

# 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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ReactomeGraph4R-package

*ReactomeGraph4R: 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.

## 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>
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---

**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>
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---

**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>
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---

**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>
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---

**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**

event.id	stId/dbId of an Event
event.displayName	displayName of an Event
species	name or taxon id or dbId or abbreviation of specified species
depth	number of depths
all.depth	if set to TRUE, all RLE(s) connected to the given Event in all depths returned
type	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")
```

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