number of steps taken and convergence status for branch and bound algorithm.

**See Also**

[simesCT](#) [pairTDP](#) [runbab](#)

**Examples**

```
#number of feature per omic data set
n_cols<-1000
n_rows<-1200

#random matrix of p-values
set.seed(1258)
pvalmat<-matrix(runif(n_rows*n_cols, min=0, max=1)^3, nrow=n_rows, ncol=n_cols)

#calculate CT parameters
gCT<-simesCT(mps=pvalmat, m=nrow(pvalmat)*ncol(pvalmat))

#calculate TDPs for a random feature set
subpmat<-pvalmat[1:400,100:750]
#Note: it can take loner to run this script if nMax is large
out<-ocean(mps=subpmat, gCT=gCT, nMax=2)
out
```

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pairTDP	<i>pairwise true discoveries proportion</i>
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**Description**

Calculates the TDP over pairs; based on SEA algorithm

**Usage**

```
pairTDP(mps, n, gCT)
```

**Arguments**

mps	Matrix or vector of pairwise associations.
n	Number of pairs; may not be the size of p if a threshold is used to remove large p-values.
gCT	Parameters of the global closed testing, output of <code>simesCT</code> function.

**Value**

Proportion of true discoveries out of n pairs of features.

**See Also**

[SEA](#), [simesCT](#)

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runbab	<i>Branch and bound algorithm implementation</i>
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**Description**

Performs B&B when the bound are not exact

**Usage**

```
runbab(sCat, ssh, ssb, nMax = 100)
```

**Arguments**

sCat	Category matrix, output of getCat function
ssh	current Heuristic as provided by SingleStep function
ssb	current Bound as provided by SingleStep function
nMax	Maximum number of steps for the algorithm, the algorithm may stop sooner if it converges.

**Value**

A list, including the heuristic and the bound for the number of true discoveries, along with number of steps taken and convergence status.

**See Also**

[getCat](#) [singleStep](#)

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simesCT	<i>Closed testing with Simes</i>
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**Description**

Calculates five parameters from closed testing with Simes local tests based on raw data. These parameter are unique per data/alpha-level combination and do not depend on feature sets. Calculation may be somewhat long depending on the size of data sets and PC configurations.

**Usage**

```
simesCT(om1, om2, mps, m, alpha = 0.05)
```

**Arguments**

om1, om2	Two omics data sets where rows are features and columns are samples.
mps, m	Optional, pre-calculated matrix/vector of pairwise associations and the size. To save time in calculation of parameters, a threshold such as the type I error may be applied to remove larger p-values. If a threshold is used, size of matrix and m will not match. m should always be the size of the matrix of associations (number of features in om1 X number of features in om2).
alpha	type I error rate, default value is 0.05.

**Value**

Vector of integers: grand H value, concentration p-value, size of concentration set z, size of the original pair-wise associations matrix and the type I error level used in calculations.

**References**

See more details in "Hommel's procedure in linear time" doi: 10.1002/bimj.201700316.

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singleStep	<i>Single step algorithm</i>
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**Description**

Calculates heuristic and upper-bound for the number of true discoveries based on the Algorithm 1 introduced in paper.

**Usage**

```
singleStep(sCat, B)
```

**Arguments**

sCat	p-categories matrix, output of getCat function.
B	Optional, to identify rows to be fixed (1) or removed (0) while splitting the search space.

**Value**

A list of two objects, the heuristic and the lower bound for true number of discoveries

**See Also**

[getCat](#)

# Index

corPs, [2](#)

getCat, [2](#), [5](#), [6](#)

ocean, [3](#)

pairTDP, [4](#), [4](#)

runbab, [4](#), [5](#)

SEA, [5](#)

simesCT, [3–5](#), [5](#)

singleStep, [5](#), [6](#)