

# Package ‘ReactomeGraph4R’

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**Title** Interface for the Reactome Graph Database

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**Description** Pathways, reactions, and biological entities in Reactome knowledge are systematically represented as an ordered network. Instances are represented as nodes and relationships between instances as edges; they are all stored in the Reactome Graph Database. This package serves as an interface to query the interconnected data from a local Neo4j database, with the aim of minimizing the usage of Neo4j Cypher queries.

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## Contents

|                                            |           |
|--------------------------------------------|-----------|
| ReactomeGraph4R-package . . . . .          | 2         |
| login . . . . .                            | 3         |
| matchDiseases . . . . .                    | 3         |
| matchHierarchy . . . . .                   | 4         |
| matchInteractors . . . . .                 | 5         |
| matchObject . . . . .                      | 6         |
| matchPaperObjects . . . . .                | 7         |
| matchPERoles . . . . .                     | 8         |
| matchPrecedingAndFollowingEvents . . . . . | 9         |
| matchReactionsInPathway . . . . .          | 10        |
| matchReferrals . . . . .                   | 11        |
| multiObjects . . . . .                     | 12        |
| unnestListCol . . . . .                    | 13        |
| <b>Index</b>                               | <b>14</b> |

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ReactomeGraph4R-package

*ReactomeGraph4R: Interface for the Reactome Graph Database*

---

## Description

Pathways, reactions, and biological entities in Reactome knowledge are systematically represented as an ordered network. Instances are represented as nodes and relationships between instances as edges; they are all stored in the Reactome Graph Database. This package serves as an interface to query the interconnected data from a local Neo4j database, with the aim of minimizing the usage of Neo4j Cypher queries.

## Author(s)

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## See Also

Useful links:

- <https://github.com/reactome/ReactomeGraph4R>
- Report bugs at <https://github.com/reactome/ReactomeGraph4R/issues>

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|       |                                         |
|-------|-----------------------------------------|
| login | <i>Log in to the local neo4j server</i> |
|-------|-----------------------------------------|

---

**Description**

Before running `login()`, you have to successfully finish the Reactome Neo4j database setup and build a connection on your local machine (details see: <https://github.com/reactome/ReactomeGraph4R>). This command is to create a `neo4r` object that is used to communicate between R and Neo4j, also to do a sanity check for the connection.

**Usage**

```
login(con = NULL)
```

**Arguments**

`con` an existed connexion object. It is not necessary to log in for the first time.

**Value**

connection to the local neo4j database

**Examples**

```
## Not run:  
# The first step to the graph database!  
login()  
  
## End(Not run)  
# you can also check the neo4r connexion object by running:  
getOption("con")
```

---

|               |                                                          |
|---------------|----------------------------------------------------------|
| matchDiseases | <i>MATCH diseases of PhysicalEntity/Reaction/Pathway</i> |
|---------------|----------------------------------------------------------|

---

**Description**

To find Diseases related to a PhysicalEntity or an Event, or get PhysicalEntities/Events associated with a Disease in reverse

**Usage**

```
matchDiseases(  
  id = NULL,  
  displayName = NULL,  
  species = NULL,  
  type = c("row", "graph")  
)
```

**Arguments**

|             |                                                                                                  |
|-------------|--------------------------------------------------------------------------------------------------|
| id          | stId or dbId of a PhysicalEntity/Event/Disease                                                   |
| displayName | displayName of a PhysicalEntity/Event/Disease                                                    |
| species     | name or taxon id or dbId or abbreviation of aspecies                                             |
| type        | return results as a list of dataframes ( <b>'row'</b> ), or as a graph object ( <b>'graph'</b> ) |

**Value**

Disease(s) related to the given PhysicalEntity/Reaction/Pathway; or instances related to the given Disease

**See Also**

Other match: [matchHierarchy\(\)](#), [matchInteractors\(\)](#), [matchObject\(\)](#), [matchPEroles\(\)](#), [matchPaperObjects\(\)](#), [matchPrecedingAndFollowingEvents\(\)](#), [matchReactionsInPathway\(\)](#), [matchReferrals\(\)](#)

**Examples**

```
disease <- "neuropathy"
# matchDiseases(displayName=disease, species="M. musculus", type="row")
# matchDiseases(id="R-HSA-162588", type="graph")
```

---

|                |                        |
|----------------|------------------------|
| matchHierarchy | <i>MATCH hierarchy</i> |
|----------------|------------------------|

---

**Description**

Reactome data are organized in a hierarchical way: Pathway-Reaction-Entity. This function retrieves the hierarchical data of a given Event (Pathway or Reaction) or Entity (PhysicalEntity or ReferenceEntity).

**Usage**

```
matchHierarchy(
  id = NULL,
  displayName = NULL,
  databaseName = "Reactome",
  species = NULL,
  type = c("row", "graph")
)
```

**Arguments**

|              |                                                                                                  |
|--------------|--------------------------------------------------------------------------------------------------|
| id           | stId or dbId of an Event/Entity; or an external id                                               |
| displayName  | displayName of Event/PhysicalEntity/ReferenceEntity                                              |
| databaseName | database name                                                                                    |
| species      | name or taxon id or dbId or abbreviation of specified species                                    |
| type         | return results as a list of dataframes ( <b>'row'</b> ), or as a graph object ( <b>'graph'</b> ) |

**Value**

hierarchical instances of the given id and databaseName

**See Also**

Other match: [matchDiseases\(\)](#), [matchInteractors\(\)](#), [matchObject\(\)](#), [matchPEroles\(\)](#), [matchPaperObjects\(\)](#), [matchPrecedingAndFollowingEvents\(\)](#), [matchReactionsInPathway\(\)](#), [matchReferrals\(\)](#)

**Examples**

```
## use the Reactome displayName of a UniProt object
uniprot.name <- "UniProt:P04637 TP53"
# matchHierarchy(displayName=uniprot.name,
#               databaseName="UniProt", type="row")
# matchHierarchy(id="R-HSA-1369062", type="graph")
```

---

|                  |                          |
|------------------|--------------------------|
| matchInteractors | <i>MATCH interactors</i> |
|------------------|--------------------------|

---

**Description**

To retrieve interactions of a given PhysicalEntity (PE), it first finds the ReferenceEntity matched with the PE, then get the Interactions having "interactor" relationship with the ReferenceEntity.

**Usage**

```
matchInteractors(
  pe.id = NULL,
  pe.displayName = NULL,
  species = NULL,
  type = c("row", "graph")
)
```

**Arguments**

|                |                                                                                                  |
|----------------|--------------------------------------------------------------------------------------------------|
| pe.id          | stId or dbId of a PhysicalEntity                                                                 |
| pe.displayName | displayName of a PhysicalEntity                                                                  |
| species        | name or taxon id or dbId or abbreviation of specified species                                    |
| type           | return results as a list of dataframes ( <b>'row'</b> ), or as a graph object ( <b>'graph'</b> ) |

**Value**

interactions of a given PhysicalEntity

**See Also**

Other match: [matchDiseases\(\)](#), [matchHierarchy\(\)](#), [matchObject\(\)](#), [matchPEroles\(\)](#), [matchPaperObjects\(\)](#), [matchPrecedingAndFollowingEvents\(\)](#), [matchReactionsInPathway\(\)](#), [matchReferrals\(\)](#)

## Examples

```
pe.id <- 996766
# matchInteractors(pe.id)
```

---

matchObject

*Basic query for database objects*


---

## Description

This function can fetch instance by setting the following arguments:

- **id**: a Reactome dbId/stId, or non-Reactome id (e.g. UniProt)
- **displayName**: a display name of a Reactome object
- **schemaClass**: a specific schema class, see [Data Schema](#)
- **property**: a property of a node or relationship, access the full list of properties: `con <- getOption("con"); con$get_`
- **relationship**: a relationship between nodes, access the full list of relationships: `con <- getOption("con"); con$get_`
- Species information can see [here](#), or run `View(matchObject(schemaClass = "Species")[['databaseObject']])` to view a full table

## Usage

```
matchObject(
  id = NULL,
  displayName = NULL,
  schemaClass = NULL,
  species = NULL,
  returnedAttributes = NULL,
  property = NULL,
  relationship = NULL,
  limit = NULL,
  databaseName = "Reactome"
)
```

## Arguments

|                    |                                                                                                   |
|--------------------|---------------------------------------------------------------------------------------------------|
| id                 | Reactome stId or dbId, or non-Reactome identifier                                                 |
| displayName        | displayName of a database object                                                                  |
| schemaClass        | schema class of a database object                                                                 |
| species            | name or taxon id or dbId or abbreviation of specified species                                     |
| returnedAttributes | specific attribute(s) to be returned. If set to NULL, all attributes returned                     |
| property           | a list of property keys and values, e.g. <code>list(isChimeric = TRUE, isInDisease = TRUE)</code> |
| relationship       | relationship type(s)                                                                              |
| limit              | the number of returned objects                                                                    |
| databaseName       | database name. All databases see <a href="#">here</a>                                             |

**Value**

Reactome database object(s) that meets all specified conditions

**See Also**

[multiObjects](#) for multiple ids

Other match: [matchDiseases\(\)](#), [matchHierarchy\(\)](#), [matchInteractors\(\)](#), [matchPERoles\(\)](#), [matchPaperObjects\(\)](#), [matchPrecedingAndFollowingEvents\(\)](#), [matchReactionsInPathway\(\)](#), [matchReferrals\(\)](#)

**Examples**

```
## fetch instance by class
# all.species <- matchObject(schemaClass = "Species")

## fetch instance by name
# matchObject(displayName = "RCOR1 [nucleoplasm]",
#             returnedAttributes=c("stId", "speciesName"))

## fetch instance by id
## Reactome id
# matchObject(id = "R-HSA-9626034")
## non-Reactome id
# matchObject(id = "P60484", databaseName = "UniProt")

## fetch instances by relationship
# matchObject(relationship="inferredTo", limit=10)

## fetch instances by property
property.list <- list(hasEHLD = TRUE, isInDisease = TRUE)
# matchObject(property = property.list,
#             returnedAttributes = c("displayName", "stId", "isInDisease", "hasEHLD"),
#             limit=20)
```

---

|                   |                                         |
|-------------------|-----------------------------------------|
| matchPaperObjects | <i>MATCH objects related to a paper</i> |
|-------------------|-----------------------------------------|

---

**Description**

Fetch Reactome instances related to a paper by its PubMed id or title

**Usage**

```
matchPaperObjects(
  pubmed.id = NULL,
  displayName = NULL,
  type = c("row", "graph")
)
```

**Arguments**

|             |                                                                                                  |
|-------------|--------------------------------------------------------------------------------------------------|
| pubmed.id   | PubMed identifier of a paper                                                                     |
| displayName | paper title                                                                                      |
| type        | return results as a list of dataframes ( <b>'row'</b> ), or as a graph object ( <b>'graph'</b> ) |

**Value**

Reactome instances associated with a paper

**See Also**

Other match: [matchDiseases\(\)](#), [matchHierarchy\(\)](#), [matchInteractors\(\)](#), [matchObject\(\)](#), [matchPEroles\(\)](#), [matchPrecedingAndFollowingEvents\(\)](#), [matchReactionsInPathway\(\)](#), [matchReferrals\(\)](#)

**Examples**

```
## fetch Reactome instances by paper title
paper <- "Chaperone-mediated autophagy at a glance"
# matchPaperObjects(displayName=paper)

## fetch Reactome instances by pubmed id
# matchPaperObjects(pubmed.id="20797626", type="graph")
# matchPaperObjects(pubmed.id="23515720", type="row")
```

---

matchPEroles

---

*MATCH roles of PhysicalEntity*


---

**Description**

This function retrieves the role(s) of a given PhysicalEntity including:

- Input
- Output
- Regulator
- Catalyst

**Usage**

```
matchPEroles(
  pe.id = NULL,
  pe.displayName = NULL,
  species = NULL,
  type = c("row", "graph")
)
```

**Arguments**

`pe.id`                    `stId` or `dbId` of a `PhysicalEntity`  
`pe.displayName`        `displayName` of a `PhysicalEntity`  
`species`                `name` or `taxon id` or `dbId` or `abbreviation` of a `species`  
`type`                    return results as a list of dataframes (**'row'**), or as a graph object (**'graph'**)

**Value**

information of the given `PhysicalEntity` and its role(s)

**See Also**

Other `match`: `matchDiseases()`, `matchHierarchy()`, `matchInteractors()`, `matchObject()`, `matchPaperObjects()`, `matchPrecedingAndFollowingEvents()`, `matchReactionsInPathway()`, `matchReferrals()`

**Examples**

```

stId <- "R-HSA-8944354"
# matchPEroles(pe.id = stId, type = "graph")

# matchPEroles(pe.displayName = "2SUM01:MITF [nucleoplasm]",
#               species = "pig", type = "row")

```

---

`matchPrecedingAndFollowingEvents`

*MATCH the preceding/following Events*

---

**Description**

This method can find preceding and following `ReactionLikeEvents` (RLEs) of a specific Event with the relationship `'precedingEvent'`. The argument `"depth"` is used to describe the "variable length relationships" in Neo4j, default is 1 (i.e. immediately connected); or you can set `all.depth = TRUE` to retrieve the whole context.

**Usage**

```

matchPrecedingAndFollowingEvents(
  event.id = NULL,
  event.displayName = NULL,
  species = NULL,
  depth = 1,
  all.depth = FALSE,
  type = c("row", "graph")
)

```

**Arguments**

|                                |                                                                                                     |
|--------------------------------|-----------------------------------------------------------------------------------------------------|
| <code>event.id</code>          | stId/dbId of an Event                                                                               |
| <code>event.displayName</code> | displayName of an Event                                                                             |
| <code>species</code>           | name or taxon id or dbId or abbreviation of specified species                                       |
| <code>depth</code>             | number of depths                                                                                    |
| <code>all.depth</code>         | if set to TRUE, all RLE(s) connected to the given Event in all depths returned                      |
| <code>type</code>              | to return results as a list of dataframes ( <b>'row'</b> ), or as a graph object ( <b>'graph'</b> ) |

**Value**

preceding/following Events connected to the given Event in specified depth(s), default depth = 1

**See Also**

Other match: [matchDiseases\(\)](#), [matchHierarchy\(\)](#), [matchInteractors\(\)](#), [matchObject\(\)](#), [matchPEroles\(\)](#), [matchPaperObjects\(\)](#), [matchReactionsInPathway\(\)](#), [matchReferrals\(\)](#)

**Examples**

```
stId <- "R-HSA-983150"
# matchPrecedingAndFollowingEvents(event.id=stId, depth=2, type="row")
```

---

matchReactionsInPathway

*MATCH Reactions in associated Pathway*

---

**Description**

This method could find all Reactions connected with a given Pathway by the relationship 'hasEvent'. Also, the input can be a Reaction, the result would then be Pathway(s) linked via 'hasEvent' together with other Reactions linked with the Pathways(s).

**Usage**

```
matchReactionsInPathway(
  event.id = NULL,
  event.displayName = NULL,
  species = NULL,
  type = c("row", "graph")
)
```

**Arguments**

`event.id`            stId or dbId of an Event  
`event.displayName`        displayName of an Event  
  
`species`            name or taxon id or dbId or abbreviation of a species  
`type`                return results as a list of dataframes (**'row'**), or as a graph object (**'graph'**)

**Value**

Reactions connected to the given Pathway/Reaction via 'hasEvent' relationships

**See Also**

Other match: [matchDiseases\(\)](#), [matchHierarchy\(\)](#), [matchInteractors\(\)](#), [matchObject\(\)](#), [matchPEroles\(\)](#), [matchPaperObjects\(\)](#), [matchPrecedingAndFollowingEvents\(\)](#), [matchReferrals\(\)](#)

**Examples**

```

reaction <- "R-HSA-1369062"
# matchReactionsInPathway(event.id=reaction, type="graph")
# matchReactionsInPathway("R-HSA-5682285", type="row")

```

---

|                |                                   |
|----------------|-----------------------------------|
| matchReferrals | <i>MATCH biological referrals</i> |
|----------------|-----------------------------------|

---

**Description**

This method retrieves Reactome objects that are connected with the given object in a *reverse* relationship. For example, to find Pathways containing the given Reaction.

**Usage**

```

matchReferrals(
  id = NULL,
  displayName = NULL,
  main = TRUE,
  depth = 1,
  all.depth = FALSE,
  species = NULL,
  type = c("row", "graph")
)

```

**Arguments**

|             |                                                                                                  |
|-------------|--------------------------------------------------------------------------------------------------|
| id          | stId or dbId of a Reactome object                                                                |
| displayName | displayName of a Reactome object                                                                 |
| main        | if set to TRUE, only <b>first-class</b> referrals returned                                       |
| depth       | number of depths                                                                                 |
| all.depth   | if set to TRUE, connected objects in all depths returned                                         |
| species     | name or taxon id or dbId or abbreviation of a species                                            |
| type        | return results as a list of dataframes ( <b>'row'</b> ), or as a graph object ( <b>'graph'</b> ) |

**Details**

For now it just focuses on biological referrals in the following Classes: "Event", "PhysicalEntity", "Regulation", "CatalystActivity", "ReferenceEntity", "Interaction", "AbstractModifiedResidue".

**Value**

referrals of the given instance

**See Also**

Other match: [matchDiseases\(\)](#), [matchHierarchy\(\)](#), [matchInteractors\(\)](#), [matchObject\(\)](#), [matchPERoles\(\)](#), [matchPaperObjects\(\)](#), [matchPrecedingAndFollowingEvents\(\)](#), [matchReactionsInPathway\(\)](#)

**Examples**

```
stId <- "R-HSA-112479"
# matchReferrals("R-HSA-112479", main=FALSE, all.depth=TRUE, type="row")
```

---

multiObjects

*Retrieve multiple Reactome objects*


---

**Description**

The [matchObject](#) function takes only one id/name at a time, this method allows you to input many ids and get an aggregated table for their detailed information. It can only accept **ids** for now.

**Usage**

```
multiObjects(ids, databaseName = "Reactome", speedUp = FALSE, cluster = 2)
```

**Arguments**

|              |                                                      |
|--------------|------------------------------------------------------|
| ids          | Reactome stIds/dbIds, or non-Reactome ids            |
| databaseName | database name                                        |
| speedUp      | set TRUE to use <a href="#">doParallel</a> method    |
| cluster      | the number of cluster in <a href="#">makeCluster</a> |

**Value**

Reactome database objects for the given ids

**See Also**

[matchObject](#) for details

**Examples**

```
## "ids" can be Reactome or non-Reactome ids
ids <- c("P02741", "P08887", "P08505", "Q9GZQ8")
#res <- multiObjects(ids, databaseName="UniProt", speedUp=TRUE)
```

---

unnestListCol

*Unnest a column of lists in a dataframe*

---

**Description**

Unnest a column of lists in a dataframe

**Usage**

```
unnestListCol(df, column = "properties")
```

**Arguments**

|        |                                         |
|--------|-----------------------------------------|
| df     | dataframe where a column to be unnested |
| column | specific column to be unnested          |

**Value**

an unnested dataframe for network visualization

**Examples**

```
# nodes <- unnestListCol(graph$nodes, "properties")
```

# Index

## \* **internal**

ReactomeGraph4R-package, [2](#)

## \* **match**

matchDiseases, [3](#)

matchHierarchy, [4](#)

matchInteractors, [5](#)

matchObject, [6](#)

matchPaperObjects, [7](#)

matchPEroles, [8](#)

matchPrecedingAndFollowingEvents,  
[9](#)

matchReactionsInPathway, [10](#)

matchReferrals, [11](#)

doParallel, [12](#)

login, [3](#)

makeCluster, [12](#)

matchDiseases, [3](#), [5](#), [7–12](#)

matchHierarchy, [4](#), [4](#), [5](#), [7–12](#)

matchInteractors, [4](#), [5](#), [5](#), [7–12](#)

matchObject, [4](#), [5](#), [6](#), [8–13](#)

matchPaperObjects, [4](#), [5](#), [7](#), [7](#), [9–12](#)

matchPEroles, [4](#), [5](#), [7](#), [8](#), [8](#), [10–12](#)

matchPrecedingAndFollowingEvents, [4](#), [5](#),  
[7–9](#), [9](#), [11](#), [12](#)

matchReactionsInPathway, [4](#), [5](#), [7–10](#), [10](#),  
[12](#)

matchReferrals, [4](#), [5](#), [7–11](#), [11](#)

multiObjects, [7](#), [12](#)

ReactomeGraph4R

(ReactomeGraph4R-package), [2](#)

ReactomeGraph4R-package, [2](#)

unnestListCol, [13](#)