

# Package ‘omXplore’

May 26, 2026

**Type** Package

**Title** Vizualization tools for 'omics' datasets with R

**Version** 1.7.0

**Description** This package contains a collection of functions (written as shiny modules) for the visualisation and the statistical analysis of omics data. These plots can be displayed individually or embedded in a global Shiny module.

Additionally, it is possible to integrate third party modules to the main interface of the package omXplore.

**License** Artistic-2.0

**Depends** R (>= 4.5.0), methods

**Imports** DT, shiny, MSnbase, PSMATCH, SummarizedExperiment, MultiAssayExperiment, shinyBS, shinyjs, shinyjqui, RColorBrewer, gplots, plotly, visNetwork, tibble, grDevices, stats, utils, htmlwidgets, vioplot, graphics, FactoMineR, dendextend, dplyr, factoextra, tidyr, nipals, Biobase

**Suggests** knitr, rmarkdown, BiocStyle, testthat, Matrix, graph

**biocViews** Software, ShinyApps, MassSpectrometry, DataRepresentation, GUI, QualityControl

**NeedsCompilation** no

**Collate** 'mod\_explore\_graphs.R' 'Infos\_adjacencyMatrix.R' 'Prostar\_1x.R' 'convert\_to\_mae.R' 'doc-data.R' 'external\_apps\_examples.R' 'get\_pep\_prot\_CC.R' 'mae\_accessors.R' 'metacell\_utils.R' 'mod\_colorLegend.R' 'modules.R' 'omXplore\_cc.R' 'omXplore\_corrmatrix.R' 'omXplore\_density.R' 'omXplore\_format\_DT.R' 'omXplore\_heatmap.R' 'omXplore\_intensity.R' 'omXplore\_PCA\_nipals.R' 'omXplore\_pca.R' 'omXplore\_plots\_tracking.R' 'omXplore\_tabExplorer.R' 'omXplore\_variance.R' 'omXplore\_view\_dataset.R' 'palette.R' 'plot\_boxplot.R' 'plot\_heatmap.R' 'plot\_pca.R' 'plot\_violin.R' 'utils.R' 'zzz.R'

**RoxygenNote** 7.3.3

**Encoding** UTF-8

**LazyData** false

**URL** <https://github.com/edyp-lab/omXplore>,  
<https://edyp-lab.github.io/omXplore/>

**BugReports** <https://github.com/edyp-lab/omXplore/issues>

**Roxygen** list(markdown = TRUE)

**VignetteBuilder** knitr

**git\_url** <https://git.bioconductor.org/packages/omXplore>

**git\_branch** devel

**git\_last\_commit** 2b3acd9

**git\_last\_commit\_date** 2026-04-28

**Repository** Bioconductor 3.24

**Date/Publication** 2026-05-25

**Author** Samuel Wiczorek [aut, cre] (ORCID:  
<https://orcid.org/0000-0002-5016-1203>),  
 Thomas Burger [aut],  
 Enora Fremy [ctb],  
 Cyril Ariztegui [ctb],  
 Manon Gaudin [ctb]

**Maintainer** Samuel Wiczorek <samuel.wiczorek@cea.fr>

## Contents

accessors	3
BuildColorStyles	5
Build_enriched_qdata	5
color-legend	6
converters	7
corrmatrix	9
density-plot	10
ds-cc	11
ds-pca	12
ds-view	14
external_app	16
FormatDataForDT	17
format_DT	18
GetPkgVersion	19
globals	20
intensity-plots	20
is.listOf	22
omXplore-modules	22
omXplore_heatmap	23
omXplore_tabExplorer	25
palette	26
pep_prot_CC	27
pkgs.require2	29
plot-variance	29
plots_tracking	30
Prostar-1x-compatible	31
q_metadata	32

<i>accessors</i>	3
sub_R25 . . . . .	33
vdata . . . . .	34

**Index** **35**

---

<i>accessors</i>	<i>Accessors functions</i>
------------------	----------------------------

---

**Description**

Functions used to access the additional plots in the instances of the class `MultiAssayExperiment`.

**Usage**

```

get_adjacencyMatrix(object, ...)

## S4 method for signature 'SummarizedExperiment'
get_adjacencyMatrix(object)

get_design(object, ...)

## S4 method for signature 'MultiAssayExperiment'
get_design(object)

get_group(object, ...)

## S4 method for signature 'MultiAssayExperiment'
get_group(object)

get_metacell(object, ...)

## S4 method for signature 'SummarizedExperiment'
get_metacell(object, slot.name = c("metacell", "qMetacell"))

get_cc(object, ...)

## S4 method for signature 'SummarizedExperiment'
get_cc(object)

get_parentProtId(object, ...)

## S4 method for signature 'SummarizedExperiment'
get_parentProtId(object)

get_colID(object, ...)

## S4 method for signature 'SummarizedExperiment'
get_colID(object)

get_type(object, ...)

## S4 method for signature 'SummarizedExperiment'

```

```

get_type(object)

get_pkg_version(object, ...)

## S4 method for signature 'SummarizedExperiment'
get_pkg_version(object)

```

### Arguments

object	An instance of class SummarizedExperiment.
...	Additional parameters
slot.name	The name of the slot dedicated to cell metadata to search. Default values are 'metacell' and 'qMetacell'

### Value

See individual method description for the return value.

If exists, the slot value requested.

A DataFrame containing the adjacency matrix of the dataset

A data.frame containing the metadata of the dataset

A data.frame containing the metadata of the dataset

A data.frame containing the metadata of the dataset

A data.frame containing the metadata of the dataset

A data.frame containing the metadata of the dataset

A data.frame containing the metadata of the dataset

A data.frame containing the metadata of the dataset

A data.frame containing the metadata of the dataset

### Examples

```

## -----
## Accessing slots from a MSnSet dataset
## -----
data(sub_R25)
se1 <- sub_R25[[1]]
parentProtId <- get_parentProtId(se1)
colID <- get_colID(se1)
type <- get_type(se1)
metacell <- get_metacell(se1)
conds <- get_group(sub_R25)

```

---

BuildColorStyles	<i>Build color style for DT tables</i>
------------------	--

---

**Description**

This function builds a list which is used for styling DT tables with the function `DT::styleEqual()`

**Usage**

```
BuildColorStyles(type)
```

**Arguments**

type                    The type of dataset. Available values are protein and peptide

**Value**

A list

**Examples**

```
NULL
```

---

Build_enriched_qdata	<i>Builds enriched assay with cell metadata info</i>
----------------------	--

---

**Description**

If the cell metadata exists in the object of class `SummarizedExperiment`, then these information are added to the quantitative data so as to use styles with the functions of the package DT.

**Usage**

```
Build_enriched_qdata(obj)
```

**Arguments**

obj                    An instance of the class `SummarizedExperiment`

**Value**

A data.frame with new columns corresponding to the cell metadata (if exists)

**Examples**

```
NULL
```

---

`color-legend`*Color legend for DaparToolshed*

---

**Description**

Shows a legend based on the tags in the package 'DaparToolshed'

**Usage**

```
custom_metacell_colors()

colorLegend_ui(id)

colorLegend_server(
  id,
  presentTags = reactive({
    NULL
  }),
  hide.white = TRUE
)

colorLegend(dataIn = SummarizedExperiment::SummarizedExperiment())
```

**Arguments**

<code>id</code>	A character(1) which is the id of the shiny module.
<code>presentTags</code>	A vector of character() which correspond to the tags.
<code>hide.white</code>	A boolean() to indicate whether the white cells must be hidden or not.
<code>dataIn</code>	An instance of the class SummarizedExperiment.

**Value**

A vector  
NA  
NA  
A shiny app

**Examples**

```
if (interactive()) {
  data(vdata)
  shiny::runApp(colorLegend(vdata[[1]]))
}
```

---

converters

*Convert to enriched MultiAssayExperiment*

---

### **Description**

The resulting object is an instance of the `MultiAssayExperiment` class.

### **Usage**

```
convert_to_mae(obj)
MSnSet_to_mae(obj)
matrix_to_mae(obj)
df_to_mae(obj)
Compute_CC(obj)
QFeatures_to_mae(obj)
SE_to_mae(obj)
MAE_to_mae(obj)
Check_se_Consistency(obj)
list_to_se(l1)
Check_List_consistency(l1)
listOfLists_to_mae(obj, colData = NULL)
listOfSE_to_mae(obj)
Check_MSnSet_Consistency(obj)
matrix_to_se(obj)
df_to_se(obj)
MSnSet_to_se(obj)
Build_X_CC(se)
listOfMSnSet_to_mae(obj)
listOfmatrix_to_mae(obj)
listOfdf_to_mae(obj)
```

**Arguments**

obj	An object compliant with the formats accepted by omXplore
ll	A list
colData	A data.frame()
se	AN instance of the class SummarizedExperiment

**Value**

An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 An instance of SimpleList  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 A boolean(1)  
 An enriched instance of the class SummarizedExperiment  
 A boolean(1)  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 A boolean(1)  
 An enriched instance of the class SummarizedExperiment  
 An enriched instance of the class SummarizedExperiment  
 An enriched instance of the class SummarizedExperiment  
 An enriched instance of the class SummarizedExperiment  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment  
 An enriched instance of the class MultiAssayExperiment

**Examples**

```

#-----
# Conversion of a MultiAssayExperiment instance
#-----
data(miniACC, package = "MultiAssayExperiment")
convert_to_mae(miniACC)

```

---

corrmatrix	<i>Displays a correlation matrix of the quantitative data of a numeric matrix.</i>
------------	--

---

### Description

Displays a correlation matrix of the quantitative data of a numeric matrix.

### Usage

```
omXplore_corrmatrix_ui(id)

omXplore_corrmatrix_server(
  id,
  dataIn = reactive({
    NULL
  }),
  i = reactive({
    1
  })
)

corrMatrix(data, rate = 0.5, showValues = FALSE)

omXplore_corrmatrix(dataIn, i)
```

### Arguments

id	A character(1) which is the id of the shiny module.
dataIn	An instance of the class SummarizedExperiment
i	An integer which is the index of the assay in the param dataIn
data	An object of class 'matrix'
rate	The rate parameter to control the exponential law for the gradient of colors
showValues	A boolean which indicates whether to show values in the correlation plot.

### Value

NA  
 NA  
 A plot  
 A shiny app

### Examples

```
if (interactive()) {
  data(vdata)
  omXplore_corrmatrix(vdata, 1)
}
```

---

density-plot	<i>Displays a correlation matrix of the quantitative data of a numeric matrix.</i>
--------------	--

---

### Description

Displays a correlation matrix of the quantitative data of a numeric matrix.

### Usage

```
omXplore_density_ui(id)

omXplore_density_server(
  id,
  dataIn = reactive({
    NULL
  }),
  i = reactive({
    1
  }),
  pal.name = reactive({
    NULL
  })
)

densityPlot(data, conds = NULL, pal.name = NULL)

omXplore_density(dataIn, i)
```

### Arguments

id	A character(1) which is the id of the shiny module.
dataIn	An instance of the class SummarizedExperiment
i	An integer which is the index of the assay in the param obj
pal.name	A character(1) which is the name of the palette from the package <a href="#">RColorBrewer::RColorBrewer</a> from which the colors are taken. Default value is 'Set1'.
data	A data.frame() of quantitative data
conds	A vector indicating the name of each sample.

### Value

NA  
 NA  
 A plot  
 A shiny app

**Examples**

```

if (interactive()) {
  data(vdata)
  shiny::runApp(omXplore_density(vdata, 1))
}

if (interactive()) {
  data(vdata)
  qdata <- SummarizedExperiment::assay(vdata[[1]])
  conds <- get_group(vdata)
  densityPlot(qdata, conds)
}

```

ds-cc

*plots\_cc\_ui and plots\_cc\_server***Description**

A shiny Module.

**Usage**

```

omXplore_cc_ui(id)

omXplore_cc_server(
  id,
  dataIn = reactive({
    NULL
  }),
  i = reactive({
    NULL
  })
)

omXplore_cc(dataIn, i)

```

**Arguments**

<code>id</code>	A character(1) which is the id of the shiny module.
<code>dataIn</code>	An instance of SummarizedExperiment class
<code>i</code>	An integer which is the index of the assay in the param obj

**Value**

A shiny module  
A shiny plot  
A shiny module  
A shiny app  
A shiny app

**Examples**

```
if (interactive()) {
  data(vdata)
  shiny::runApp(omXplore_cc(vdata, 1))
}
```

---

ds-pca

*my\_PCA*


---

**Description**

Process a PCA, using `nipals` or `FactoMineR`, on a quantitative dataset.

This method plots a bar plot which represents the distribution of the number of missing values (NA) per lines (ie proteins).

- `wrapper_pca()`
- `plotPCA_Eigen_hc()`: plots the eigen values of PCA with the `plotly` library
- `plotPCA_Eigen()`: plots the eigen values of PCA
- `plotPCA_Var()`
- `plotPCA_Ind()`

**Usage**

```
my_PCA(
  X,
  scale.unit = TRUE,
  ncp = min(12, nrow(X) - 1, ncol(X)),
  ind.sup = NULL,
  quanti.sup = NULL,
  quali.sup = NULL,
  row.w = NULL,
  col.w = NULL,
  graph = FALSE,
  axes = c(1, 2),
  approach = "FM",
  gramschmidt = TRUE
)
```

```
omXplore_pca_ui(id)
```

```
omXplore_pca_server(id, dataIn, i)
```

```
omXplore_pca(dataIn, i)
```

```
wrapper_pca(
  qdata,
  group,
  var.scaling = TRUE,
  ncp = NULL,
```

```

  approach = "FM",
  gramschmidt = TRUE
)

plotPCA_Eigen(res.pca)

plotPCA_Var(res.pca, chosen.axes = c(1, 2))

plotPCA_Ind(res.pca, chosen.axes = c(1, 2))

plotPCA_Eigen_hc(res.pca)

```

### Arguments

X	a data.frame() of quantitative data
scale.unit	See FactoMineR::PCA()
ncp	See FactoMineR::PCA()
ind.sup	See FactoMineR::PCA()
quanti.sup	See FactoMineR::PCA()
quali.sup	See FactoMineR::PCA()
row.w	See FactoMineR::PCA()
col.w	See FactoMineR::PCA()
graph	See FactoMineR::PCA()
axes	See FactoMineR::PCA()
approach	a string corresponding to the package to use for PCA (if no NA, default is "FM" for FactoMineR)
gramschmidt	A boolean indicating whether to use Gram-Schmidt orthogonalization or not.
id	A character(1) which is the id of the shiny module.
dataIn	An instance of the class MultiAssayExperiment.
i	An integer which is the index of the assay in the param obj
qdata	A data.frame() of quantitative data
group	A vector with the name of samples
var.scaling	A boolean indicating whether to scale the data or not
res.pca	The result of the function FactoMineR::PCA()
chosen.axes	See the parameter 'axes' of the function factoextra::fviz_pca_var()

### Value

res.pca a "PCA" "list" object

NA

NA

A shiny app

The result of the function FactoMineR::PCA()

A plot

A plot

A plot

A plot

**Author(s)**

Samuel Wiczorek, Enora Fremy

**Examples**

```

if (interactive()) {
  data(vdata)
  obj <- vdata[[1]]
  res.pca <- my_PCA(SummarizedExperiment::assay(obj), approach = "FM")
  plotPCA_Eigen(res.pca)
  plotPCA_Var(res.pca)
  plotPCA_Eigen_hc(res.pca)
  plotPCA_Ind(res.pca)
}

if (interactive()) {
  data(vdata)
  library(shiny)
  library(dplyr)
  library(plotly)
  # Replace missing values for the example
  sel <- is.na(SummarizedExperiment::assay(vdata, 1))
  SummarizedExperiment::assay(vdata[[1]])[sel] <- 0
  SummarizedExperiment::assay(vdata[[1]])[1, 1] <- NA
  omXplore_pca(vdata, 1)
}

if (interactive()) {
  data(vdata)
  obj <- vdata[[1]]
  res.pca <- wrapper_pca(qdata = SummarizedExperiment::assay(obj), group = get_group(obj))
  plotPCA_Eigen(res.pca)
  plotPCA_Var(res.pca)
  plotPCA_Eigen_hc(res.pca)
  plotPCA_Ind(res.pca)
}

```

---

ds-view

*Bar plot of missing values per lines using plotly.*

---

**Description**

This method plots a bar plot which represents the distribution of the number of missing values (NA) per lines (i.e. proteins).

**Usage**

```

view_dataset_ui(id)

view_dataset_server(
  id,
  dataIn = reactive({

```

```

      NULL
    }),
    addons = list(),
    verbose = FALSE
  )

view_dataset(dataIn = NULL, addons = NULL)

```

### Arguments

id	A character(1) for the 'id' of the shiny module. It must be the same as for the '*_ui' function.
dataIn	An instance of the class MultiAssayExperiment.
addons	A list to configure the other shiny apps to integrate. Each item correspond to one package: <ul style="list-style-type: none"> <li>• the name of the slot is the name of the package</li> <li>• the content of the slot is a vector composed of the generic name of the shiny app. Each of the apps listed here must be an exported app of the package. For example, given the value <code>addons = list(testPkg = c('foo', 'foo2'))</code>. That means that the package called "testPkg" must provide the four functions: <code>foo1_ui()</code>, <code>foo1_server()</code> and <code>foo2_ui()</code>, <code>foo2_server()</code></li> </ul>
verbose	A boolean for verbose mode. Default is FALSE.

### Details

- distribution of the missing values per line,
- a bar plot which represents the distribution of the number of missing values (NA) per lines (i.e. proteins) and per conditions,
- Histogram of missing values.
- Variance : Builds a densityplot of the CV of entities in numeric matrix. The CV is calculated for each condition present in the dataset (see the slot 'Condition' in the `colData()` DataFrame)
- Heatmap:

The function [heatmapD\(\)](#)

The function `[]` is inspired from the function 'heatmap.2' that displays a numeric matrix. For more information, please refer to the help of the heatmap.2 function.

### Value

NA

NA

NA

A shiny application which wraps the functions `view_dataset_ui()` and the `view_dataset_server()`

### Missing values

#' - distribution of the missing values per line,

- a bar plot which represents the distribution of the number of missing values (NA) per lines (ie proteins) and per conditions,
- Histogram of missing values.

**Author(s)**

Samuel Wieczorek, Enora Fremy

**Examples**

```

if (interactive()) {
  library(shiny)
  library(omXplore)
  data(vdata)
  addons <- list(omXplore = c("extFoo1", "extFoo2"))
  runApp(omXplore::view_dataset(vdata, addons))

  omXplore::view_dataset(vdata)
}

if (interactive()) {
  data(vdata)
  view_dataset(vdata)
}

```

---

external\_app

*External module example*

---

**Description**

Example for an external shiny module, well structured to be run within a workflow for MagellanNTK

**Usage**

```

extFoo1_ui(id)

extFoo1_server(
  id,
  dataIn = reactive({
    NULL
  }),
  i = reactive({
    NULL
  })
)

extFoo1(dataIn, i)

extFoo2_ui(id)

extFoo2_server(
  id,
  dataIn = reactive({
    NULL
  }),
  i = reactive({

```

```

        NULL
      })
    )

    extFoo2(dataIn, i)

```

**Arguments**

`id` A character(1) which is the id of the shiny module.

`dataIn` An object of instance MultiAssayExperiment

`i` An integer which is the index of the assay in the param obj

**Value**

NA

NA

NA

A shiny app

NA

NA

A shiny app

**Examples**

```

if (interactive()) {
  data(vdata)
  app1 <- extFoo1(vdata, 1)
  app2 <- extFoo2(vdata, 1)
  shiny::runApp(app1)
  shiny::runApp(app2)
}

```

---

FormatDataForDT

*Constructs a dataset suitable to use with the module format\_DT.*


---

**Description**

This function builds the skeleton of a dataset which can be used by the module formatDT. It creates additional columns to be used to style the table. to colors cells.

**Usage**

```
FormatDataForDT(se, digits = 2)
```

**Arguments**

`se` An instance of the class SummarizedExperiment

`digits` An 'integer(1)' to specify the number of digits to display in the tables for numerical values. Default is 2.

**Value**

A data.frame

**Examples**

NULL

---

format\_DT

*formatDT\_ui and formatDT\_server*

---

**Description**

A shiny Module.

See DT package homepage for more details about styling tables. If no style is precised, this module show the raw data. If any style is given, then the dataset must be well configured (I.e. it must contain the correct columns )

**Usage**

```
formatDT_ui(id)

formatDT_server(
  id,
  data = reactive({
    NULL
  }),
  data_nostyle = reactive({
    NULL
  }),
  withDLBtns = FALSE,
  showRownames = FALSE,
  dt_style = reactive({
    NULL
  }),
  filename = "Prostar_export",
  selection = "single"
)

formatDT(data)
```

**Arguments**

id	shiny id
data	A data.frame
data_nostyle	A data.frame() to be bind to the main data with no custom style
withDLBtns	A boolean to indicate whether to display download buttons or not.
showRownames	A boolean to indicate whether to show rownames.
dt_style	A list composed of:

- data : a data.frame
  - colors : a named vector
- filename      A character(1) which is the default filename for download.
- selection     A character(1) which indicates the type of selection. Default is 'single'.

**Value**

NA

NA

NA

NA

**Examples**

```
if (interactive()) {  
  data(vdata)  
  formatDT(vdata)  
}
```

---

GetPkgVersion	<i>Package version</i>
---------------	------------------------

---

**Description**

Gets the version number of a package

**Usage**

```
GetPkgVersion(pkg)
```

**Arguments**

pkg            The name of the package

**Value**

A character(1) with the name of the package and its version number.

**Examples**

```
GetPkgVersion("omXplore")
```

---

`globals`*Global variables*

---

**Description**

Defines the global variables for the package omXplore

**Usage**

```
globals()
```

**Value**

A list

**Examples**

```
globals()
```

---

`intensity-plots`*Displays different intensity plots.*

---

**Description**

Displays different intensity plots.

**Usage**

```
omXplore_intensity_ui(id)

omXplore_intensity_server(
  id,
  dataIn = reactive({
    NULL
  }),
  i = reactive({
    1
  }),
  track.indices = reactive({
    NULL
  }),
  remoteReset = reactive({
    NULL
  }),
  is.enabled = reactive({
    TRUE
  })
)
```

```
omXplore_intensity(dataIn, i = NULL, withTracking = FALSE)
boxPlot(dataIn, conds, legend = NULL, pal = NULL, subset = NULL)
violinPlot(data, conds, subset = NULL, pal.name = "Set1")
```

### Arguments

<code>id</code>	A character(1) which is the id of the shiny module.
<code>dataIn</code>	xxxx
<code>i</code>	An integer which is the index of the assay in the param obj
<code>track.indices</code>	A vector of integers which are the indices of lines to track.
<code>remoteReset</code>	An integer to activate the reset functions
<code>is.enabled</code>	A boolean that indicates whether the widgets should be enabled or disabled. Default is TRUE
<code>withTracking</code>	A boolean(1) indicating whether the tracking option is activated or not.
<code>conds</code>	A vector indicating the name of each sample.
<code>legend</code>	A vector of the conditions (one condition per sample).
<code>pal</code>	A basis palette for the boxes which length must be equal to the number of unique conditions in the dataset.
<code>subset</code>	A integer() vector of index indicating the indices of rows in the dataset to highlight
<code>data</code>	xxxx
<code>pal.name</code>	A character(1) which is the name of the palette from the package <a href="#">RColorBrewer::RColorBrewer</a> from which the colors are taken. Default value is 'Set1'.

### Value

NA  
 NA  
 A shiny app  
 A boxplot  
 A violin plot

### Author(s)

Samuel Wieczorek, Anais Courtier, Enora Fremy

### Examples

```
if (interactive()) {
  data(vdata)
  shiny::runApp(omXplore_intensity(vdata, 1))

  data(sub_R25)
  conds <- legend <- SummarizedExperiment::colData(sub_R25)$group
  pal <- ExtendPalette(length(unique(conds)))
  boxPlot(sub_R25[[1]], conds, legend, pal, seq_len(10))
}
```

```

    shiny::runApp(omXplore_intensity(sub_R25, 1, withTracking = TRUE))
  }

```

---

is.listOf	<i>Checks the class of a list's slots</i>
-----------	---

---

### Description

Checks if all slots of the given list are of the same class.

### Usage

```
is.listOf(object, obj.class = NULL)
```

### Arguments

object	A list
obj.class	The name of the class to search in items of the list.

### Value

A character (1) with the name of the package or NULL

### Examples

```

ll <- as.list(LETTERS[1:3])
is.listOf(ll, "data.frame")
is.listOf(ll, "character")

```

---

omXplore-modules	<i>Shiny modules used by omXplore</i>
------------------	---------------------------------------

---

### Description

These functions are relative to external modules that can be added into omXplore UI:

- `listShinyApps()`: Show the shiny modules recognized by omXplore and ready to be integrated in the UI of the function `view_dataset()`
- `listPlotModules()`: Show the shiny modules function names (only prefixes) recognized by omXplore and ready to use in the UI.
- `addModules()`: Add external shiny module(s) to the R global environment in such a way (specific prefix renaming of the functions) that it can be discovered by the function `view_dataset()` of the package omXplore during its launch.

**Usage**

```
addModules(addons = list())

listShinyApps(location = "both")

listPlotModules(location = "both")
```

**Arguments**

addons	A list in which each item: <ul style="list-style-type: none"> <li>• is named by the name of a package containing the modules to add,</li> <li>• contains the name of the shiny modules to integrate (without <code>'_ui'</code> nor <code>'_server'</code> suffixes)</li> </ul>
location	A character(0) to indicate which modules to list. Available values are: <code>'builtin'</code> , <code>'external'</code> and <code>'both'</code> (default).

**Value**

```
NA
A vector
A vector
```

**Examples**

```
listShinyApps()
listPlotModules()

#####
# Integration of a module in the package 'mypackage'
#####
addons <- list(omXplore = c("extFoo1", "extFoo2"))
addModules(addons)
```

---

omXplore_heatmap	<i>Displays a correlation matrix of the quantitative data of a numeric matrix.</i>
------------------	--

---

**Description**

This function is a wrapper to `heatmap.2()` that displays assay data in an instance of `SummarizedExperiment`. For more details, see `heatmap.2()`.

**Usage**

```
omXplore_heatmap_ui(id)

omXplore_heatmap_server(
  id,
  dataIn = reactive({
```

```

        NULL
    }},
    i = reactive({
        NULL
    })
)

omXplore_heatmap(dataIn, i)

heatmapD(
  qdata,
  conds,
  distance = "euclidean",
  cluster = "complete",
  dendro = FALSE
)

mv.heatmap(
  x,
  col = grDevices::heat.colors(100),
  srtCol = NULL,
  labCol = NULL,
  labRow = NULL,
  key = TRUE,
  key.title = NULL,
  main = NULL,
  ylab = NULL
)

heatmapForMissingValues(
  x,
  col = NULL,
  srtCol = NULL,
  labCol = NULL,
  labRow = NULL,
  key = TRUE,
  key.title = NULL,
  main = NULL,
  ylab = NULL
)

```

### Arguments

id	A character(1) which is the id of the shiny module.
dataIn	An instance of a class MultiAssayExperiment.
i	An integer which is the index of the assay in the param obj
qdata	A data.frame() of quantitative data.
conds	A vector indicating the name of each sample.
distance	The distance used by the clustering algorithm to compute the dendrogram.
cluster	the clustering algorithm used to build the dendrogram.
dendro	A boolean to indicate if the dendrogram has to be displayed

x	A matrix or array containing the quantitative data.
col	Colors used for the image. Defaults to heat colors (heat.colors).
srtCol	Angle of column conds, in degrees from horizontal
labCol	Character vectors with column conds to use.
labRow	Character vectors with row conds to use.
key	Logical indicating whether a color-key should be shown.
key.title	Main title of the color key. If set to NA no title will be plotted.
main	Main title; default to none.
ylab	y-axis title; default to none.

**Value**

NA  
 NA  
 A shiny app  
 A heatmap  
 A heatmap  
 A heatmap

**Author(s)**

Florence Combes, Samuel Wiczorek, Enora Fremy

**Examples**

```
if (interactive()) {
  data(vdata)
  omXplore_heatmap(vdata, 1)
}
```

---

omXplore\_tabExplorer *Explore MultiAssayExperiment objects.*

---

**Description**

Explore MultiAssayExperiment objects.

**Usage**

```
omXplore_tabExplorer_ui(id)

omXplore_tabExplorer_server(
  id,
  dataIn = reactive({
    NULL
  }),
  i = reactive({
```

```

      NULL
    }},
    digits = reactive({
      3
    })
  )

  omXplore_tabExplorer(dataIn, i)

```

### Arguments

<code>id</code>	A character(1) which is the id of the shiny module.
<code>dataIn</code>	An instance of the class <code>MultiAssayExperiment</code>
<code>i</code>	An integer which is the index of the assay in the param obj
<code>digits</code>	An integer for the number of digits shown in the table

### Value

NA  
 NA  
 NA  
 A shiny app

### Examples

```

if (interactive()) {
  data(vdata)
  shiny::runApp(omXplore_tabExplorer(vdata, 1))
}

NULL

```

---

palette

*Palette for samples.*

---

### Description

Builds a vector of `#conditions` colors for a set of samples. One color is given for a given condition. This function extends a base palette from the package `RColorBrewer::RColorBrewer` to 'n' colors. The colors in the returned palette are always in the same order

### Usage

```

SampleColors(conds, pal.name = "Set1")

ExtendPalette(n, pal.name = "Set1")

GetColorsForConditions(conds, pal = NULL)

```

**Arguments**

conds	A character() of conditions. The length of the vector is the number of samples in the dataset.
pal.name	A character(1) which is the name of the palette from the package <a href="#">RColorBrewer::RColorBrewer</a> from which the colors are taken. Default value is 'Set1'.
n	The number of desired colors in the palette
pal	A vector of HEX color code that form the basis palette from which to build the complete color vector for the conditions.

**Value**

A vector  
 A vector  
 A vector

**Examples**

```
if (interactive()) {
  #-----
  # Builds a palette for a dataset with 3 conditions
  # of 3 samples each.
  #-----

  conds <- c(rep("A", 3), rep("B", 3), rep("C", 3))
  SampleColors(conds)
  SampleColors(conds, "Dark2")

  #-----
  # Extend the default palette to 12 colors
  #-----

  ExtendPalette(12)

  data(vdata)
  conds <- get_group(vdata)
  GetColorsForConditions(conds, ExtendPalette(2))
}
```

---

 pep\_prot\_CC

*Display a CC*


---

**Description**

Display a CC  
 Connected Components infos

**Usage**

```
buildGraph(cc = NULL, meta = NULL)

display.CC.visNet(g = NULL, layout = "layout_with_fr")

plotCCJitter(df)

GetCCInfos(cc)
```

**Arguments**

<code>cc</code>	A list of connected component
<code>meta</code>	A <code>data.frame()</code>
<code>g</code>	An instance of a graph
<code>layout</code>	A character(1) which is the layout used in <code>visNetwork</code> . Default value is 'layout_with_fr'
<code>df</code>	A <code>data.frame()</code>

**Value**

A list

A plot

A plot

A list of three items:

- `One_One`: the number of cc composed of one protein and one peptide
- `One_Multi`: the number of cc composed of one protein and several peptides
- `Multi_Multi`: the number of cc composed of several proteins and several (shared) peptides.

**Author(s)**

Thomas Burger, Samuel Wieczorek

**Examples**

```
if (interactive()) {
  data(sub_R25)
  se <- sub_R25[[1]]
  g <- buildGraph(get_cc(se)[[1]])
  display.CC.visNet(g)
}

if (interactive()) {
  data(sub_R25)
  GetCCInfos(get_cc(sub_R25[[1]]))
}
```

---

pkgs.require2	<i>Loads packages</i>
---------------	-----------------------

---

**Description**

Checks if a package is available to load it

**Usage**

```
pkgs.require2(ll.deps)
```

**Arguments**

ll.deps            A character() vector which contains packages names

**Value**

NA

**Author(s)**

Samuel Wieczorek

**Examples**

```
pkgs.require2("omXplore")
```

---

plot-variance	<i>Variance plot</i>
---------------	----------------------

---

**Description**

A shiny module which plots the variance of samples

**Usage**

```
omXplore_variance_ui(id)
```

```
omXplore_variance_server(id, dataIn, i, pal.name = NULL)
```

```
CVDist(dataIn, conds, pal.name = NULL)
```

```
omXplore_variance(dataIn, i)
```

**Arguments**

id	A character(1) which is the id of the shiny module.
dataIn	An matrix
i	An integer which is the index of the assay in the param obj
pal.name	A character(1) which is the name of the palette from the package <a href="#">RColorBrewer::RColorBrewer</a> from which the colors are taken. Default value is 'Set1'.
conds	A vector indicating the name of each sample.

**Value**

NA  
 NA  
 A plot  
 A shiny app

**Examples**

```
if (interactive()) {
  data(vdata)
  shiny::runApp(omXplore_variance(vdata, 1))
}
```

---

plots\_tracking

*plots\_tracking\_ui and plots\_tracking\_server*


---

**Description**

This shiny module provides a tool to select

**Usage**

```
plots_tracking_ui(id)

plots_tracking_server(
  id,
  dataIn = reactive({
    NULL
  }),
  remoteReset = reactive({
    NULL
  }),
  is.enabled = reactive({
    TRUE
  })
)

plots_tracking(obj)
```

**Arguments**

id	shiny id
remoteReset	A boolean(1) which indicates whether to show the 'Reset' button or not.
is.enabled	xxx
obj	An instance of the class MultiAssayExperiment

**Value**

NA  
 A list (same structure as the parameter params)  
 A shiny app

**Examples**

```
if (interactive()) {
  data(vdata)
  shiny::runApp(plots_tracking(vdata[[1]]))
}
NULL
```

---

Prostar-1x-compatible *Convert datasets exported by the package Prostar*

---

**Description**

Convert datasets exported by the package Prostar

**Usage**

```
SE_Compatibility_with_Prostar_1x(obj, se)
MAE_Compatibility_with_Prostar_1x(obj, mae)
```

**Arguments**

obj	An instance of the class MSnSet
se	An instance of the class SummarizedExperiment
mae	An instance of the class MultiAssayExperiment

**Value**

An enriched instance of the class SummarizedExperiment  
 An enriched instance of the class MultiAssayExperiment

**Examples**

```
data(sub_R25)
```

---

q\_metadata

*Quantitative metadata vocabulary for entities*


---

### Description

This function gives the vocabulary used for the quantitative metadata of each entity in each condition.

### Usage

```
metacell.def(level)
```

```
Parent(level, node = NULL)
```

```
Children(level, parent = NULL)
```

```
GetMetacellTags(metacells = NULL, level = NULL, onlyPresent = FALSE)
```

### Arguments

level	A string corresponding to the type of object
node	The name of the node for which one wants its parent
parent	The name of the parent node
metacells	A data.frame() representing the cell metadata
onlyPresent	A boolean(1)

### Value

A data.frame containing the different tags and corresponding colors for the level given in parameter

A list

A vector

A vector

A vector

### Glossary

Peptide-level vocabulary

├─ 'Any' │ │ │ │─ 1.0 'Quantified' │ │ │ │ │ │ │ │─ 1.1 "Quant. by direct id" (color 4, white) │ │ │ │ │ │ │ │─ 1.2 "Quant. by recovery" (color 3, lightgrey) │ │ │ │ │ │ │ │─ 2.0 "Missing" (no color) │ │ │ │ │ │ │ │─ 2.1 "Missing POV" (color 1) │ │ │ │ │ │ │ │─ 2.2 'Missing MEC' (color 2) │ │ │ │ │ │ │ │─ 3.0 'Imputed' │ │ │ │ │ │ │ │─ 3.1 'Imputed POV' (color 1) │ │ │ │ │ │ │ │─ 3.2 'Imputed MEC' (color 2)

Protein-level vocabulary: └─ 'Any' │ │ │ │─ 1.0 'Quantified' │ │ │ │ │ │ │ │─ 1.1 "Quant. by direct id" (color 4, white) │ │ │ │ │ │ │ │─ 1.2 "Quant. by recovery" (color 3, lightgrey) │ │ │ │ │ │ │ │─ 2.0 "Missing" │ │ │ │ │ │ │ │─ 2.1 "Missing POV" (color 1) │ │ │ │ │ │ │ │─ 2.2 'Missing MEC' (color 2) │ │ │ │ │ │ │ │─ 3.0 'Imputed' │ │ │ │ │ │ │ │─ 3.1 'Imputed POV' (color 1) │ │ │ │ │ │ │ │─ 3.2 'Imputed MEC' (color 2) │ │ │ │ │ │ │ │─ 4.0 'Combined tags' (color 3bis, lightgrey)

**Conversion to the glossary**

A generic conversion  
Conversion for Proline datasets  
Conversion from Maxquant datasets

**Author(s)**

Thomas Burger, Samuel Wieczorek  
Samuel Wieczorek

**Examples**

```
if (interactive()) {
  metacell.def("protein")
  metacell.def("peptide")
}

Parent("protein", "Missing")
Parent("protein", "Missing POV")
Parent("protein", c("Missing POV", "Missing MEC"))
Parent("protein", c("Missing", "Missing POV", "Missing MEC"))

Children("protein", "Missing")
Children("protein", "Missing POV")
Children("protein", c("Missing POV", "Missing MEC"))
Children("protein", c("Missing", "Missing POV", "Missing MEC"))
data(vdata)
metacells <- get_metacell(vdata[[1]])
level <- get_type(vdata[[1]])
GetMetacellTags(metacells, level)
```

---

sub\_R25

*Feature example data*

---

**Description**

sub\_R25 is a protein subset of the dataset 'Exp1\_R25\_pept' in the package DAPARdata.

**Format**

An instance of the class MultiAssayExperiment

**Value**

An enriched instance of the class MultiAssayExperiment

**Source**

sub\_R25 was built from the source code available in [inst/scripts/build\\_datasets.R](#)  
The DAPARdata package: <https://github.com/edyp-lab/DAPARdata>

---

vdata

*Feature example data*

---

**Description**

vdata is a small object for testing and demonstration.

**Format**

An instance of the class `MultiAssayExperiment`

**Value**

An enriched instance of the class `MultiAssayExperiment`

**Source**

vdata was built from the source code available in [inst/scripts/build\\_datasets.R](#)

# Index

- \* **datasets**
  - vdata, 34
- \* **data**
  - vdata, 34
- \* **internal**
  - ds-cc, 11
  - plots\_tracking, 30
- accessors, 3
- addModules (omXplore-modules), 22
- boxPlot (intensity-plots), 20
- Build\_enriched\_qdata, 5
- Build\_X\_CC (converters), 7
- BuildColorStyles, 5
- buildGraph (pep\_prot\_CC), 27
- Check\_List\_consistency (converters), 7
- Check\_MSnSet\_Consistency (converters), 7
- Check\_se\_Consistency (converters), 7
- Children (q\_metadata), 32
- color-legend, 6
- colorLegend (color-legend), 6
- colorLegend\_server (color-legend), 6
- colorLegend\_ui (color-legend), 6
- Compute\_CC (converters), 7
- convert\_to\_mae (converters), 7
- converters, 7
- corrMatrix (corrmatrix), 9
- corrmatrix, 9
- custom\_metacell\_colors (color-legend), 6
- CVDist (plot-variance), 29
- density-plot, 10
- densityPlot (density-plot), 10
- df\_to\_mae (converters), 7
- df\_to\_se (converters), 7
- display.CC.visNet (pep\_prot\_CC), 27
- ds-cc, 11
- ds-pca, 12
- ds-view, 14
- ExtendPalette (palette), 26
- external\_app, 16
- extFoo1 (external\_app), 16
- extFoo1\_server (external\_app), 16
- extFoo1\_ui (external\_app), 16
- extFoo2 (external\_app), 16
- extFoo2\_server (external\_app), 16
- extFoo2\_ui (external\_app), 16
- format\_DT, 18
- FormatDataForDT, 17
- formatDT (format\_DT), 18
- formatDT\_server (format\_DT), 18
- formatDT\_ui (format\_DT), 18
- get\_adjacencyMatrix (accessors), 3
- get\_adjacencyMatrix, SummarizedExperiment-method (accessors), 3
- get\_cc (accessors), 3
- get\_cc, SummarizedExperiment-method (accessors), 3
- get\_colID (accessors), 3
- get\_colID, SummarizedExperiment-method (accessors), 3
- get\_design (accessors), 3
- get\_design, MultiAssayExperiment-method (accessors), 3
- get\_group (accessors), 3
- get\_group, MultiAssayExperiment-method (accessors), 3
- get\_metacell (accessors), 3
- get\_metacell, SummarizedExperiment-method (accessors), 3
- get\_parentProtId (accessors), 3
- get\_parentProtId, SummarizedExperiment-method (accessors), 3
- get\_pkg\_version (accessors), 3
- get\_pkg\_version, SummarizedExperiment-method (accessors), 3
- get\_type (accessors), 3
- get\_type, SummarizedExperiment-method (accessors), 3
- GetCCInfos (pep\_prot\_CC), 27
- GetColorsForConditions (palette), 26
- GetMetacellTags (q\_metadata), 32
- GetPkgVersion, 19
- globals, 20

- heatmapD (omXplore\_heatmap), 23
- heatmapD(), 15
- heatmapForMissingValues  
(omXplore\_heatmap), 23
- intensity-plots, 20
- is.listOf, 22
- list\_to\_se (converters), 7
- listOfdf\_to\_mae (converters), 7
- listOfLists\_to\_mae (converters), 7
- listOfmatrix\_to\_mae (converters), 7
- listOfMSnSet\_to\_mae (converters), 7
- listOfSE\_to\_mae (converters), 7
- listPlotModules (omXplore-modules), 22
- listShinyApps (omXplore-modules), 22
- MAE\_Compatibility\_with\_Prostar\_1x  
(Prostar-1x-compatible), 31
- MAE\_to\_mae (converters), 7
- matrix\_to\_mae (converters), 7
- matrix\_to\_se (converters), 7
- metacell.def (q\_metadata), 32
- MSnSet\_to\_mae (converters), 7
- MSnSet\_to\_se (converters), 7
- mv.heatmap (omXplore\_heatmap), 23
- my\_PCA (ds-pca), 12
- omXplore-modules, 22
- omXplore\_cc (ds-cc), 11
- omXplore\_cc\_server (ds-cc), 11
- omXplore\_cc\_ui (ds-cc), 11
- omXplore\_corrmatrix (corrmatrix), 9
- omXplore\_corrmatrix\_server  
(corrmatrix), 9
- omXplore\_corrmatrix\_ui (corrmatrix), 9
- omXplore\_density (density-plot), 10
- omXplore\_density\_server (density-plot),  
10
- omXplore\_density\_ui (density-plot), 10
- omXplore\_heatmap, 23
- omXplore\_heatmap\_server  
(omXplore\_heatmap), 23
- omXplore\_heatmap\_ui (omXplore\_heatmap),  
23
- omXplore\_intensity (intensity-plots), 20
- omXplore\_intensity\_server  
(intensity-plots), 20
- omXplore\_intensity\_ui  
(intensity-plots), 20
- omXplore\_pca (ds-pca), 12
- omXplore\_pca\_server (ds-pca), 12
- omXplore\_pca\_ui (ds-pca), 12
- omXplore\_tabExplorer, 25
- omXplore\_tabExplorer\_server  
(omXplore\_tabExplorer), 25
- omXplore\_tabExplorer\_ui  
(omXplore\_tabExplorer), 25
- omXplore\_variance (plot-variance), 29
- omXplore\_variance\_server  
(plot-variance), 29
- omXplore\_variance\_ui (plot-variance), 29
- palette, 26
- Parent (q\_metadata), 32
- pep\_prot\_CC, 27
- pkgs.require2, 29
- plot-variance, 29
- plotCCJitter (pep\_prot\_CC), 27
- plotPCA\_Eigen (ds-pca), 12
- plotPCA\_Eigen\_hc (ds-pca), 12
- plotPCA\_Ind (ds-pca), 12
- plotPCA\_Var (ds-pca), 12
- plots\_tracking, 30
- plots\_tracking\_server (plots\_tracking),  
30
- plots\_tracking\_ui (plots\_tracking), 30
- Prostar-1x-compatible, 31
- q\_metadata, 32
- QFeatures\_to\_mae (converters), 7
- RColorBrewer::RColorBrewer, 10, 21, 26,  
27, 30
- SampleColors (palette), 26
- SE\_Compatibility\_with\_Prostar\_1.x  
(Prostar-1x-compatible), 31
- SE\_to\_mae (converters), 7
- sub\_R25, 33
- vdata, 34
- view\_dataset (ds-view), 14
- view\_dataset\_server (ds-view), 14
- view\_dataset\_ui (ds-view), 14
- violinPlot (intensity-plots), 20
- wrapper\_pca (ds-pca), 12