

# Package: MultiFit (via r-universe)

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**Type** Package

**Title** Multiscale Fisher's Independence Test for Multivariate  
Dependence

**Version** 1.1.1

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**Description** Test for independence of two random vectors, learn and  
report the dependency structure. For more information, see  
Gorsky, Shai and Li Ma, Multiscale Fisher's Independence Test  
for Multivariate Dependence, Biometrika, accepted, January  
2022.

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**Imports** Rcpp (>= 1.0.8), data.table

**LinkingTo** Rcpp, RcppArmadillo

**Suggests** png, qgraph, knitr, rmarkdown

**RoxygenNote** 6.1.1

**VignetteBuilder** knitr

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

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**Description**

Perform multiscale test of independence for multivariate vectors. See vignettes for further examples.

**Usage**

```
MultiFIT(xy, x = NULL, y = NULL, p_star = NULL, R_max = NULL,
  R_star = 1, rank.transform = TRUE, ranking.approximation = FALSE,
  M = 10, apply.stopping.rule = FALSE, alpha = 0.05,
  test.method = "Fisher", correct = TRUE, min.tbl.tot = 25L,
  min.row.tot = 10L, min.col.tot = 10L, p.adjust.methods = c("H",
  "Hcorrected"), compute.all.holm = TRUE, return.all.pvs = TRUE,
  verbose = FALSE)
```

**Arguments**

<code>xy</code>	A list, whose first element corresponds to the matrix <code>x</code> as below, and its second element corresponds to the matrix <code>y</code> as below. If <code>xy</code> is not specified, <code>x</code> and <code>y</code> need to be assigned.
<code>x</code>	A matrix, number of columns = dimension of random vector, number of rows = number of observations.
<code>y</code>	A matrix, number of columns = dimension of random vector, number of rows = number of observations.
<code>p_star</code>	Numeric, cuboids associated with tests whose p-value is below <code>p_star</code> will be halved and further tested.
<code>R_max</code>	A positive integer (or Inf), the maximal number of resolutions to scan (algorithm will stop at a lower resolution if all tables in it do not meet the criteria specified at <code>min.tbl.tot</code> , <code>min.row.tot</code> and <code>min.col.tot</code> )
<code>R_star</code>	A positive integer, if set to an integer between 0 and <code>R_max</code> , all tests up to and including resolution <code>R_star</code> will be performed (algorithm will stop at a lower resolution than requested if all tables in it do not meet the criteria specified at <code>min.tbl.tot</code> , <code>min.row.tot</code> and <code>min.col.tot</code> ). For higher resolutions only the children of tests with p-value lower than <code>p_star</code> will be considered.
<code>rank.transform</code>	Logical, if TRUE, marginal rank transform is performed on all margins of <code>x</code> and <code>y</code> . If FALSE, all margins are scaled to 0-1 scale. When FALSE, the average and top statistics of the negative logarithm of the p-values are only computed for the univariate case.
<code>ranking.approximation</code>	Logical, if FALSE, select only tests with p-values more extreme than <code>p_star</code> to halve and further test. FWER control not guaranteed. If TRUE, choose at each resolution the <code>M</code> tests with the most extreme p-values to further halve and test.

<code>M</code>	A positive integer (or Inf), the number of top ranking tests to continue to split at each resolution. FWER control not guaranteed for this method.
<code>apply.stopping.rule</code>	Logical. If TRUE, an adjusted p-value is computed for each resolution,
<code>alpha</code>	Numeric. Threshold below which resolution-specific p-values trigger early stopping.
<code>test.method</code>	String, choose "Fisher" for Fisher's exact test (slowest), "chi.sq" for Chi-squared test, "LR" for likelihood-ratio test and "norm.approx" for approximating the hypergeometric distribution with a normal distribution (fastest).
<code>correct</code>	Logical, if TRUE compute mid-p corrected p-values for Fisher's exact test, or Yates corrected p-values for the Chi-squared test, or Williams corrected p-values for the likelihood-ratio test.
<code>min.tbl.tot</code>	Non-negative integer, the minimal number of observations per table below which a p-value for a given table will not be computed.
<code>min.row.tot</code>	Non-negative integer, the minimal number of observations for row totals in the 2x2 contingency tables below which a contingency table will not be tested.
<code>min.col.tot</code>	Non-negative integer, the minimal number of observations for column totals in the 2x2 contingency tables below which a contingency table will not be tested.
<code>p.adjust.methods</code>	String, choose between "H" for Holm, "Hcorrected" for Holm with the correction as specified in correct.
<code>compute.all.holm</code>	Logical, if FALSE, only global p-value is computed (may be a little faster when any tests are performed). If TRUE adjusted p-values are computed for all tests.
<code>return.all.pvs</code>	Logical, if TRUE, a data frame with all p-values is returned (not applicable when stopping rule is applied)
<code>verbose</code>	Logical.

## Value

`p.values.holistic`, a named numerical vector containing the holistic p-values of for the global null hypothesis (i.e. x independent of y).

`p.values.resolution.specific`, a named numerical vector containing the resolution specific p-values of for the global null hypothesis (i.e. x independent of y).

`res.by.res.pvs`, a data frame that contains the raw and Bonferroni adjusted resolution specific p-values.

`all.pvs`, a data frame that contains all p-values and adjusted p-values that are computed. Returned if `return.all.pvs` is TRUE.

`all`, a nested list. Each entry is named and contains data about a resolution that was tested. Each resolution is a list in itself, with `cuboids`, a summary of all tested cuboids in a resolution, `tables`, a summary of all 2x2 contingency tables in a resolution, `pv`, a numerical vector containing the p-values from the tests of independence on 2x2 contingency table in tables that meet the criteria defined by `min.tbl.tot`, `min.row.tot` and `min.col.tot`. The length of `pv` is equal to the number of rows of `tables`. `pv.correct`, similar to the above `pv`, corrected p-values are computed

and returned when correct is TRUE. `rank.tests`, logical vector that indicates whether or not a test was ranked among the top M tests in a resolution. The length of `rank.tests` is equal to the number of rows of tables. `parent.cuboids`, an integer vector, indicating which cuboids in a resolution are associated with the ranked tests, and will be further halved in the next higher resolution. `parent.tests`, a logical vector of the same length as the number of rows of tables, indicating whether or not a test was chosen as a parent test (same tests may have multiple children).

## Examples

```
set.seed(1)
n = 300
Dx = Dy = 2
x = matrix(0, nrow = n, ncol = Dx)
y = matrix(0, nrow = n, ncol = Dy)
x[,1] = rnorm(n)
x[,2] = runif(n)
y[,1] = rnorm(n)
y[,2] = sin(5 * pi * x[, 2]) + 1 / 5 * rnorm(n)
fit = MultiFIT(x = x, y = y, verbose = TRUE)
w = MultiSummary(x = x, y = y, fit = fit, alpha = 0.0001)
```

---

MultiSummary

*Summary of significant tests*

---

## Description

Provide a post-hoc summary of significant tests. See vignettes for further examples.

## Usage

```
MultiSummary(xy, x = NULL, y = NULL, fit, alpha = 0.05,
  only.rk = NULL, use.pval = NULL, plot.tests = TRUE, pch = NULL,
  rd = 2, plot.margin = FALSE)
```

## Arguments

<code>xy</code>	A list, whose first element corresponds to the matrix <code>x</code> as below, and its second element corresponds to the matrix <code>y</code> as below. if <code>xy</code> is not specified, <code>x</code> and <code>y</code> need to be assigned.
<code>x</code>	A matrix, number of columns = dimension of random vector, number of rows = number of observations.
<code>y</code>	A matrix, number of columns = dimension of random vector, number of rows = number of observations.
<code>fit</code>	An object generated by <code>MultiFIT</code> .
<code>alpha</code>	Numeric, only tests with adjusted p-values less than <code>alpha</code> are presented in the output.

<code>only.rk</code>	Positive integer vector. Show only tests that are ranked according to <code>only.rk</code> and have adjusted p-value below <code>alpha</code> . If left as <code>NULL</code> , all tests with adjusted p-values less than <code>alpha</code> are presented in the output.
<code>use.pval</code>	String, choose between "H" (for Holm), "Hcorrected" (for Holm on corrected p-values) or "MH" for modified Holm. If left <code>NULL</code> , the order of preference is "MH", "Hcorrected" and then "H", according to which is present in the object <code>fit</code> .
<code>plot.tests</code>	Logical, plot the marginal scatter plots that are associated with the presented significant tests.
<code>pch</code>	Point style for plots. If left as <code>NULL</code> , a default combination of crosses and bullets is applied.
<code>rd</code>	Numeric, number of figures to round to when presenting ranges of variables.
<code>plot.margin</code>	Logical, plot the marginal scatter plot of the margins that are associated with each significant test, without highlighting which points are conditioned on and are in the discretized 2x2 contingency table.

### Value

List whose elements are `significant.tests`, a data frame that summarizes the main features of the tests and their overall ranking by p-value and `original.scale.cuboids`, a list whose number of elements is equal to the number of significant tests (the same number of rows of the data frame `significant.tests`). Each element corresponds to a test and is a list whose elements are the marginal ranges of the associated cuboid.

### Examples

```
set.seed(1)
n = 300
Dx = Dy = 2
x = matrix(0, nrow = n, ncol = Dx)
y = matrix(0, nrow = n, ncol = Dy)
x[,1] = rnorm(n)
x[,2] = runif(n)
y[,1] = rnorm(n)
y[,2] = sin(5 * pi * x[, 2]) + 1 / 5 * rnorm(n)
fit = MultiFIT(x = x, y = y, verbose = TRUE)
w = MultiSummary(x = x, y = y, fit = fit, alpha = 0.0001)
```

---

MultiTree

---

*Plot tree structure of tests on 2x2 contingency tables*


---

### Description

Plot a post-hoc tree of all tests or all significant tests on 2x2 discretized contingency tables. See vignettes for examples.

**Usage**

```
MultiTree(xy, x = NULL, y = NULL, fit, show.all = FALSE,
          max.node.size = 5, min.node.size = 2.5, use.pval = NULL,
          images.path = NULL, node.name = "node", filename = NULL,
          filetype = "pdf")
```

**Arguments**

<code>xy</code>	A list (optional), whose first element corresponds to the matrix <code>x</code> as below, and its second element corresponds to the matrix <code>y</code> as below. if <code>xy</code> is not specified, <code>x</code> and <code>y</code> need to be assigned. If <code>xy</code> , <code>x</code> and <code>y</code> are missing or <code>NULL</code> , the tree nodes are blank. If <code>xy</code> or <code>x</code> and <code>y</code> are provided, nodes are png images of the marginal scatter plots that are associated with each test.
<code>x</code>	A matrix (optional), number of columns = dimension of random vector, number of rows = number of observations. If <code>xy</code> , <code>x</code> and <code>y</code> are missing or <code>NULL</code> , the tree nodes are blank. If <code>xy</code> or <code>x</code> and <code>y</code> are provided, nodes are png images of the marginal scatter plots that are associated with each test.
<code>y</code>	A matrix (optional), number of columns = dimension of random vector, number of rows = number of observations. If <code>xy</code> , <code>x</code> and <code>y</code> are missing or <code>NULL</code> , the tree nodes are blank. If <code>xy</code> or <code>x</code> and <code>y</code> are provided, nodes are png images of the marginal scatter plots that are associated with each test.
<code>fit</code>	An object generated by <code>multiFit</code> .
<code>show.all</code>	Logical. If <code>TRUE</code> , all tests are shown. If <code>FALSE</code> only tests who were ranked in each resolution amongst the top <code>M</code> ranking tests are shown. See <code>?multiFit</code> for an explanation about the parameter <code>M</code> and see documentation for further information.
<code>max.node.size</code>	Numeric. Maximal node size. All nodes are scaled between <code>min.node.size</code> and <code>max.node.size</code> , where larger nodes are associated smaller p-values of the corresponding tests on 2x2 contingency tables.
<code>min.node.size</code>	Numeric. Minimal node size. All nodes are scaled between <code>min.node.size</code> and <code>max.node.size</code> , where larger nodes are associated smaller p-values of the corresponding tests on 2x2 contingency tables.
<code>use.pval</code>	String, choose between "H" (for Holm), "Hcorrected" (for Holm on corrected p-values) or "MH" for modified Holm. If left <code>NULL</code> , the order of preference is "MH", "Hcorrected" and then "H", according to which is present in the object <code>fit</code> .
<code>images.path</code>	String, path to save png images of nodes to. If not specified, images are saved to <code>tempdir()</code> .
<code>node.name</code>	String, prefix for file names for nodes pngs.
<code>filename</code>	String, file name for tree output. If left <code>NULL</code> , file name is prefixed by <code>multiTree</code> and ends with system time. See documentation of <code>qgraph::qgraph</code> for further information.
<code>filetype</code>	String, default is <code>pdf</code> , See documentation of <code>qgraph::qgraph</code> for further information.

**Value**

The main output of `multiTree` is a pdf file with the directed acyclic graph showing tests as nodes.

In addition, the function returns a list. Its elements are: `qgraph.object`, the graphical object generated by the `qgraph` function. See the `qgraph` package documentation for further details. `qgraph.call`, the call for the tree generating function. Arguments for the call: `adj`, the adjacency matrix, `nodes.size`, a numeric vector with the scaled sizes of the nodes, `images`, the file names of the nodes images (may be `NULL`), `filename` as passed to `multiTree` and passed over to `qgraph`, and `filetype` as passed to `multiTree` and passed over to `qgraph`.

Other elements of the returned list are `pvs.attributes`, the attributes summarizing the data and the tests performed as stored in `fit`, and `n.nodes`, the number of nodes.

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