# Package: BiocGenerics (via r-universe)

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**Title** S4 generic functions used in Bioconductor

**Description** The package defines many S4 generic functions used in Bioconductor.

**biocViews** Infrastructure

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**Collate** S3-classes-as-S4-classes.R utils.R normarg-utils.R replaceSlots.R aperm.R append.R as.data.frame.R as.list.R as.vector.R cbind.R do.call.R duplicated.R eval.R Extremes.R format.R funprog.R get.R grep.R is.unsorted.R lapply.R mapply.R match.R mean.R nrow.R order.R paste.R rank.R rep.R row_colnames.R saveRDS.R sort.R start.R subset.R t.R table.R tapply.R unique.R unlist.R unsplit.R which.R which.min.R relist.R boxplot.R image.R density.R IQR.R mad.R residuals.R var.R weights.R xtabs.R setops.R annotation.R combine.R containsOutOfMemoryData.R dbconn.R dge.R dims.R fileName.R longForm.R normalize.R Ontology.R organism_species.R paste2.R path.R plotMA.R plotPCA.R score.R strand.R toTable.R type.R updateObject.R testPackage.R zzz.R

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---

`BiocGenerics-package` *S4 generic functions for Bioconductor*

---

## Description

S4 generic functions needed by many Bioconductor packages.

## Details

We divide the generic functions defined in the **BiocGenerics** package in 2 categories:

1. Functions already defined in base R or in CRAN package **generics**, and explicitly promoted to S4 generics in **BiocGenerics**
2. S4 generics specific to Bioconductor.

**(1) Functions defined in base R or CRAN package generics and explicitly promoted to S4 generics in the BiocGenerics package:**

Generics for functions defined in package **base**:

- BiocGenerics::aperm
- BiocGenerics::append
- BiocGenerics::as.data.frame
- BiocGenerics::as.list
- BiocGenerics::as.vector
- BiocGenerics::rbind, BiocGenerics::cbind
- BiocGenerics::do.call
- BiocGenerics::duplicated, BiocGenerics::anyDuplicated
- BiocGenerics::eval
- Extremes: BiocGenerics::pmax, BiocGenerics::pmin, BiocGenerics::pmax.int, BiocGenerics::pmin.int
- BiocGenerics::format
- funprog: BiocGenerics::Reduce, BiocGenerics::Filter, BiocGenerics::Find, BiocGenerics::Map, BiocGenerics::Position
- BiocGenerics::get, BiocGenerics::mget
- BiocGenerics::grep, BiocGenerics::grepl
- BiocGenerics::is.unsorted
- BiocGenerics::lapply, BiocGenerics::sapply
- BiocGenerics::mapply
- BiocGenerics::match, BiocGenerics::%in%
- BiocGenerics::nrow, BiocGenerics::ncol, BiocGenerics::NROW, BiocGenerics::NCOL
- BiocGenerics::order
- BiocGenerics::paste
- BiocGenerics::rank
- BiocGenerics::rep.int
- BiocGenerics::rownames, BiocGenerics::rownames<-, BiocGenerics::colnames, BiocGenerics::colnames<
- BiocGenerics::saveRDS
- BiocGenerics::sort
- BiocGenerics::start, BiocGenerics::start<-, BiocGenerics::end, BiocGenerics::end<-, BiocGenerics::width, BiocGenerics::width<-, BiocGenerics::pos
- BiocGenerics::subset
- BiocGenerics::t
- BiocGenerics::table
- BiocGenerics::tapply
- BiocGenerics::unique
- BiocGenerics::unlist
- BiocGenerics::unsplit
- BiocGenerics::which
- BiocGenerics::which.min, BiocGenerics::which.max

Generics for functions defined in package **utils**:

- BiocGenerics::relist

Generics for functions defined in package **graphics**:

- BiocGenerics::boxplot
- BiocGenerics::image

Generics for functions defined in package **stats**:

- BiocGenerics::density
- BiocGenerics::residuals
- BiocGenerics::weights
- BiocGenerics::xtabs

Generics for functions defined in CRAN package **generics**:

- setops: BiocGenerics::union, BiocGenerics::intersect, BiocGenerics::setdiff, BiocGenerics::setequal

**(2) S4 generics specific to Bioconductor:**

- annotation, annotation<-
- combine
- containsOutOfMemoryData
- dbconn, dbfile
- counts, counts<-, design, design<-, dispTable, dispTable<-, sizeFactors, sizeFactors<-, conditions, conditions<-, estimateSizeFactors, estimateDispersions, plotDispEsts
- dims, nrows, ncols,
- fileName
- longForm
- normalize
- Ontology
- organism, organism<-, species, species<-
- paste2
- path, path<-, basename, basename<-, dirname, dirname<-
- plotMA
- plotPCA
- score, score<-
- strand, strand<-, invertStrand
- toTable
- type, type<-
- updateObject

## Note

More generics can be added on request by sending an email to the Bioc-devel mailing list:

http://bioconductor.org/help/mailing-list/

Things that should NOT be added to the **BiocGenerics** package:

- Internal generic primitive functions like `length`, `dim`, `` `dim<-` ``, etc... See ?`InternalMethods` for the complete list. There are a few exceptions though, that is, the **BiocGenerics** package may actually redefine a few of those internal generic primitive functions as S4 generics when for example the signature of the internal generic primitive is not appropriate (this is the case for `BiocGenerics::cbind`).
- S3 and S4 group generic functions like `Math`, `Ops`, etc... See ?`groupGeneric` and ?`S4groupGeneric` for the complete list.
- Generics already defined in the **stats4** package.

## Author(s)

The Bioconductor Dev Team

## See Also

- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- `setGeneric` and `setMethod` for defining generics and methods.

## Examples

```
## List all the symbols defined in this package:
ls('package:BiocGenerics')
```

---

# annotation *Accessing annotation information*

---

## Description

Get or set the annotation information contained in an object.

## Usage

```
annotation(object, ...)
annotation(object, ...) <- value
```

## Arguments

| | |
|---|---|
| `object` | An object containing annotation information. |
| `...` | Additional arguments, for use in specific methods. |
| `value` | The annotation information to set on `object`. |

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- annotation,eSet-method in the **Biobase** package for an example of a specific annotation method (defined for eSet objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
annotation
showMethods("annotation")

library(Biobase)
showMethods("annotation")
selectMethod("annotation", "eSet")
```

---

# aperm *Transposing an array-like object*

---

## Description

Transpose an array-like object by permuting its dimensions.

This is a multidimensional generalization of the t() operator used for 2D-transposition.

NOTE: This man page is for the aperm *S4 generic function* defined in the **BiocGenerics** package. See ?base::aperm for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
aperm(a, perm, ...)
```

## Arguments

| | |
|---|---|
| a | An array-like object. |
| perm, ... | See ?base::aperm for a description of these arguments. |

## Value

A transposed version of array-like object a, with subscripts permuted as indicated by the perm vector.

## See Also

- base::aperm for the default aperm method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- aperm,SVT_SparseArray-method in the **SparseArray** package for an example of a specific aperm method (defined for SVT_SparseArray objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
aperm  # note the dispatch on the 'a' arg only
showMethods("aperm")
selectMethod("aperm", "ANY")  # the default method
```

---

# append *Append elements to a vector-like object*

---

## Description

Append (or insert) elements to (in) a vector-like object.

NOTE: This man page is for the append *S4 generic function* defined in the **BiocGenerics** package. See ?base::append for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like or data-frame-like) not supported by the default method.

## Usage

```
append(x, values, after=length(x))
```

## Arguments

| | |
|---|---|
| x | The vector-like object to be modified. |
| values | The vector-like object containing the values to be appended to x. values would typically be of the same class as x, but not necessarily. |
| after | A subscript, after which the values are to be appended. |

## Value

See ?base::append for the value returned by the default method.

Specific methods defined in Bioconductor packages will typically return an object of the same class as x and of length length(x) + length(values).

## See Also

- base::append for the default append method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- append,Vector,Vector-method in the **S4Vectors** package for an example of a specific append method (defined for Vector objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
append  # note the dispatch on the 'x' and 'values' args only
showMethods("append")
selectMethod("append", c("ANY", "ANY"))  # the default method
```

---

# as.data.frame *Coerce to a data frame*

---

## Description

Generic function to coerce to a data frame, if possible.

NOTE: This man page is for the as.data.frame *S4 generic function* defined in the **BiocGenerics** package. See ?base::as.data.frame for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
as.data.frame(x, row.names=NULL, optional=FALSE, ...)
```

## Arguments

| | |
|---|---|
| x | The object to coerce. |
| row.names, optional, ... | See ?base::as.data.frame for a description of these arguments. |

## Value

An ordinary data frame.

See ?base::as.data.frame for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- base::as.data.frame for the default as.data.frame method.
- toTable for an alternative to as.data.frame.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- as.data.frame,DataFrame-method in the **S4Vectors** package, and as.data.frame,IntegerRanges-method in the **IRanges** package, for examples of specific as.data.frame methods (defined for DataFrame and IntegerRanges objects, respectively).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
as.data.frame  # note the dispatch on the 'x' arg only
showMethods("as.data.frame")
selectMethod("as.data.frame", "ANY")  # the default method
```

---

# as.list *Coerce to a list*

---

## Description

Generic function to coerce to a list, if possible.

NOTE: This man page is for the as.list *S4 generic function* defined in the **BiocGenerics** package. See ?base::as.list for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
as.list(x, ...)
```

## Arguments

| | |
|---|---|
| x | The object to coerce. |
| ... | Additional arguments, for use in specific methods. |

## Value

An ordinary list.

## See Also

- base::as.list for the default as.list method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- as.list,List-method in the **S4Vectors** package for an example of a specific as.list method (defined for List objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
as.list
showMethods("as.list")
selectMethod("as.list", "ANY")  # the default method

library(S4Vectors)
showMethods("as.list")
## The as.list() method for List objects:
selectMethod("as.list", "List")
```

---

`as.vector` *Coerce an object into a vector*

---

## Description

Attempt to coerce an object into a vector of the specified mode. If the mode is not specified, attempt to coerce to whichever vector mode is considered more appropriate for the class of the supplied object.

NOTE: This man page is for the `as.vector` *S4 generic function* defined in the **BiocGenerics** package. See `?base::as.vector` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
as.vector(x, mode="any")
```

## Arguments

| | |
|---|---|
| `x` | The object to coerce. |
| `mode` | See `?base::as.vector` for a description of this argument. |

## Value

A vector.

See `?base::as.vector` for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- `base::as.vector` for the default `as.vector` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- as.vector,Rle-method in the **S4Vectors** package, and as.vector,AtomicList-method in the **IRanges** packages, for examples of specific `as.vector` methods (defined for Rle and AtomicList objects, respectively).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
as.vector  # note the dispatch on the 'x' arg only
showMethods("as.vector")
selectMethod("as.vector", "ANY")  # the default method
```

---

boxplot *Box plots*

---

## Description

Produce box-and-whisker plot(s) of the given (grouped) values.

NOTE: This man page is for the `boxplot` *S4 generic function* defined in the **BiocGenerics** package. See `?graphics::boxplot` for the default method (defined in the **graphics** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
boxplot(x, ...)
```

## Arguments

| | |
|---|---|
| `x, ...` | See `?graphics::boxplot`. |

## Value

See `?graphics::boxplot` for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- `graphics::boxplot` for the default `boxplot` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- boxplot,AffyBatch-method in the **affy** package for an example of a specific `boxplot` method (defined for AffyBatch objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
boxplot
showMethods("boxplot")
selectMethod("boxplot", "ANY")  # the default method

library(affy)
showMethods("boxplot")
## The boxplot() method for AffyBatch objects:
selectMethod("boxplot", "AffyBatch")
```

cbind *Combine objects by rows or columns*

## Description

rbind and cbind take one or more objects and combine them by columns or rows, respectively.

NOTE: This man page is for the rbind and cbind *S4 generic functions* defined in the **BiocGenerics** package. See ?base::cbind for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like or matrix-like) not supported by the default methods.

## Usage

```
rbind(..., deparse.level=1)
cbind(..., deparse.level=1)
```

## Arguments

| | |
|---|---|
| ... | One or more vector-like or matrix-like objects. These can be given as named arguments. |
| deparse.level | See ?base::cbind for a description of this argument. |

## Value

See ?base::cbind for the value returned by the default methods.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input objects.

## See Also

- base::cbind for the default rbind and cbind methods.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- rbind,RectangularData-method and cbind,DataFrame-method in the **S4Vectors** package for examples of specific rbind and cbind methods (defined for RectangularData derivatives and DataFrame objects, respectively).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
rbind  # note the dispatch on the '...' arg only
showMethods("rbind")
selectMethod("rbind", "ANY")  # the default method

cbind  # note the dispatch on the '...' arg only
showMethods("cbind")
```

```
selectMethod("cbind", "ANY")  # the default method

library(S4Vectors)
showMethods("rbind")
## The rbind() method for RectangularData derivatives:
selectMethod("rbind", "RectangularData")
## The cbind() method for DataFrame objects:
selectMethod("cbind", "DataFrame")
```

---

`combine` *Combining or merging different Bioconductor data structures*

---

## Description

The `combine` generic function handles methods for combining or merging different Bioconductor data structures. It should, given an arbitrary number of arguments of the same class (possibly by inheritance), combine them into a single instance in a sensible way (some methods may only combine 2 objects, ignoring `...` in the argument list; because Bioconductor data structures are complicated, check carefully that `combine` does as you intend).

## Usage

```
combine(x, y, ...)
```

## Arguments

| | |
|---|---|
| `x` | One of the values. |
| `y` | A second value. |
| `...` | Any other objects of the same class as `x` and `y`. |

## Details

There are two basic combine strategies. One is an intersection strategy. The returned value should only have rows (or columns) that are found in all input data objects. The union strategy says that the return value will have all rows (or columns) found in any one of the input data objects (in which case some indication of what to use for missing values will need to be provided).

These functions and methods are currently under construction. Please let us know if there are features that you require.

## Value

A single value of the same class as the most specific common ancestor (in class terms) of the input values. This will contain the appropriate combination of the data in the input values.

## Methods

The following methods are defined in the **BiocGenerics** package:

`combine(x=ANY, missing)` Return the first (x) argument unchanged.

`combine(data.frame, data.frame)` Combines two `data.frame` objects so that the resulting `data.frame` contains all rows and columns of the original objects. Rows and columns in the returned value are unique, that is, a row or column represented in both arguments is represented only once in the result. To perform this operation, `combine` makes sure that data in shared rows and columns are identical in the two data.frames. Data differences in shared rows and columns usually cause an error. `combine` issues a warning when a column is a `factor` and the levels of the factor in the two data.frames are different.

`combine(matrix, matrix)` Combined two `matrix` objects so that the resulting `matrix` contains all rows and columns of the original objects. Both matricies must have `dimnames`. Rows and columns in the returned value are unique, that is, a row or column represented in both arguments is represented only once in the result. To perform this operation, `combine` makes sure that data in shared rows and columns are all equal in the two matricies.

Additional `combine` methods are defined in the **Biobase** package for AnnotatedDataFrame, AssayData, MIAME, and eSet objects.

## Author(s)

Biocore

## See Also

- combine,AnnotatedDataFrame,AnnotatedDataFrame-method, combine,AssayData,AssayData-method, combine,MIAME,MIAME-method, and combine,eSet,eSet-method in the **Biobase** package for additional `combine` methods.
- `merge` for merging two data frames (or data-frame-like) objects.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
combine
showMethods("combine")
selectMethod("combine", c("ANY", "missing"))
selectMethod("combine", c("data.frame", "data.frame"))
selectMethod("combine", c("matrix", "matrix"))

## ---------------------------------------------------------------------
## COMBINING TWO DATA FRAMES
## ---------------------------------------------------------------------
x <- data.frame(x=1:5,
        y=factor(letters[1:5], levels=letters[1:8]),
        row.names=letters[1:5])
```

```
y <- data.frame(z=3:7,
        y=factor(letters[3:7], levels=letters[1:8]),
        row.names=letters[3:7])
combine(x,y)

w <- data.frame(w=4:8,
       y=factor(letters[4:8], levels=letters[1:8]),
       row.names=letters[4:8])
combine(w, x, y)

# y is converted to 'factor' with different levels
df1 <- data.frame(x=1:5,y=letters[1:5], row.names=letters[1:5])
df2 <- data.frame(z=3:7,y=letters[3:7], row.names=letters[3:7])
try(combine(df1, df2)) # fails
# solution 1: ensure identical levels
y1 <- factor(letters[1:5], levels=letters[1:7])
y2 <- factor(letters[3:7], levels=letters[1:7])
df1 <- data.frame(x=1:5,y=y1, row.names=letters[1:5])
df2 <- data.frame(z=3:7,y=y2, row.names=letters[3:7])
combine(df1, df2)
# solution 2: force column to be 'character'
df1 <- data.frame(x=1:5,y=I(letters[1:5]), row.names=letters[1:5])
df2 <- data.frame(z=3:7,y=I(letters[3:7]), row.names=letters[3:7])
combine(df1, df2)

## ---------------------------------------------------------------------
## COMBINING TWO MATRICES
## ---------------------------------------------------------------------
m <- matrix(1:20, nrow=5, dimnames=list(LETTERS[1:5], letters[1:4]))
combine(m[1:3,], m[4:5,])
combine(m[1:3, 1:3], m[3:5, 3:4]) # overlap
```

---

`containsOutOfMemoryData` *Does an object contain out-of-memory data?*

---

## Description

Some objects in Bioconductor can use on-disk or other out-of-memory representation for their data, typically (but not necessarily) when the data is too big to fit in memory. For example the data in a TxDb object is stored in an SQLite database, and the data in an HDF5Array object is stored in an HDF5 file.

The `containsOutOfMemoryData()` function determines whether an object contains out-of-memory data or not.

Note that objects with out-of-memory data are usually not compatible with a serialization/unserialization roundtrip. More concretely, `base::saveRDS()`/`base::readRDS()` tend to silently break them!

See `?saveHDF5SummarizedExperiment` in the **HDF5Array** package for a more extensive discussion about this.

## Usage

```
containsOutOfMemoryData(object)
```

## Arguments

| | |
|---|---|
| object | The object to be tested. |

## Details

An object can store *some* of its data on disk and *some* of it in memory. This is the case for example when a SummarizedExperiment object (or derivative) has some of its assays on disk (e.g. in HDF5Matrix objects) and others in memory (e.g. in ordinary matrices and/or SparseMatrix objects).

Of course in this case, `containsOutOfMemoryData()` will still return `TRUE`. In other words, `containsOutOfMemoryData(obj` will only return `FALSE` when all the data in `object` resides in memory, that is, when the object can safely be serialized.

## Value

`TRUE` or `FALSE`.

## Note

TO DEVELOPERS:

The **BiocGenerics** package also defines the following:

- A default `containsOutOfMemoryData()` method that returns `TRUE` if `object` is an S4 object with at least one slot for which `containsOutOfMemoryData()` is `TRUE` (recursive definition), and `FALSE` otherwise.
- A `containsOutOfMemoryData()` method for `list` objects that returns `TRUE` if `object` has at least one list element for which `containsOutOfMemoryData()` is `TRUE` (recursive definition), and `FALSE` otherwise.
- A `containsOutOfMemoryData()` method for `environment` objects that returns `TRUE` if `object` contains at least one object for which `containsOutOfMemoryData()` is `TRUE` (recursive definition), and `FALSE` otherwise.
- The OutOfMemoryObject class. This is a virtual S4 class with no slots that any class defined in Bioconductor that represents out-of-memory objects should extend.
- A `containsOutOfMemoryData()` method for OutOfMemoryObject derivatives that returns `TRUE`.

Therefore, if you implement a class that uses an out-of-memory representation, make sure that it contains the OutOfMemoryObject class. This will make `containsOutOfMemoryData()` return `TRUE` on your objects, so you don't need to define a `containsOutOfMemoryData()` method for them.

## See Also

- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
containsOutOfMemoryData
showMethods("containsOutOfMemoryData")

## The default method:
selectMethod("containsOutOfMemoryData", "ANY")

## The method for list objects:
selectMethod("containsOutOfMemoryData", "list")

## The method for OutOfMemoryObject derivatives:
selectMethod("containsOutOfMemoryData", "OutOfMemoryObject")

m <- matrix(0, nrow=7, ncol=10)
m[sample(length(m), 20)] <- runif(20)
containsOutOfMemoryData(m)  # FALSE

library(SparseArray)
svt <- as(m, "SparseArray")
svt
containsOutOfMemoryData(m)  # FALSE
containsOutOfMemoryData(list(m, svt))  # FALSE

library(HDF5Array)
M <- as(m, "HDF5Array")
M
containsOutOfMemoryData(M)  # TRUE
containsOutOfMemoryData(list(m, svt, M))  # TRUE
```

---

# dbconn *Accessing SQLite DB information*

---

## Description

Get a connection object or file path for a SQLite DB

## Usage

```
dbconn(x)
dbfile(x)
```

## Arguments

| | |
|---|---|
| x | An object with a SQLite connection. |

## Value

`dbconn` returns a connection object to the SQLite DB containing `x`'s data.

`dbfile` returns a path (character string) to the SQLite DB (file) containing `x`'s data.

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- dbconn,AnnotationDb-method in the **AnnotationDbi** package for an example of a specific dbconn method (defined for dbconn objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
dbconn
showMethods("dbconn")
dbfile
showMethods("dbfile")

library(AnnotationDbi)
showMethods("dbconn")
selectMethod("dbconn", "AnnotationDb")
```

---

# density *Kernel density estimation*

---

## Description

The generic function density computes kernel density estimates.

NOTE: This man page is for the density *S4 generic function* defined in the **BiocGenerics** package. See ?stats::density for the default method (defined in the **stats** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
density(x, ...)
```

## Arguments

| | |
|---|---|
| x, ... | See ?stats::density. |

## Value

See ?stats::density for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- `stats::density` for the default `density` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- density,flowClust-method in the **flowClust** package for an example of a specific `density` method (defined for flowClust objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
density
showMethods("density")
selectMethod("density", "ANY")  # the default method
```

---

dge *Accessors and generic functions used in the context of count datasets*

---

## Description

These generic functions provide basic interfaces to operations on and data access to count datasets.

## Usage

```
counts(object, ...)
counts(object, ...) <- value
design(object, ...)
design(object, ...) <- value
dispTable(object, ...)
dispTable(object, ...)  <- value
sizeFactors(object, ...)
sizeFactors(object, ...) <- value
conditions(object, ...)
conditions(object, ...) <- value
estimateSizeFactors(object, ...)
estimateDispersions(object, ...)
plotDispEsts(object, ...)
```

## Arguments

| | |
|---|---|
| `object` | Object of class for which methods are defined, e.g., `CountDataSet`, `DESeqSummarizedExperiment` or `ExonCountSet`. |
| `value` | Value to be assigned to corresponding components of `object`; supported types depend on method implementation. |
| `...` | Further arguments, perhaps used by metohds |

## Details

For the details, please consult the manual pages of the methods in the **DESeq**, **DESeq2**, and **DEXSeq** packages and the package vignettes.

## Author(s)

W. Huber, S. Anders

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

---

# dims *Get the dimensions of each element of a list-like object*

---

## Description

Get the dimensions, number of rows, or number of columns, of each element of a list-like object.

Note that these functions are the *vectorized versions* of corresponding functions `dim()`, `nrow()`, and `ncol()`, in the same fashion that `lengths()` is the *vectorized version* of `length`.

## Usage

```
dims(x, use.names=TRUE)
nrows(x, use.names=TRUE)
ncols(x, use.names=TRUE)
```

## Arguments

| | |
|---|---|
| `x` | List-like object (or environment) where all the list elements are expected to be array-like objects with the *same number of dimensions*. |
| `use.names` | Logical indicating if the names on `x` should be propagated to the returned matrix (as its rownames) or vector (as its names). |

## Value

For `dims()`: Typically a numeric matrix with one row per list element in `x` and one column per dimension in these list elements (they're all expected to have the same number of dimensions). The i-th row in the returned matrix is a vector containing the dimensions of the i-th list element in `x`. More formally:

```
dims(x)[i, ] is dim(x[[i]])
```

for any valid i. By default the names on x, if any, are propagated as the rownames of the returned matrix, unless use.names is set to FALSE.

For nrows() or ncols(): A numeric vector with one element per list element in x. The i-th element in the returned vector is the number of rows (or columns) of the i-th list element in x. More formally:

```
nrows(x)[i] is nrow(x[[i]]) and ncols(x)[i] is ncol(x[[i]])
```

for any valid i. By default the names on x, if any, are propagated on the returned vector, unless use.names is set to FALSE.

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- dims,DataFrameList-method in the **IRanges** package for an example of a specific dims method (defined for DataFrameList objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
dims
showMethods("dims")

library(IRanges)
showMethods("dims")
selectMethod("dims", "DataFrameList")
```

---

# do.call *Execute a function call*

---

## Description

do.call constructs and executes a function call from a name or a function and a list of arguments to be passed to it.

NOTE: This man page is for the do.call *S4 generic function* defined in the **BiocGenerics** package. See ?base::do.call for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
do.call(what, args, quote=FALSE, envir=parent.frame())
```

### Arguments

| | |
|---|---|
| what | The default method expects either a function or a non-empty character string naming the function to be called. See ?base::do.call for the details.<br>Specific methods can support other objects. Please refer to the documentation of a particular method for the details. |
| args | The default method expects a *list* of arguments to the function call (the names attribute of args gives the argument names). See ?base::do.call for the details.<br>Specific methods can support other objects. Please refer to the documentation of a particular method for the details. |
| quote, envir | See ?base::do.call for a description of these arguments. |

### Value

The result of the (evaluated) function call.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

### See Also

- base::do.call for the default do.call method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

### Examples

```
do.call  # note the dispatch on the 'what' and 'args' args only
showMethods("do.call")
selectMethod("do.call", c("ANY", "ANY"))  # the default method
```

---

## duplicated *Determine duplicate elements*

---

### Description

Determines which elements of a vector-like or data-frame-like object are duplicates of elements with smaller subscripts, and returns a logical vector indicating which elements (rows) are duplicates.

NOTE: This man page is for the duplicated and anyDuplicated *S4 generic functions* defined in the **BiocGenerics** package. See ?base::duplicated for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like or data-frame-like) not supported by the default method.

## Usage

```
duplicated(x, incomparables=FALSE, ...)
anyDuplicated(x, incomparables=FALSE, ...)
```

## Arguments

| | |
|---|---|
| x | A vector-like or data-frame-like object. |
| incomparables, ... | See ?base::duplicated for a description of these arguments. |

## Value

The default duplicated method (see ?base::duplicated) returns a logical vector of length N where N is:

- length(x) when x is a vector;
- nrow(x) when x is a data frame.

Specific duplicated methods defined in Bioconductor packages must also return a logical vector of the same length as x when x is a vector-like object, and a logical vector with one element for each row when x is a data-frame-like object.

The default anyDuplicated method (see ?base::duplicated) returns a single non-negative integer and so must the specific anyDuplicated methods defined in Bioconductor packages.

anyDuplicated should always behave consistently with duplicated.

## See Also

- base::duplicated for the default duplicated and anyDuplicated methods.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- duplicated,Rle-method in the **S4Vectors** package for an example of a specific duplicated method (defined for Rle objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
duplicated
showMethods("duplicated")
selectMethod("duplicated", "ANY")  # the default method

anyDuplicated
showMethods("anyDuplicated")
selectMethod("anyDuplicated", "ANY")  # the default method
```

eval *Evaluate an (unevaluated) expression*

## Description

eval evaluates an R expression in a specified environment.

NOTE: This man page is for the eval *S4 generic function* defined in the **BiocGenerics** package. See ?base::eval for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
eval(expr, envir=parent.frame(),
           enclos=if (is.list(envir) || is.pairlist(envir))
                  parent.frame() else baseenv())
```

## Arguments

| | |
|---|---|
| expr | An object to be evaluated. May be any object supported by the default method (see ?base::eval) or by the additional methods defined in Bioconductor packages. |
| envir | The *environment* in which expr is to be evaluated. May be any object supported by the default method (see ?base::eval) or by the additional methods defined in Bioconductor packages. |
| enclos | See ?base::eval for a description of this argument. |

## Value

See ?base::eval for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- base::eval for the default eval method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- eval,expression,Vector-method in the **IRanges** package for an example of a specific eval method (defined for when the expr and envir arguments are an expression and a Vector object, respectively).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
eval  # note the dispatch on 'expr' and 'envir' args only
showMethods("eval")
selectMethod("eval", c("ANY", "ANY"))  # the default method
```

## evalq *Evaluate an (unevaluated) expression*

### Description

evalq evaluates an R expression (the quoted form of its first argument) in a specified environment.

NOTE: This man page is for the evalq wrapper defined in the **BiocGenerics** package. See ?base::evalq for the function defined in the **base** package. This wrapper correctly delegates to the eval generic, rather than base::eval.

### Usage

```
evalq(expr, envir=parent.frame(),
           enclos=if (is.list(envir) || is.pairlist(envir))
                  parent.frame() else baseenv())
```

### Arguments

| | |
|---|---|
| expr | Quoted to form the expression that is evaluated. |
| envir | The *environment* in which expr is to be evaluated. May be any object supported by methods on the eval generic. |
| enclos | See ?base::evalq for a description of this argument. |

### Value

See ?base::evalq.

### See Also

- base::evalq for the base evalq function.

### Examples

```
evalq  # note just a copy of the original evalq
```

## Extremes *Maxima and minima*

### Description

pmax, pmin, pmax.int and pmin.int return the parallel maxima and minima of the input values.

NOTE: This man page is for the pmax, pmin, pmax.int and pmin.int *S4 generic functions* defined in the **BiocGenerics** package. See ?base::pmax for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like or matrix-like) not supported by the default methods.

## Usage

```
pmax(..., na.rm=FALSE)
pmin(..., na.rm=FALSE)

pmax.int(..., na.rm=FALSE)
pmin.int(..., na.rm=FALSE)
```

## Arguments

| | |
|---|---|
| ... | One or more vector-like or matrix-like objects. |
| na.rm | See ?base::pmax for a description of this argument. |

## Value

See ?base::pmax for the value returned by the default methods.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input objects.

## See Also

- base::pmax for the default pmax, pmin, pmax.int and pmin.int methods.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- pmax,Rle-method in the **S4Vectors** package for an example of a specific pmax method (defined for Rle objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
pmax
showMethods("pmax")
selectMethod("pmax", "ANY")  # the default method

pmin
showMethods("pmin")
selectMethod("pmin", "ANY")  # the default method

pmax.int
showMethods("pmax.int")
selectMethod("pmax.int", "ANY")  # the default method

pmin.int
showMethods("pmin.int")
selectMethod("pmin.int", "ANY")  # the default method
```

## fileName *Accessing the file name of an object*

### Description

Get the file name of an object.

### Usage

```
fileName(object, ...)
```

### Arguments

| | |
|---|---|
| object | An object with a file name. |
| ... | Additional arguments, for use in specific methods. |

### See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- fileName,MSmap-method in the **MSnbase** package for an example of a specific fileName method (defined for MSmap objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

### Examples

```
fileName
showMethods("fileName")

library(MSnbase)
showMethods("fileName")
selectMethod("fileName", "MSmap")
```

## format *Format an R object for pretty printing*

### Description

Turn an R object into a character vector used for pretty printing.

NOTE: This man page is for the format *S4 generic function* defined in the **BiocGenerics** package. See ?base::format for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

### Usage

```
format(x, ...)
```

### Arguments

| | |
|---|---|
| x | The object to format. |
| ... | Additional arguments, for use in specific methods. |

### Value

A character vector that provides a "compact representation" of x. This character vector is typically used by print.data.frame to display the columns of a data.frame object. See ?base::print.data.frame for more information.

### See Also

- base::format for the default format method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

### Examples

```
format
showMethods("format")
selectMethod("format", "ANY")  # the default method
```

---

## funprog *Common higher-order functions in functional programming languages*

---

### Description

Reduce uses a binary function to successively combine the elements of a given list-like or vector-like object and a possibly given initial value. Filter extracts the elements of a list-like or vector-like object for which a predicate (logical) function gives true. Find and Position give the first or last such element and its position in the object, respectively. Map applies a function to the corresponding elements of given list-like or vector-like objects.

NOTE: This man page is for the Reduce, Filter, Find, Map and Position *S4 generic functions* defined in the **BiocGenerics** package. See ?base::Reduce for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically list-like or vector-like) not supported by the default methods.

## Usage

```
Reduce(f, x, init, right=FALSE, accumulate=FALSE, simplify=TRUE)
Filter(f, x)
Find(f, x, right=FALSE, nomatch=NULL)
Map(f, ...)
Position(f, x, right=FALSE, nomatch=NA_integer_)
```

## Arguments

| | |
|---|---|
| f, init, right, accumulate, nomatch, simplify | See ?base::Reduce for a description of these arguments. |
| x | A list-like or vector-like object. |
| ... | One or more list-like or vector-like objects. |

## Value

See ?base::Reduce for the value returned by the default methods.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default methods.

## See Also

- base::Reduce for the default Reduce, Filter, Find, Map and Position methods.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- Reduce,List-method in the **S4Vectors** package for an example of a specific Reduce method (defined for List objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
Reduce  # note the dispatch on the 'x' arg only
showMethods("Reduce")
selectMethod("Reduce", "ANY")  # the default method

Filter  # note the dispatch on the 'x' arg only
showMethods("Filter")
selectMethod("Filter", "ANY")  # the default method

Find  # note the dispatch on the 'x' arg only
showMethods("Find")
selectMethod("Find", "ANY")  # the default method

Map  # note the dispatch on the '...' arg only
showMethods("Map")
selectMethod("Map", "ANY")  # the default method

Position  # note the dispatch on the 'x' arg only
```

```
showMethods("Position")
selectMethod("Position", "ANY")  # the default method
```

---

`get` *Return the value of a named object*

---

## Description

Search for an object with a given name and return it.

NOTE: This man page is for the `get` and `mget` *S4 generic functions* defined in the **BiocGenerics** package. See `?base::get` for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (list-like or environment-like) not supported by the default methods.

## Usage

```
get(x, pos=-1, envir=as.environment(pos), mode="any",
    inherits=TRUE)

mget(x, envir=as.environment(pos), mode="any", ifnotfound,
     inherits=FALSE, pos=-1L)
```

## Arguments

| | |
|---|---|
| `x` | For `get`: A variable name (or, more generally speaking, a *key*), given as a single string.<br>For `mget`: A vector of variable names (or *keys*). |
| `envir` | Where to look for the key(s). Typically a list-like or environment-like object. |
| `pos`, `mode`, `inherits`, `ifnotfound` | See `?base::get` for a description of these arguments. |

## Details

See `?base::get` for details about the default methods.

## Value

For `get`: The value corresponding to the specified key.

For `mget`: The list of values corresponding to the specified keys. The returned list must have one element per key, and in the same order as in `x`.

See `?base::get` for the value returned by the default methods.

## See Also

- base::get for the default get and mget methods.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- get,ANY,Bimap,missing-method in the **AnnotationDbi** package for an example of a specific get method (defined for Bimap objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
get  # note the dispatch on the 'x', 'pos' and 'envir' args only
showMethods("get")
selectMethod("get", c("ANY", "ANY", "ANY"))  # the default method

mget  # note the dispatch on the 'x' and 'envir' args only
showMethods("mget")
selectMethod("mget", c("ANY", "ANY"))  # the default method
```

---

# grep *Pattern Matching and Replacement*

---

## Description

Search for matches to argument 'pattern' within each element of a character vector.

NOTE: This man page is for the grep and grepl *S4 generic functions* defined in the **BiocGenerics** package. See ?base::grep for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
grep(pattern, x, ignore.case = FALSE, perl = FALSE, value = FALSE,
          fixed = FALSE, useBytes = FALSE, invert = FALSE)
grepl(pattern, x, ignore.case = FALSE, perl = FALSE,
          fixed = FALSE, useBytes = FALSE)
```

## Arguments

| | |
|---|---|
| pattern | The pattern for searching in x, such as a regular expression. |
| x | The character vector (in the general sense) to search. |
| ignore.case, perl, value, fixed, useBytes, invert | See ?base::grep for a description of these arguments. |

## Value

See ?base::grep for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- base::grep for the default grep and grepl methods.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
grep  # note the dispatch on 'pattern' and 'x' args only
showMethods("grep")
selectMethod("grep", "ANY")  # the default method
```

---

# image *Display a color image*

---

## Description

Creates a grid of colored or gray-scale rectangles with colors corresponding to the values in z. This can be used to display three-dimensional or spatial data aka *images*.

NOTE: This man page is for the image *S4 generic function* defined in the **BiocGenerics** package. See ?graphics::image for the default method (defined in the **graphics** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
image(x, ...)
```

## Arguments

| | |
|---|---|
| x, ... | See ?graphics::image. |

## Details

See ?graphics::image for the details.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- graphics::image for the default image method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- image,AffyBatch-method in the **affy** package for an example of a specific image method (defined for AffyBatch objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
image
showMethods("image")
selectMethod("image", "ANY")  # the default method

library(affy)
showMethods("image")
## The image() method for AffyBatch objects:
selectMethod("image", "AffyBatch")
```

---

IQR *The Interquartile Range*

---

## Description

Compute the interquartile range for a vector.

NOTE: This man page is for the `IQR` *S4 generic function* defined in the **BiocGenerics** package. See `?stats::IQR` for the default method (defined in the **stats** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
IQR(x, na.rm = FALSE, type = 7)
```

## Arguments

`x`, `na.rm`, `type` See `?stats::IQR`.

## Value

See `?stats::IQR` for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- `stats::IQR` for the default `IQR` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
IQR
showMethods("IQR")
selectMethod("IQR", "ANY")  # the default method
```

`is.unsorted` *Test if a vector-like object is not sorted*

## Description

Test if a vector-like object is not sorted, without the cost of sorting it.

NOTE: This man page is for the `is.unsorted` *S4 generic function* defined in the **BiocGenerics** package. See `?base::is.unsorted` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like) not supported by the default method.

## Usage

```
is.unsorted(x, na.rm=FALSE, strictly=FALSE, ...)
```

## Arguments

| | |
|---|---|
| `x` | A vector-like object. |
| `na.rm`, `strictly` | See `?base::is.unsorted` for a description of these arguments. |
| `...` | Additional arguments, for use in specific methods.<br>Note that `base::is.unsorted` (the default method) only takes the `x`, `na.rm`, and `strictly` arguments. |

## Value

See `?base::is.unsorted` for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## Note

TO DEVELOPERS:

The `is.unsorted` method for specific vector-like objects should adhere to the same underlying order used by the `order`, `sort`, and `rank` methods for the same objects.

## See Also

- `base::is.unsorted` for the default `is.unsorted` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- is.unsorted,GenomicRanges-method in the **GenomicRanges** package for an example of a specific `is.unsorted` method (defined for GenomicRanges objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
is.unsorted  # note the dispatch on the 'x' arg only
showMethods("is.unsorted")
selectMethod("is.unsorted", "ANY")  # the default method
```

---

`lapply` *Apply a function over a list-like or vector-like object*

---

## Description

`lapply` returns a list of the same length as `X`, each element of which is the result of applying `FUN` to the corresponding element of `X`.

`sapply` is a user-friendly version and wrapper of `lapply` by default returning a vector, matrix or, if `simplify="array"`, an array if appropriate, by applying `simplify2array()`. `sapply(x, f, simplify=FALSE, USE.NAMES=FALSE)` is the same as `lapply(x, f)`.

NOTE: This man page is for the `lapply` and `sapply` *S4 generic functions* defined in the **BiocGenerics** package. See `?base::lapply` for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically list-like or vector-like) not supported by the default methods.

## Usage

```
lapply(X, FUN, ...)
sapply(X, FUN, ..., simplify=TRUE, USE.NAMES=TRUE)
```

## Arguments

| | |
|---|---|
| `X` | A list-like or vector-like object. |
| `FUN, ..., simplify, USE.NAMES` | See `?base::lapply` for a description of these arguments. |

## Value

See `?base::lapply` for the value returned by the default methods.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default methods. In particular, `lapply` and `sapply(simplify=FALSE)` should always return a list.

## See Also

- `base::lapply` for the default `lapply` and `sapply` methods.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- lapply,List-method in the **S4Vectors** package for an example of a specific `lapply` method (defined for List objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
lapply  # note the dispatch on the 'X' arg only
showMethods("lapply")
selectMethod("lapply", "ANY")  # the default method

sapply  # note the dispatch on the 'X' arg only
showMethods("sapply")
selectMethod("sapply", "ANY")  # the default method
```

---

`longForm` *Turn object into long form*

---

## Description

A generic function that returns the *long form* of an object.

## Usage

```
longForm(object, ...)
```

## Arguments

| | |
|---|---|
| `object` | A data object. |
| `...` | Additional arguments, for use in specific methods. |

## Value

The *long form* version of the original object. This is typically a data-frame-like object.

## See Also

- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- longForm,MultiAssayExperiment-method in the **MultiAssayExperiment** package for an example of a specific `longForm` method (defined for MultiAssayExperiment objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
longForm
showMethods("longForm")

library(MultiAssayExperiment)
showMethods("longForm")
selectMethod("longForm", "MultiAssayExperiment")
```

mad *Median Absolute Deviation*

## Description

Compute the median absolute deviation for a vector, dispatching only on the first argument, x.

NOTE: This man page is for the mad *S4 generic function* defined in the **BiocGenerics** package. See ?stats::mad for the default method (defined in the **stats** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
mad(x, center = median(x), constant = 1.4826,
    na.rm = FALSE, low = FALSE, high = FALSE)
```

## Arguments

x, center, constant, na.rm, low, high
: See ?stats::mad.

## Value

See ?stats::mad for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- stats::mad for the default mad method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
mad
showMethods("mad")
selectMethod("mad", "ANY")  # the default method
```

mapply *Apply a function to multiple list-like or vector-like arguments*

## Description

mapply is a multivariate version of sapply. mapply applies FUN to the first elements of each ... argument, the second elements, the third elements, and so on. Arguments are recycled if necessary.

NOTE: This man page is for the mapply *S4 generic function* defined in the **BiocGenerics** package. See ?base::mapply for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically list-like or vector-like) not supported by the default methods.

## Usage

```
mapply(FUN, ..., MoreArgs=NULL, SIMPLIFY=TRUE, USE.NAMES=TRUE)
```

## Arguments

| | |
|---|---|
| FUN, MoreArgs, SIMPLIFY, USE.NAMES | See ?base::mapply for a description of these arguments. |
| ... | One or more list-like or vector-like objects of strictly positive length, or all of zero length. |

## Value

See ?base::mapply for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- base::mapply for the default mapply method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
mapply  # note the dispatch on the '...' arg only
showMethods("mapply")
selectMethod("mapply", "ANY")  # the default method
```

`match` *Value matching*

## Description

`match` returns a vector of the positions of (first) matches of its first argument in its second.

`%in%` is a binary operator that returns a logical vector of the length of its left operand indicating if the elements in it have a match or not.

NOTE: This man page is for the `match` and `%in%` *S4 generic functions* defined in the **BiocGenerics** package. See `?base::match` for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like) not supported by the default methods.

## Usage

```
match(x, table, nomatch=NA_integer_, incomparables=NULL, ...)

x %in% table
```

## Arguments

| | |
|---|---|
| `x`, `table` | Vector-like objects (typically of the same class, but not necessarily). |
| `nomatch`, `incomparables` | See `?base::match` for a description of these arguments. |
| `...` | Additional arguments, for use in specific methods. |

## Value

The same as the default methods (see `?base::match` for the value returned by the default methods).

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default methods.

## Note

The default `base::match` method (defined in the **base** package) doesn't have the `...` argument. We've added it to the generic function defined in the **BiocGenerics** package in order to allow specific methods to support additional arguments if needed.

## See Also

- `base::match` for the default `match` and `%in%` methods.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- match,Hits,Hits-method and %in%,Rle,ANY-method in the **S4Vectors** package for examples of specific `match` and `%in%` methods (defined for Hits and Rle objects, respectively).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

### Examples

```
match  # note the dispatch on the 'x' and 'table' args only
showMethods("match")
selectMethod("match", c("ANY", "ANY"))  # the default method

`%in%`
showMethods("%in%")
selectMethod("%in%", c("ANY", "ANY"))  # the default method
```

---

## mean *Arithmetic Mean*

---

### Description

Generic function for the (trimmed) arithmetic mean.

NOTE: This man page is for the `mean` *S4 generic function* defined in the **BiocGenerics** package. See `?base::mean` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like) not supported by the default method.

### Usage

```
mean(x, ...)
```

### Arguments

| | |
|---|---|
| `x` | typically a vector-like object |
| `...` | see `mean` |

### Value

See `?base::mean` for the value returned by the default method.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input object.

### See Also

- `base::mean` for the default `mean` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- mean,Rle-method in the **S4Vectors** package for an example of a specific `mean` method (defined for Rle objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

### Examples

```
mean
showMethods("mean")
selectMethod("mean", "ANY")  # the default method
```

`normalize` *Normalize an object*

## Description

A generic function which normalizes an object containing microarray data or other data. Normalization is intended to remove from the intensity measures any systematic trends which arise from the microarray technology rather than from differences between the probes or between the target RNA samples hybridized to the arrays.

## Usage

```
normalize(object, ...)
```

## Arguments

| | |
|---|---|
| `object` | A data object, typically containing microarray data. |
| `...` | Additional arguments, for use in specific methods. |

## Value

An object containing the normalized data.

## See Also

- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- normalize,AffyBatch-method in the **affy** package and normalize,MSnExp-method in the **MSnbase** package for examples of specific `normalize` methods (defined for AffyBatch and MSnExp objects, respectively).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
normalize
showMethods("normalize")

library(affy)
showMethods("normalize")
selectMethod("normalize", "AffyBatch")
```

nrow *The number of rows/columns of an array-like object*

## Description

Return the number of rows or columns present in an array-like object.

NOTE: This man page is for the `nrow`, `ncol`, `NROW` and `NCOL` *S4 generic functions* defined in the **BiocGenerics** package. See `?base::nrow` for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically matrix- or array-like) not supported by the default methods.

## Usage

```
nrow(x)
ncol(x)
NROW(x)
NCOL(x)
```

## Arguments

| | |
|---|---|
| x | A matrix- or array-like object. |

## Value

A single integer or `NULL`.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default methods.

## See Also

- `base::nrow` for the default `nrow`, `ncol`, `NROW` and `NCOL` methods.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- nrow,DataFrame-method in the **S4Vectors** package for an example of a specific `nrow` method (defined for DataFrame objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
nrow
showMethods("nrow")
selectMethod("nrow", "ANY")  # the default method

ncol
showMethods("ncol")
selectMethod("ncol", "ANY")  # the default method
```

```
NROW
showMethods("NROW")
selectMethod("NROW", "ANY")  # the default method

NCOL
showMethods("NCOL")
selectMethod("NCOL", "ANY")  # the default method
```

---

`Ontology` *Generic Ontology getter*

---

## Description

Get the Ontology of an object.

## Usage

```
Ontology(object)
```

## Arguments

| | |
|---|---|
| `object` | An object with an Ontology. |

## See Also

- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- Ontology,GOTerms-method in the **AnnotationDbi** package for an example of a specific `Ontology` method (defined for GOTerms objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
Ontology
showMethods("Ontology")

library(AnnotationDbi)
showMethods("Ontology")
selectMethod("Ontology", "GOTerms")
```

order *Ordering permutation*

## Description

order returns a permutation which rearranges its first argument into ascending or descending order, breaking ties by further arguments.

NOTE: This man page is for the order *S4 generic function* defined in the **BiocGenerics** package. See ?base::order for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like) not supported by the default method.

## Usage

```
order(..., na.last=TRUE, decreasing=FALSE, method=c("auto", "shell", "radix"))
```

## Arguments

| | |
|---|---|
| ... | One or more vector-like objects, all of the same length. |
| na.last, decreasing, method | See ?base::order for a description of these arguments. |

## Value

The default method (see ?base::order) returns an integer vector of length N where N is the common length of the input objects. This integer vector represents a permutation of N elements and can be used to rearrange the first argument in ... into ascending or descending order (by subsetting it).

Specific methods defined in Bioconductor packages should also return an integer vector representing a permutation of N elements.

## Note

TO DEVELOPERS:

Specific order methods should preferably be made "stable" for consistent behavior across platforms and consistency with base::order(). Note that C qsort() is *not* "stable" so order methods that use qsort() at the C-level need to ultimately break ties by position, which can easily be done by adding a little extra code at the end of the comparison function passed to qsort().

order(x, decreasing=TRUE) is *not* always equivalent to rev(order(x)).

order, sort, and rank methods for specific vector-like objects should adhere to the same underlying order that should be conceptually defined as a binary relation on the set of all possible vector values. For completeness, this binary relation should also be incarnated by a <= method.

## See Also

- base::order for the default order method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- order,IntegerRanges-method in the **IRanges** package for an example of a specific order method (defined for IntegerRanges objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
order
showMethods("order")
selectMethod("order", "ANY")  # the default method
```

---

# organism_species *Organism and species accessors*

---

## Description

Get or set the organism and/or species of an object.

## Usage

```
organism(object)
organism(object) <- value

species(object)
species(object) <- value
```

## Arguments

| | |
|---|---|
| object | An object to get or set the organism or species of. |
| value | The organism or species to set on object. |

## Value

organism should return the *scientific name* (i.e. genus and species, or genus and species and subspecies) of the organism. Preferably in the format "Genus species" (e.g. "Homo sapiens") or "Genus species subspecies" (e.g. "Homo sapiens neanderthalensis").

species should of course return the species of the organism. Unfortunately there is a long history of misuse of this accessor in Bioconductor so its usage is now discouraged (starting with BioC 3.1).

## Note

TO DEVELOPERS:

`species` has been historically misused in many places in Bioconductor and is redundant with `organism`. So implementing the `species` accessor is now discouraged (starting with BioC 3.1). The `organism` accessor (returning the *scientific name*) should be implemented instead.

## See Also

- http://bioconductor.org/packages/release/BiocViews.html#___Organism for browsing the annotation packages currently available in Bioconductor by organism.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- organism,character-method and organism,chromLocation-method in the **annotate** package for examples of specific `organism` methods (defined for character and chromLocation objects).
- species,AnnotationDb-method in the **AnnotationDbi** package for an example of a specific `species` method (defined for AnnotationDb objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
## organism() getter:
organism
showMethods("organism")

library(annotate)
showMethods("organism")
selectMethod("organism", "character")
selectMethod("organism", "chromLocation")

## organism() setter:
`organism<-`
showMethods("organism<-")

## species() getter:
species
showMethods("species")

library(AnnotationDbi)
selectMethod("species", "AnnotationDb")

## species() setter:
`species<-`
showMethods("species<-")
```

paste *Concatenate strings*

## Description

paste concatenates vectors of strings or vector-like objects containing strings.

NOTE: This man page is for the paste *S4 generic function* defined in the **BiocGenerics** package. See ?base::paste for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like objects containing strings) not supported by the default method.

## Usage

```
paste(..., sep=" ", collapse=NULL, recycle0=FALSE)
```

## Arguments

| | |
|---|---|
| ... | One or more vector-like objects containing strings. |
| sep, collapse, recycle0 | See ?base::paste for a description of these arguments. |

## Value

See ?base::paste for the value returned by the default method.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input objects.

## See Also

- base::paste for the default paste method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- paste,Rle-method in the **S4Vectors** package for an example of a specific paste method (defined for Rle objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
paste
showMethods("paste")
selectMethod("paste", "ANY")  # the default method
```

paste2 *Concatenate strings (binary form)*

## Description

paste2() is a simplified version of paste0() that takes only two arguments and behaves like the arithmetic binary operations (+, *, etc...) in terms of:

1. recycling;
2. propagation of NAs;
3. propagation of names;
4. propagation of dimensions and dimnames.

add_prefix() and add_suffix() are simple wrappers around paste2() provided for convenience and code readability.

## Usage

```
paste2(x, y)

add_prefix(x, prefix="")
add_suffix(x, suffix="")
```

## Arguments

x, y, prefix, suffix
: Vector- or array-like objects containing strings.

## Details

Unlike paste0(), paste2() only takes two arguments: x and y. It's defined as an S4 generic that dispatches on its two arguments and with methods for ordinary vectors and arrays. Bioconductor packages can define methods for other vector-like or array-like objects that contain strings.

paste2() follows the same rules as arithmetic operations (+, *, etc...) with regard to:

1. Recycling: The longer argument "wins" i.e. the shorter argument is recycled to the length of the longer (with a warning if the length of the latter is not a multiple of the length of the former). There's one important exception to this rule: if one of the two arguments has length 0, then no recycling is performed and a zero-length vector is returned.
2. Propagation of NAs.
3. Propagation of names: The longer argument also wins. If the two arguments have the same length, then:
   - the names on the first argument are propagated, if any;
   - otherwise the names on the second argument are propagated, if any.

4. Propagation of dimensions and dimnames: If `x` and `y` are both arrays, then they must be *conformable* i.e. have the same dimensions. In this case the result of `paste2(x, y)` is also an array of same dimensions. Furthermore it will have `dimnames(x)` on it if `dimnames(x)` is not `NULL`, otherwise it will have `dimnames(y)` on it.

`add_prefix(x, prefix="")` and `add_suffix(x, suffix="")` are convenience wrappers that just do `paste2(prefix, x)` and `paste2(x, suffix)`, respectively.

## Value

If `x` and `y` are both vectors, a character vector *parallel* to the longer vector is returned.

If one of `x` or `y` is an array and the other one a vector, an array *parallel* to the input array is returned.

If `x` and `y` are both arrays (in which case they must be *conformable*), an array *parallel* to `x` and `y` is returned.

## See Also

- `base::paste0` in base R.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- paste2,DelayedArray,DelayedArray-method in the **DelayedArray** package for an example of a specific `paste2` method (defined for DelayedArray objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
## ---------------------------------------------------------------------
## The paste2() generic and methods
## ---------------------------------------------------------------------

paste2  # note the dispatch on 'x' and 'y'
showMethods("paste2")

## ---------------------------------------------------------------------
## paste0() vs paste2()
## ---------------------------------------------------------------------

## Propagation of NAs and names:
x <- c(A="foo", B="bar", C=NA)
paste0(x, "XX")  # NAs are not propagated, names are lost
paste2(x, "XX")  # NAs and names are propagated
paste2(x, setNames(1:6, letters[1:6]))  # longer argument "wins"

## If 'x' or 'y' has length 0:
paste0(x, character(0))  # unname(x)
paste2(x, character(0))  # character(0)

## Propagation of dimensions and dimnames:
m <- matrix(1:12, ncol=3, dimnames=list(NULL, LETTERS[1:3]))
y <- c("a", NA, "c")
```

```
paste0(m, y)  # dimensions and dimnames are lost
paste2(m, y)  # dimensions are preserved and dimnames are propagated

## ------------------------------------------------------------------------
## add_prefix() and add_suffix()
## ------------------------------------------------------------------------

m2 <- add_prefix(m, "ID")  # same as paste2("ID", m)
add_suffix(m2, ".fasta")   # same as paste2(m2, ".fasta")
```

---

path *Accessing the path of an object*

---

## Description

Get or set the path of an object.

## Usage

```
path(object, ...)
path(object, ...) <- value

basename(path, ...)
basename(path, ...) <- value

dirname(path, ...)
dirname(path, ...) <- value

## The purpose of the following methods is to make the basename() and
## dirname() getters work out-of-the-box on any object for which the
## path() getter works.

## S4 method for signature 'ANY'
basename(path, ...)

## S4 method for signature 'ANY'
dirname(path, ...)

## The purpose of the following replacement methods is to make the
## basename() and dirname() setters work out-of-the-box on any object
## for which the path() getter and setter work.

## S4 replacement method for signature 'character'
basename(path, ...) <- value

## S4 replacement method for signature 'ANY'
basename(path, ...) <- value
```

```
## S4 replacement method for signature 'character'
dirname(path, ...) <- value

## S4 replacement method for signature 'ANY'
dirname(path, ...) <- value
```

## Arguments

| | |
|---|---|
| object | An object containing paths. Even though it will typically contain a single path, object can actually contain an arbitrary number of paths. |
| ... | Additional arguments, for use in specific methods. |
| value | For path<-, the paths to set on object.<br>For basename<- or dirname<-, the basenames or dirnames to set on path. |
| path | A character vector *or an object containing paths*. |

## Value

A character vector for path(object), basename(path), and dirname(path). Typically of length 1 but not necessarily. Possibly with names on it for path(object).

## See Also

- base::basename for the functions the basename and dirname generics are based on.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- path,RsamtoolsFile-method in the **Rsamtools** package for an example of a specific path method (defined for RsamtoolsFile objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
## ---------------------------------------------------------------------
## GENERIC FUNCTIONS AND DEFAULT METHODS
## ---------------------------------------------------------------------

path
showMethods("path")

`path<-`
showMethods("path<-")

basename
showMethods("basename")

`basename<-`
showMethods("basename<-")

dirname
```

```
showMethods("dirname")

`dirname`
showMethods("dirname<-")


## Default basename() and dirname() getters:
selectMethod("basename", "ANY")
selectMethod("dirname", "ANY")

## Default basename() and dirname() setters:
selectMethod("basename<-", "character")
selectMethod("basename<-", "ANY")
selectMethod("dirname<-", "character")
selectMethod("dirname<-", "ANY")

## ---------------------------------------------------------------------
## OBJECTS CONTAINING PATHS
## ---------------------------------------------------------------------

## Let's define a simple class to represent objects that contain paths:
setClass("A", slots=c(stuff="ANY", path="character"))

a <- new("A", stuff=runif(5),
              path=c(one="path/to/file1", two="path/to/file2"))

## path() getter:
setMethod("path", "A", function(object) object@path)

path(a)

## Because the path() getter works on 'a', now the basename() and
## dirname() getters also work:
basename(a)
dirname(a)

## path() setter:
setReplaceMethod("path", "A",
    function(object, ..., value)
    {
        if (length(list(...)) != 0L) {
            dots <- match.call(expand.dots=FALSE)[[3L]]
            stop(BiocGenerics:::unused_arguments_msg(dots))
        }
        object@path <- value
        object
    }
)

a <- new("A", stuff=runif(5))
path(a) <- c(one="path/to/file1", two="path/to/file2")
path(a)
```

```
## Because the path() getter and setter work on 'a', now the basename()
## and dirname() setters also work:
basename(a) <- toupper(basename(a))
path(a)
dirname(a) <- "~/MyDataFiles"
path(a)
```

---

plotMA *MA-plot: plot differences versus averages for high-throughput data*

---

## Description

A generic function which produces an MA-plot for an object containing microarray, RNA-Seq or other data.

## Usage

```
plotMA(object, ...)
```

## Arguments

| | |
|---|---|
| object | A data object, typically containing count values from an RNA-Seq experiment or microarray intensity values. |
| ... | Additional arguments, for use in specific methods. |

## Value

Undefined. The function exists for its side effect, producing a plot.

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- plotMA in the **limma** package for a function with the same name that is not dispatched through this generic function.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
showMethods("plotMA")

suppressWarnings(
  if(require("DESeq2"))
    example("plotMA", package="DESeq2", local=TRUE)
)
```

plotPCA *PCA-plot: Principal Component Analysis plot*

## Description

A generic function which produces a PCA-plot.

## Usage

```
plotPCA(object, ...)
```

## Arguments

| | |
|---|---|
| object | A data object, typically containing gene expression information. |
| ... | Additional arguments, for use in specific methods. |

## Value

Undefined. The function exists for its side effect, producing a plot.

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- plotPCA in the **DESeq2** package for an example method that uses this generic.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
showMethods("plotPCA")

suppressWarnings(
  if(require("DESeq2"))
    example("plotPCA", package="DESeq2", local=TRUE)
)
```

rank *Ranks the values in a vector-like object*

## Description

Returns the ranks of the values in a vector-like object. Ties (i.e., equal values) and missing values can be handled in several ways.

NOTE: This man page is for the rank *S4 generic function* defined in the **BiocGenerics** package. See ?base::rank for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
rank(x, na.last=TRUE,
     ties.method=c("average", "first", "last", "random", "max", "min"),
     ...)
```

## Arguments

| | |
|---|---|
| x | A vector-like object. |
| na.last, ties.method | See ?base::rank for a description of these arguments. |
| ... | Additional arguments, for use in specific methods.<br>Note that base::rank (the default method) only takes the x, na.last, and ties.method arguments. |

## Value

See ?base::rank for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## Note

TO DEVELOPERS:

See note in ?BiocGenerics::order about "stable" order.

order, sort, and rank methods for specific vector-like objects should adhere to the same underlying order that should be conceptually defined as a binary relation on the set of all possible vector values. For completeness, this binary relation should also be incarnated by a <= method.

## See Also

- base::rank for the default rank method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.

- rank,Vector-method in the **S4Vectors** package for an example of a specific `rank` method (defined for Vector objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
rank  # note the dispatch on the 'x' arg only
showMethods("rank")
selectMethod("rank", "ANY")  # the default method
```

---

`relist` *Re-listing an unlist()ed object*

---

## Description

`relist` is a generic function with a few methods in order to allow easy inversion of `unlist(x)`.

NOTE: This man page is for the `relist` *S4 generic function* defined in the **BiocGenerics** package. See `?utils::relist` for the default method (defined in the **utils** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
relist(flesh, skeleton)
```

## Arguments

| | |
|---|---|
| `flesh` | A vector-like object. |
| `skeleton` | A list-like object. Only the "shape" (i.e. the lengths of the individual list elements) of `skeleton` matters. Its exact content is ignored. |

## Value

A list-like object with the same "shape" as `skeleton` and that would give `flesh` back if unlist()ed.

## See Also

- `utils::relist` for the default `relist` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- relist,ANY,List-method in the **IRanges** package for an example of a specific `relist` method (defined for when `skeleton` is a List object).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
relist
showMethods("relist")
selectMethod("relist", c("ANY", "ANY"))  # the default method
```

rep *Replicate elements of a vector-like object*

## Description

rep.int replicates the elements in x.

NOTE: This man page is for the rep.int *S4 generic function* defined in the **BiocGenerics** package. See ?base::rep.int for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like) not supported by the default method.

## Usage

```
rep.int(x, times)
```

## Arguments

| | |
|---|---|
| x | The object to replicate (typically vector-like). |
| times | See ?base::rep.int for a description of this argument. |

## Value

See ?base::rep.int for the value returned by the default method.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input object.

## See Also

- base::rep.int for the default rep.int method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- rep.int,Rle-method in the **S4Vectors** package for an example of a specific rep.int method (defined for Rle objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
rep.int
showMethods("rep.int")
selectMethod("rep.int", "ANY")  # the default method
```

`residuals` *Extract model residuals*

## Description

`residuals` is a generic function which extracts model residuals from objects returned by modeling functions.

NOTE: This man page is for the `residuals` *S4 generic function* defined in the **BiocGenerics** package. See `?stats::residuals` for the default method (defined in the **stats** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
residuals(object, ...)
```

## Arguments

| | |
|---|---|
| `object, ...` | See `?stats::residuals`. |

## Value

Residuals extracted from the object `object`.

## See Also

- `stats::residuals` for the default `residuals` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- residuals,PLMset-method in the **affyPLM** package for an example of a specific `residuals` method (defined for PLMset objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
residuals
showMethods("residuals")
selectMethod("residuals", "ANY")  # the default method
```

`row+colnames` *Row and column names*

## Description

Get or set the row or column names of a matrix-like object.

NOTE: This man page is for the `rownames`, `` `rownames<-` ``, `colnames`, and `` `colnames<-` `` *S4 generic functions* defined in the **BiocGenerics** package. See `?base::rownames` for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically matrix-like) not supported by the default methods.

## Usage

```
rownames(x, do.NULL=TRUE, prefix="row")
rownames(x) <- value

colnames(x, do.NULL=TRUE, prefix="col")
colnames(x) <- value
```

## Arguments

| | |
|---|---|
| `x` | A matrix-like object. |
| `do.NULL, prefix` | See `?base::rownames` for a description of these arguments. |
| `value` | Either `NULL` or a character vector equal of length equal to the appropriate dimension. |

## Value

The getters will return `NULL` or a character vector of length `nrow(x)` for `rownames` and length `ncol(x)` for `colnames(x)`.

See `?base::rownames` for more information about the default methods, including how the setters are expected to behave.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default methods.

## See Also

- `base::rownames` for the default `rownames`, `` `rownames<-` ``, `colnames`, and `` `colnames<-` `` methods.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- rownames,DataFrame-method in the **S4Vectors** package for an example of a specific `rownames` method (defined for DataFrame objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
## rownames() getter:
rownames  # note the dispatch on the 'x' arg only
showMethods("rownames")
selectMethod("rownames", "ANY")  # the default method

## rownames() setter:
`rownames<-`
showMethods("rownames<-")
selectMethod("rownames<-", "ANY")  # the default method

## colnames() getter:
colnames  # note the dispatch on the 'x' arg only
showMethods("colnames")
selectMethod("colnames", "ANY")  # the default method

## colnames() setter:
`colnames<-`
showMethods("colnames<-")
selectMethod("colnames<-", "ANY")  # the default method
```

---

`S3-classes-as-S4-classes`

*S3 classes as S4 classes*

---

## Description

Some old-style (aka S3) classes are turned into formally defined (aka S4) classes by the **BiocGenerics** package. This allows S4 methods defined in Bioconductor packages to use them in their signatures.

## Details

S3 classes currently turned into S4 classes:

- connection class and subclasses: connection, file, url, gzfile, bzfile, unz, pipe, fifo, sockconn, terminal, textConnection, gzcon. Addtitionally the character_OR_connection S4 class is defined as the union of classes character and connection.
- others: AsIs, dist

## See Also

setOldClass and setClassUnion in the **methods** package.

saveRDS *The saveRDS() S4 generic and default method*

## Description

Generic function to write a single R object to a file.

NOTE: This man page is for the saveRDS *S4 generic function* and default method defined in the **BiocGenerics** package. See ?base::saveRDS for the corresponding function defined in base R.

## Usage

```
saveRDS(object, file="", ascii=FALSE, version=NULL,
        compress=TRUE, refhook=NULL)
```

## Arguments

object, file, ascii, version, compress, refhook
: See ?base::saveRDS for a description of these arguments.

## Details

The default saveRDS method defined in this package is a thin wrapper around base::saveRDS that issues a warning if the object to serialize contains out-of-memory data. See ?containsOutOfMemoryData for more information.

Bioconductor packages can override this default method with more specialized methods.

## Value

An invisible NULL.

## See Also

- base::saveRDS in the **base** package for the default saveRDS method.
- containsOutOfMemoryData for determining whether an object contains out-of-memory data or not.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- saveRDS,SummarizedExperiment-method in the **SummarizedExperiment** package for an example of a specific saveRDS method (defined for SummarizedExperiment objects and derivatives).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
saveRDS  # note the dispatch on the 'object' arg only
showMethods("saveRDS")
selectMethod("saveRDS", "ANY")  # the default method
```

score *Score accessor*

## Description

Get or set the score value contained in an object.

## Usage

```
score(x, ...)
score(x, ...) <- value
```

## Arguments

| | |
|---|---|
| x | An object to get or set the score value of. |
| ... | Additional arguments, for use in specific methods. |
| value | The score value to set on x. |

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- score,GenomicRanges-method in the **GenomicRanges** package for an example of a specific score method (defined for GenomicRanges objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
score
showMethods("score")

`score<-`
showMethods("score<-")

library(GenomicRanges)

showMethods("score")
selectMethod("score", "GenomicRanges")

showMethods("score<-")
selectMethod("score<-", "GenomicRanges")
```

setops *Set operations*

## Description

Performs *set* union, intersection, (asymmetric!) difference, and equality on two or more vector-like objects.

NOTE: This man page is for the `union`, `intersect`, `setdiff`, and `setequal` *S4 generic functions* defined in the **BiocGenerics** package. See `?generics::union` for the default methods (defined in CRAN package **generics**). Bioconductor packages can define specific methods for objects (typically vector-like) not supported by the default methods.

## Usage

```
union(x, y, ...)
intersect(x, y, ...)
setdiff(x, y, ...)
setequal(x, y, ...)
```

## Arguments

| | |
|---|---|
| x, y | Vector-like objects (typically of the same class, but not necessarily). |
| ... | Additional arguments, for use in specific methods. |

## Value

See `?generics::union` in CRAN package **generics** for the value returned by the default methods.

Specific methods defined in Bioconductor packages will typically act as *endomorphisms*, that is, they'll return an object of the same class as the input objects.

## Note

The default S4 methods for these S4 generics are the `union`, `intersect`, `setdiff`, and `setequal` functions defined in CRAN package **generics**, which are themselves *S3 generic functions*. These S3 generics in turn have default methods that simply call the corresponding base R function i.e. the `union`, `intersect`, `setdiff`, or `setequal` function defined in the **base** package. See for example `generics:::union.default`.

Note that the base R functions only take 2 arguments. However, the S3 generics in CRAN package **generics**, and the S4 generics in **BiocGenerics**, add `...` (a.k.a. ellipsis) to the argument list. This allow these generics to be called with an arbitrary number of effective arguments.

For `union` or `intersect`, this means that Bioconductor packages can implement *N-ary* union or intersection operations, that is, methods that compute the union or intersection of more than 2 objects.

However, for `setdiff` and `setequal`, which are conceptually binary-only operation, the presence of the ellipsis typically allows methods to support extra arguments to control/alter the behavior of

the operation. Like for example the `ignore.strand` argument supported by the `setdiff` method for GenomicRanges objects (defined in the **GenomicRanges** package). (Note that the `union` and `intersect` methods for those objects also support the `ignore.strand` argument.)

## See Also

- `generics::union` for the default `union`, `intersect`, `setdiff`, and `setequal` S4 methods defined in CRAN package **generics**.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- union,GenomicRanges,GenomicRanges-method in the **GenomicRanges** package for examples of specific `union`, `intersect`, and `setdiff` methods (defined for GenomicRanges objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
## ---------------------------------------------------------------------
## union()
## ---------------------------------------------------------------------

## S4 generic:
union  # note the dispath on 'x' and 'y'

showMethods("union")

## The default S4 method is an S3 generic function defined in
## CRAN package generics:
selectMethod("union", c("ANY", "ANY"))

## The default S3 method just calls base::union():
generics:::union.default

## ---------------------------------------------------------------------
## intersect()
## ---------------------------------------------------------------------

## S4 generic:
intersect  # note the dispath on 'x' and 'y'

showMethods("intersect")

## The default S4 method is an S3 generic function defined in
## CRAN package generics:
selectMethod("intersect", c("ANY", "ANY"))

## The default S3 method just calls base::intersect():
generics:::intersect.default

## ---------------------------------------------------------------------
## setdiff()
```

```
## ------------------------------------------------------------------------

## S4 generic:
setdiff  # note the dispath on 'x' and 'y'

showMethods("setdiff")

## The default S4 method is an S3 generic function defined in
## CRAN package generics:
selectMethod("setdiff", c("ANY", "ANY"))

## The default S3 method just calls base::setdiff():
generics:::setdiff.default

## ------------------------------------------------------------------------
## setequal()
## ------------------------------------------------------------------------

## S4 generic:
setequal  # note the dispath on 'x' and 'y'

showMethods("setequal")

## The default S4 method is an S3 generic function defined in
## CRAN package generics:
selectMethod("setequal", c("ANY", "ANY"))

## The default S3 method just calls base::setequal():
generics:::setequal.default
```

---

`sort` *Sorting a vector-like object*

---

## Description

Sort a vector-like object into ascending or descending order.

NOTE: This man page is for the `sort` *S4 generic function* defined in the **BiocGenerics** package. See `?base::sort` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
sort(x, decreasing=FALSE, ...)
```

## Arguments

| | |
|---|---|
| `x` | A vector-like object. |
| `decreasing, ...` | See `?base::sort` for a description of these arguments. |

## Value

See ?base::sort for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## Note

TO DEVELOPERS:

See note in ?BiocGenerics::order about "stable" order.

order, sort, and rank methods for specific vector-like objects should adhere to the same underlying order that should be conceptually defined as a binary relation on the set of all possible vector values. For completeness, this binary relation should also be incarnated by a <= method.

## See Also

- base::sort for the default sort method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- sort,Vector-method in the **S4Vectors** package for an example of a specific sort method (defined for Vector objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
sort  # note the dispatch on the 'x' arg only
showMethods("sort")
selectMethod("sort", "ANY")  # the default method
```

---

# start *The start(), end(), width(), and pos() generic getters and setters*

---

## Description

Get or set the start, end, width, or single positions stored in an object.

NOTE: This man page is for the start, `start<-`, end, `end<-`, width, `width<-`, and pos *S4 generic functions* defined in the **BiocGenerics** package. See ?stats::start for the start and end S3 generics defined in the **stats** package.

## Usage

```
start(x, ...)
start(x, ...) <- value

end(x, ...)
end(x, ...) <- value

width(x)
width(x, ...) <- value

pos(x)
```

## Arguments

| | |
|---|---|
| x | For the start(), end(), and width() getters/setters: an object containing start, end, and width values.<br>For the pos{} getter: an object containing single positions. |
| ... | Additional arguments, for use in specific methods. |
| value | The start, end, or width values to set on x. |

## Value

See specific methods defined in Bioconductor packages.

## See Also

- stats::start in the **stats** package for the start and end S3 generics.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- start,IRanges-method in the **IRanges** package for examples of specific start, end, and width methods (defined for IRanges objects).
- pos,UnstitchedIPos-method in the **IRanges** package for an example of a specific pos method (defined for UnstitchedIPos objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
## start() getter:
start
showMethods("start")

library(IRanges)
showMethods("start")
selectMethod("start", "IRanges")  # start() getter for IRanges objects

## start() setter:
`start<-`
```

```
showMethods("start<-")
selectMethod("start<-", "IRanges")  # start() setter for IRanges objects

## end() getter:
end
showMethods("end")
selectMethod("end", "IRanges")  # end() getter for IRanges objects

## end() setter:
`end<-`
showMethods("end<-")
selectMethod("end<-", "IRanges")  # end() setter for IRanges objects

## width() getter:
width
showMethods("width")
selectMethod("width", "IRanges")  # width() getter for IRanges objects

## width() setter:
`width<-`
showMethods("width<-")
selectMethod("width<-", "IRanges")  # width() setter for IRanges objects

## pos() getter:
pos
showMethods("pos")
selectMethod("pos", "UnstitchedIPos")  # pos() getter for UnstitchedIPos
                                       # objects
```

---

strand *Accessing strand information*

---

## Description

Get or set the strand information contained in an object.

## Usage

```
strand(x, ...)
strand(x, ...) <- value

unstrand(x)

invertStrand(x)
## S4 method for signature 'ANY'
invertStrand(x)
```

## Arguments

| | |
|---|---|
| x | An object containing strand information. |
| ... | Additional arguments, for use in specific methods. |
| value | The strand information to set on x. |

## Details

All the strand methods defined in the **GenomicRanges** package use the same set of 3 values (called the "standard strand levels") to specify the strand of a genomic location: +, -, and *. * is used when the exact strand of the location is unknown, or irrelevant, or when the "feature" at that location belongs to both strands.

Note that unstrand is not a generic function, just a convenience wrapper to the generic strand() setter (strand<-) that does:

```
strand(x) <- "*"
x
```

The default method for invertStrand does:

```
strand(x) <- invertStrand(strand(x))
x
```

## Value

If x is a vector-like object, strand(x) will typically return a vector-like object *parallel* to x, that is, an object of the same length as x where the i-th element describes the strand of the i-th element in x.

unstrand(x) and invertStrand(x) return a copy of x with the strand set to "*" for unstrand or inverted for invertStrand (i.e. "+" and "-" switched, and "*" untouched).

## See Also

- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- strand,GRanges-method in the **GenomicRanges** package for an example of a specific strand method (defined for GRanges objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
strand
showMethods("strand")

`strand<-`
showMethods("strand<-")

unstrand
```

```
invertStrand
showMethods("invertStrand")
selectMethod("invertStrand", "ANY")  # the default method

library(GenomicRanges)

showMethods("strand")
selectMethod("strand", "missing")
strand()

showMethods("strand<-")
```

---

`subset` *Subsetting vector-like, matrix-like and data-frame-like objects*

---

## Description

Return subsets of vector-like, matrix-like or data-frame-like objects which meet conditions.

NOTE: This man page is for the `subset` *S4 generic function* defined in the **BiocGenerics** package. See `?base::subset` for the `subset` S3 generic defined in the **base** package.

## Usage

```
subset(x, ...)
```

## Arguments

| | |
|---|---|
| `x` | A vector-like, matrix-like or data-frame-like object to be subsetted. |
| `...` | Additional arguments (e.g. `subset`, `select`, `drop`), for use in specific methods. See `?base::subset` for more information. |

## Value

An object similar to `x` containing just the selected elements (for a vector-like object), or the selected rows and columns (for a matrix-like or data-frame-like object).

## See Also

- `base::subset` in the **base** package for the `subset` S3 generic.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- subset,RectangularData-method in the **S4Vectors** package for an example of a specific `subset` method (defined for RectangularData derivatives).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
subset
showMethods("subset")
selectMethod("subset", "ANY")  # the default method

library(S4Vectors)
showMethods("subset")
## The subset() method for RectangularData derivatives:
selectMethod("subset", "RectangularData")
```

---

t *Matrix Transpose*

---

## Description

Given a rectangular object `x`, `t` returns the transpose of `x`.

NOTE: This man page is for the `t` *S4 generic function* defined in the **BiocGenerics** package. See `?base::t` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically array-like) not supported by the default method.

## Usage

```
t(x)
```

## Arguments

| | |
|---|---|
| `x` | A matrix-like or other rectangular object. |

## Value

See `?base::t` for the value returned by the default method.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input object.

## See Also

- `base::t` for the default `t` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- t,Hits-method in the **S4Vectors** package for an example of a specific `t` method (defined for Hits objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
t
showMethods("t")
selectMethod("t", "ANY")  # the default method
```

`table` *Cross tabulation and table creation*

## Description

`table` uses the cross-classifying factors to build a contingency table of the counts at each combination of factor levels.

NOTE: This man page is for the `table` *S4 generic function* defined in the **BiocGenerics** package. See `?base::table` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
table(...)
```

## Arguments

| | |
|---|---|
| `...` | One or more objects which can be interpreted as factors (including character strings), or a list (or data frame) whose components can be so interpreted. |

## Value

See `?base::table` for the value returned by the default method.

Specific methods defined in Bioconductor packages should also return the type of object returned by the default method.

## See Also

- `base::table` for the default `table` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- table,Rle-method in the **S4Vectors** package for an example of a specific `table` method (defined for Rle objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
table
showMethods("table")
selectMethod("table", "ANY")  # the default method
```

tapply *Apply a function over a ragged array*

## Description

tapply applies a function to each cell of a ragged array, that is to each (non-empty) group of values given by a unique combination of the levels of certain factors.

NOTE: This man page is for the tapply *S4 generic function* defined in the **BiocGenerics** package. See ?base::tapply for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically list-like or vector-like) not supported by the default method.

## Usage

```
tapply(X, INDEX, FUN=NULL, ..., default=NA, simplify=TRUE)
```

## Arguments

| | |
|---|---|
| X | The default method expects an atomic object, typically a vector. See ?base::tapply for the details.<br>Specific methods can support other objects (typically list-like or vector-like). Please refer to the documentation of a particular method for the details. |
| INDEX | The default method expects a list of one or more factors, each of same length as X. See ?base::tapply for the details.<br>Specific methods can support other objects (typically list-like). Please refer to the documentation of a particular method for the details. |
| FUN, ..., default, simplify | See ?base::tapply for a description of these arguments. |

## Value

See ?base::tapply for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- base::tapply for the default tapply method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- tapply,Vector,ANY-method in the **IRanges** package for an example of a specific tapply method (defined for Vector objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
tapply  # note the dispatch on the 'X' and 'INDEX' args only
showMethods("tapply")
selectMethod("tapply", c("ANY", "ANY"))  # the default method
```

---

`testPackage` *Run RUnit package unit tests*

---

## Description

`testPackage` helps developers implement unit tests using the **RUnit** testing conventions.

## Usage

```
testPackage(pkgname=NULL, subdir="unitTests", pattern="^test_.*\\.R$",
    path=getwd())
```

## Arguments

| | |
|---|---|
| `pkgname` | The name of the package whose installed unit tests are to be run. A missing or NULL value implies that the `testPackage` command will look for tests within the package source directory indicated by `path`. |
| `subdir` | A character(1) vector providing the subdirectory in which unit tests are located. The directory is searched first in the (installed or source) package root, or in a subdirectory `inst/` below the root. |
| `pattern` | A character(1) regular expression describing the file names to be evaluated; typically used to restrict tests to a subset of all test files. |
| `path` | A character(1) directory path indicating, when `pkgname` is missing or NULL, where unit tests will be searched. `path` can be any location at or below the package root. |

## Details

This function is not exported from the package namespace, and must be invoked using triple colons, `BiocGenerics:::testPackage()`; it is provided primarily for the convenience of developers.

When invoked with missing or NULL `pkgname` argument, the function assumes that it has been invoked from within the package source tree (or that the source tree is located above `path`), and finds unit tests in `subdir="unitTests"` in either the base or `inst/` directories at the root of the package source tree. This mode is useful when developing unit tests, since the package does not have to be re-installed to run an updated test.

When invoked with `pkgname` set to the name of an installed package, unit tests are searched for in the installed package directory.

## Value

The function returns the result of `RUnit::runTestSuite` invoked on the unit tests specified in the function call.

## See Also

http://bioconductor.org/developers/how-to/unitTesting-guidelines/

## Examples

```
## Run unit tests found in the library location where
## BiocGenerics is installed
BiocGenerics:::testPackage("BiocGenerics")
## Not run: ## Run unit tests for the package whose source tree implied
## by getwd()
BiocGenerics:::testPackage()

## End(Not run)
```

---

# toTable *An alternative to as.data.frame()*

---

## Description

`toTable()` is an *S4 generic function* provided as an alternative to `as.data.frame()`.

## Usage

```
toTable(x, ...)
```

## Arguments

| | |
|---|---|
| `x` | The object to turn into a data frame. |
| `...` | Additional arguments, for use in specific methods. |

## Value

A data frame.

## See Also

- The `as.data.frame` *S4 generic* defined in the **BiocGenerics** package.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- toTable,Bimap-method in the **AnnotationDbi** package for an example of a specific `toTable` method (defined for Bimap objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
toTable
showMethods("toTable")

library(AnnotationDbi)
showMethods("toTable")
selectMethod("toTable", "Bimap")
```

---

type *Accessing the type of an object*

---

## Description

Get or set the *type* of an object.

Note that type and type<- are defined as *S4 generic functions* and what *type* means exactly (and what type() returns) depends on the objects for which the type and/or type<- methods are defined.

## Usage

```
type(x)
type(x) <- value

## Methods defined in the BiocGenerics package:

## S4 method for signature 'vector'
type(x)
## S4 method for signature 'array'
type(x)
## S4 method for signature 'factor'
type(x)  # returns "character"
## S4 method for signature 'data.frame'
type(x)

## S4 replacement method for signature 'vector'
type(x) <- value
## S4 replacement method for signature 'array'
type(x) <- value
```

## Arguments

| | |
|---|---|
| x | Any object for which the type() getter or setter is defined. Note that objects will either: not support the getter or setter at all, or support only the getter, or support both the getter and setter. |
| value | The type to set on x (assuming x supports the type() setter). value is typically (but not necessarily) expected to be a single string (i.e. a character vector of length 1). |

## Details

On an ordinary vector, matrix, or array `x`, `type(x)` returns `typeof(x)`.

On a data frame `x` where all the columns are ordinary vectors or factors, `type(x)` is *semantically equivalent* to `typeof(as.matrix(x))`. However, the actual implementation is careful to avoid turning the full data frame `x` into a matrix, as this would tend to be very inefficient in general.

Note that for a matrix-like or array-like object, `type(x)` returns the type of the *elements* in the object. See `?S4Arrays::type` for more information.

## Value

`type(x)` is expected to return the type of `x` as a single string i.e. as a character vector of length 1.

## See Also

- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- type,ANY-method in the **S4Arrays** package for the default `type` method.
- type,DataFrame-method in the **S4Arrays** package, and type,PairwiseAlignments-method in the **pwalign** package, for examples of specific `type` methods (defined for DataFrame and PairwiseAlignments objects, respectively).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
type
showMethods("type")

`type<-`
showMethods("type<-")

## The BiocGenerics package defines methods for ordinary vectors, arrays,
## and data frames:
m <- matrix(11:22, nrow=3)
type(m)           # equivalent to 'typeof(m)' or 'storage.mode(m)'
type(m) <- "raw"  # equivalent to 'storage.mode(m) <- "raw"'
m
type(m)

selectMethod("type", "array")

selectMethod("type<-", "array")

df <- data.frame(a=44:49, b=letters[1:6], c=c(TRUE, FALSE))
stopifnot(identical(type(df), typeof(as.matrix(df))))

## Examples of methods defined in other packages:

library(S4Arrays)
showMethods("type")
```

```
selectMethod("type", "ANY")  # the default "type" method

library(pwalign)
showMethods("type")
## The type() method for PairwiseAlignments objects:
selectMethod("type", "PairwiseAlignments")
```

---

unique *Extract unique elements*

---

## Description

unique returns an object of the same class as x (typically a vector-like, data-frame-like, or array-like object) but with duplicate elements/rows removed.

NOTE: This man page is for the unique *S4 generic function* defined in the **BiocGenerics** package. See ?base::unique for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-like or data-frame-like) not supported by the default method.

## Usage

```
unique(x, incomparables=FALSE, ...)
```

## Arguments

| | |
|---|---|
| x | A vector-like, data-frame-like, or array-like object. |
| incomparables, ... | See ?base::unique for a description of these arguments. |

## Value

See ?base::unique for the value returned by the default method.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input object.

unique should always behave consistently with BiocGenerics::duplicated.

## See Also

- base::unique for the default unique method.
- BiocGenerics::duplicated for determining duplicate elements.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- unique,Rle-method in the **S4Vectors** package for an example of a specific unique method (defined for Rle objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
unique
showMethods("unique")
selectMethod("unique", "ANY")  # the default method
```

---

`unlist` *Flatten list-like objects*

---

## Description

Given a list-like object `x`, `unlist` produces a vector-like object obtained by concatenating (conceptually thru `c`) all the top-level elements in `x` (each of them being expected to be a vector-like object, typically).

NOTE: This man page is for the `unlist` *S4 generic function* defined in the **BiocGenerics** package. See `?base::unlist` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
unlist(x, recursive=TRUE, use.names=TRUE)
```

## Arguments

| | |
|---|---|
| `x` | A list-like object. |
| `recursive, use.names` | See `?base::unlist` for a description of these arguments. |

## Value

See `?base::unlist` for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- `base::unlist` for the default `unlist` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- unlist,List-method in the **S4Vectors** package for an example of a specific `unlist` method (defined for List objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
unlist  # note the dispatch on the 'x' arg only
showMethods("unlist")
selectMethod("unlist", "ANY")  # the default method
```

unsplit *Unsplit a list-like object*

## Description

Given a list-like object `value` and grouping `f`, `unsplit` produces a vector-like object `x` by conceptually reversing the split operation `value <- split(x, f)`.

NOTE: This man page is for the `unsplit` *S4 generic function* defined in the **BiocGenerics** package. See `?base::unsplit` for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
unsplit(value, f, drop=FALSE)
```

## Arguments

| | |
|---|---|
| `value` | A list-like object. |
| `f` | A factor or other grouping object that corresponds to the `f` symbol in `value <- split(x, f)`. |
| `drop` | See `?base::unsplit` for a description of this argument. |

## Value

See `?base::unsplit` for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- `base::unsplit` for the default `unsplit` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- unsplit,List-method in the **IRanges** package for an example of a specific `unsplit` method (defined for List objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
unsplit  # note the dispatch on the 'value' and 'f' args only
showMethods("unsplit")
selectMethod("unsplit", "ANY")  # the default method
```

`updateObject` *Update an object to its current class definition*

## Description

`updateObject` is a generic function that returns an instance of `object` updated to its current class definition.

## Usage

```
updateObject(object, ..., verbose=FALSE)

## Related utilities:
updateObjectFromSlots(object, objclass=class(object)[[1L]], ...,
                      verbose=FALSE)
getObjectSlots(object)
```

## Arguments

| | |
|---|---|
| `object` | Object to be updated for `updateObject` and `updateObjectFromSlots`. Object for slot information to be extracted from for `getObjectSlots`. |
| `...` | Additional arguments, for use in specific `updateObject` methods. |
| `verbose` | `TRUE` or `FALSE`, indicating whether information about the update should be reported. Use `message` to report this information. |
| `objclass` | Optional character string naming the class of the object to be created. |

## Details

Updating objects is primarily useful when an object has been serialized (e.g., stored to disk) for some time (e.g., months), and the class definition has in the mean time changed. Because of the changed class definition, the serialized instance is no longer valid.

`updateObject` requires that the class of the returned object be the same as the class of the argument `object`, and that the object is valid (see `validObject`). By default, `updateObject` has the following behaviors:

`updateObject(ANY, ..., verbose=FALSE)` By default, `updateObject` uses heuristic methods to determine whether the object should be the 'new' S4 type (introduced in R 2.4.0), but is not. If the heuristics indicate an update is required, the `updateObjectFromSlots` function tries to update the object. The default method returns the original S4 object or the successfully updated object, or issues an error if an update is required but not possible. The optional named argument `verbose` causes a message to be printed describing the action. Arguments `...` are passed to `updateObjectFromSlots`.

`updateObject(environment, ..., verbose=FALSE)` Visit each element in `environment`, applying `updateObject(environment[[elt]], ..., verbose=verbose)`

`updateObject(formula, ..., verbose=FALSE)` Do nothing; the environment of the formula may be too general (e.g., `R_GlobalEnv`) to attempt an update.

`updateObject(envRefClass, ..., verbose=FALSE)` Attempt to update objects from fields using a strategy like `updateObjectFromSlots` Method 1.

`updateObjectFromSlots(object, objclass=class(object), ..., verbose=FALSE)` is a utility function that identifies the intersection of slots defined in the `object` instance and `objclass` definition. Under Method 1, the corresponding elements in `object` are then updated (with `updateObject(elt, ..., verbose=verbose)`) and used as arguments to a call to `new(class, ...)`, with `...` replaced by slots from the original object. If this fails, then Method 2 tries `new(class)` and assigns slots of `object` to the newly created instance.

`getObjectSlots(object)` extracts the slot names and contents from `object`. This is useful when `object` was created by a class definition that is no longer current, and hence the contents of `object` cannot be determined by accessing known slots.

## Value

`updateObject` returns a valid instance of `object`.

`updateObjectFromSlots` returns an instance of class `objclass`.

`getObjectSlots` returns a list of named elements, with each element corresponding to a slot in `object`.

## See Also

- `updateObjectTo` in the **Biobase** package for updating an object to the class definition of a template (might be useful for updating a virtual superclass).
- `validObject` for testing the validity of an object.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
updateObject
showMethods("updateObject")
selectMethod("updateObject", "ANY")  # the default method

library(Biobase)
## update object, same class
data(sample.ExpressionSet)
obj <- updateObject(sample.ExpressionSet)

setClass("UpdtA", representation(x="numeric"), contains="data.frame")
setMethod("updateObject", "UpdtA",
    function(object, ..., verbose=FALSE)
    {
        if (verbose)
            message("updateObject object = 'A'")
        object <- callNextMethod()
        object@x <- -object@x
        object
```

```
    }
)

a <- new("UpdtA", x=1:10)
## See steps involved
updateObject(a)

removeMethod("updateObject", "UpdtA")
removeClass("UpdtA")
```

---

var *Variance and Standard Deviation*

---

## Description

var and sd compute the variance and standard deviation of a vector x.

NOTE: This man page is for the var and sd, *S4 generic functions* defined in the **BiocGenerics** package. See ?stats::var and ?stats::sd for the default methods (defined in the **stats** package). Bioconductor packages can define specific methods for objects (typically array-like) not supported by the default method.

## Usage

```
var(x, y = NULL, na.rm = FALSE, use)
sd(x, na.rm = FALSE)
```

## Arguments

| | |
|---|---|
| x | a vector-like object |
| y | a vector-like object, or NULL |
| na.rm, use | see var |

## Value

See ?stats::var and ?stats::sd for the value returned by the default methods.

Specific methods defined in Bioconductor packages will typically return an object of the same class as the input object.

## See Also

- stats::var and stats::sd for the default methods.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

### Examples

```
var
showMethods("var")
selectMethod("var", "ANY")  # the default method
```

---

`weights` *Extract model weights*

---

### Description

`weights` is a generic function which extracts fitting weights from objects returned by modeling functions.

NOTE: This man page is for the `weights` *S4 generic function* defined in the **BiocGenerics** package. See `?stats::weights` for the default method (defined in the **stats** package). Bioconductor packages can define specific methods for objects not supported by the default method.

### Usage

```
weights(object, ...)
```

### Arguments

| | |
|---|---|
| `object, ...` | See `?stats::weights`. |

### Value

Weights extracted from the object `object`.

See `?stats::weights` for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

### See Also

- `stats::weights` for the default `weights` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- weights,PLMset-method in the **affyPLM** package for an example of a specific `weights` method (defined for PLMset objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

### Examples

```
weights
showMethods("weights")
selectMethod("weights", "ANY")  # the default method
```

which *Which values in an object are considered TRUE?*

## Description

Give the indices of the values in a vector-, array-, or list-like object that are considered TRUE, allowing for array indices in the case of an array-like object.

NOTE: This man page is for the which *S4 generic function* defined in the **BiocGenerics** package. See ?base::which for the default method (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-, array-, or list-like) not supported by the default methods.

## Usage

```
which(x, arr.ind=FALSE, useNames=TRUE)
```

## Arguments

| | |
|---|---|
| x | An object, typically with a vector-, array-, or list-like semantic. |
| arr.ind, useNames | See ?base::which for a description of these arguments. |

## Value

See ?base::which for the value returned by the default method.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default method.

## See Also

- base::which for the default which method.
- showMethods for displaying a summary of the methods defined for a given generic function.
- selectMethod for getting the definition of a specific method.
- which,DelayedArray-method in the **DelayedArray** package for an example of a specific which method (defined for DelayedArray objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
which
showMethods("which")
selectMethod("which", "ANY")  # the default method

library(DelayedArray)
showMethods("which")
## The which() method for DelayedArray objects:
selectMethod("which", "DelayedArray")
```

which.min *What's the index of the first min or max value in an object?*

## Description

Determines the location (i.e. index) of the (first) minimum or maximum value in an object.

NOTE: This man page is for the `which.min` and `which.max` *S4 generic functions* defined in the **BiocGenerics** package. See ?`base::which.min` for the default methods (defined in the **base** package). Bioconductor packages can define specific methods for objects (typically vector-, array-, or list-like) not supported by the default methods.

## Usage

```
which.min(x, ...)
which.max(x, ...)
```

## Arguments

| | |
|---|---|
| `x` | An object, typically with a vector-, array-, or list-like semantic. |
| `...` | Additional arguments, for use in specific methods. |

## Value

See ?`base::which.min` for the value returned by the default methods.

Specific methods defined in Bioconductor packages should behave as consistently as possible with the default methods.

## Note

The default methods (defined in the **base** package) only take a single argument. We've added the `...` argument to the generic functions defined in the **BiocGenerics** package so they can be called with an arbitrary number of effective arguments. This typically allows methods to add extra arguments for controlling/altering the behavior of the operation. Like for example the `global` argument supported by the `which.max` method for NumericList objects (defined in the **IRanges** package).

## See Also

- `base::which.min` for the default `which.min` and `which.max` methods.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- which.max,NumericList-method in the **IRanges** package for an example of a specific `which.max` method (defined for NumericList objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
which.min
showMethods("which.min")
selectMethod("which.min", "ANY")  # the default method

which.max
showMethods("which.max")
selectMethod("which.max", "ANY")  # the default method

library(IRanges)
showMethods("which.max")
## The which.max() method for NumericList objects:
selectMethod("which.max", "NumericList")
```

---

# xtabs *Cross tabulation*

---

## Description

xtabs creates a contingency table (optionally a sparse matrix) from cross-classifying factors, usually contained in a data-frame-like object, using a formula interface.

NOTE: This man page is for the xtabs *S4 generic function* defined in the **BiocGenerics** package. See ?stats::xtabs for the default method (defined in the **stats** package). Bioconductor packages can define specific methods for objects not supported by the default method.

## Usage

```
xtabs(formula=~., data=parent.frame(), subset, sparse=FALSE,
      na.action, na.rm=FALSE, addNA=FALSE, exclude=if(!addNA)c(NA, NaN),
      drop.unused.levels=FALSE)
```

## Arguments

| | |
|---|---|
| formula, subset, sparse, na.action, na.rm, addNA, exclude, drop.unused.levels | See ?stats::xtabs for a description of these arguments. |
| data | A data-frame-like object. |

## Value

See ?stats::xtabs for the value returned by the default method.

Specific methods defined in Bioconductor packages should also return the type of object returned by the default method.

## See Also

- `stats::xtabs` for the default `xtabs` method.
- `showMethods` for displaying a summary of the methods defined for a given generic function.
- `selectMethod` for getting the definition of a specific method.
- xtabs,DataFrame-method in the **S4Vectors** package for an example of a specific `xtabs` method (defined for DataFrame objects).
- BiocGenerics for a summary of all the generics defined in the **BiocGenerics** package.

## Examples

```
xtabs  # note the dispatch on the 'data' arg only
showMethods("xtabs")
selectMethod("xtabs", "ANY")  # the default method

library(S4Vectors)
showMethods("xtabs")
## The xtabs() method for DataFrame objects:
selectMethod("xtabs", "DataFrame")
```

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