

Package ‘BiocGenerics’

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Title S4 generic functions used in Bioconductor

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

biocViews Infrastructure

URL <https://bioconductor.org/packages/BiocGenerics>

BugReports <https://github.com/Bioconductor/BiocGenerics/issues>

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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
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git_url <https://git.bioconductor.org/packages/BiocGenerics>

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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.

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

<code>a</code>	An array-like object.
<code>perm, ...</code>	See <code>?base::aperm</code> 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
```


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 matrices 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 matrices.

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])
```



```
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, mode="any", ifnotfound, inherits=FALSE)
```

Arguments

`x` For `get`: A variable name (or, more generally speaking, a *key*), given as a single string.
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.

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 names:
x <- c(A="foo", B="bar")
paste0(x, "XX") # names are lost
paste2(x, "XX") # 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]))
paste0(m, letters[1:4]) # dimensions and dimnames are lost
paste2(m, letters[1:4]) # 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	<i>Accessing the path of an object</i>
------	--

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 <code>path<-</code> , the paths to set on object. For <code>basename<-</code> or <code>dirname<-</code> , the basenames or dirnames to set on path.
path	A character vector <i>or an object containing paths</i> .

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)

```

Description

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

Usage

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

Arguments

<code>object</code>	A data object, typically containing count values from an RNA-Seq experiment or microarray intensity values.
<code>...</code>	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	<i>Ranks the values in a vector-like object</i>
------	---

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.

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

<code>flesh</code>	A vector-like object.
<code>skeleton</code>	A list-like object. Only the "shape" (i.e. the lengths of the individual list elements) of <code>skeleton</code> 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()`d.

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

<code>x</code>	The object to replicate (typically vector-like).
<code>times</code>	See <code>?base::rep.int</code> 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	<i>Extract model residuals</i>
-----------	--------------------------------

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

<code>x</code>	A matrix-like object.
<code>do.NULL</code> , <code>prefix</code>	See <code>?base::rownames</code> for a description of these arguments.
<code>value</code>	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 **Bioc-Generics** 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. Additionally 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	<i>Score accessor</i>
-------	-----------------------

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")
```

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

<code>x, y</code>	Vector-like objects (typically of the same class, but not necessarily).
<code>...</code>	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 allows 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 operations, 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 dispatch 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 dispatch 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 <code>start()</code> , <code>end()</code> , and <code>width()</code> getters/setters: an object containing start, end, and width values. For the <code>pos{}</code> 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 <code>?base::subset</code> 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	<i>Cross tabulation and table creation</i>
-------	--

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.
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.
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	<i>Run RUnit package unit tests</i>
-------------	-------------------------------------

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	<i>An alternative to as.data.frame()</i>
---------	--

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 <code>type()</code> 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 <code>type()</code> 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 **pwalgn** 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 <code>value <- split(x, f)</code> .
drop	See <code>?base::unsplit</code> 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	<i>Update an object to its current class definition</i>
--------------	---

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

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

<code>x</code>	a vector-like object
<code>y</code>	a vector-like object, or <code>NULL</code>
<code>na.rm, use</code>	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

<code>x</code>	An object, typically with a vector-, array-, or list-like semantic.
<code>...</code>	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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