Package 'RaggedExperiment'

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Title Representation of Sparse Experiments and Assays Across Samples

Version 1.4.0

Description This package provides a flexible representation of copy number, mutation, and other data that fit into the ragged array schema for genomic location data. The basic representation of such data provides a rectangular flat table interface to the user with range information in the rows and samples/specimen in the columns.

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biocViews Infrastructure, DataRepresentation

BugReports https://github.com/Bioconductor/RaggedExperiment/issues

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R topics documented:

RaggedExperiment-J	oackag	ge .						•		 													 	2
assay-functions	••••	•••						•		 													 	2
RaggedExperiment-o	lass	•••						•		 													 	5
sparseSummarizedE	cperim	nent		•	•	•	 •	•	•	 	•	•		•	•		•	•	•	•	•	•	 	10

13

Index

RaggedExperiment-package

RaggedExperiment: Range-based data representation package

Description

RaggedExperiment allows the user to represent, copy number, mutation, and other types of rangebased data formats where optional information about samples can be provided. At the backbone of this package is the GRangesList class. The RaggedExperiment class uses this representation and presents the data in a couple of different ways:

- rowRanges
- colData

The rowRanges method will return the internal GRangesList representation of the dataset. A distinction between the SummarizedExperiment and the RaggedExperiment classes is that the RaggedExperiment class allows for ragged ranges, meaning that there may be a different number of ranges or rows per sample.

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See Also

Useful links:

• Report bugs at https://github.com/Bioconductor/RaggedExperiment/issues

assay-functions Create simplified representation of ragged assay data.

Description

These methods transform assay() from the default (i.e., sparseAssay()) representation to various forms of more dense representation. compactAssay() collapses identical ranges across samples into a single row. disjoinAssay() creates disjoint (non-overlapping) regions, simplifies values within each sample in a user-specified manner, and returns a matrix of disjoint regions x samples.

This method transforms assay() from the default (i.e., sparseAssay()) representation to a reduced representation summarizing each original range overlapping ranges in query. Reduction in each cell can be tailored to indivdual needs using the simplifyReduce functional argument.

assay-functions

Usage

```
sparseAssay(x, i = 1, withDimnames = TRUE, background = NA_integer_)
compactAssay(x, i = 1, withDimnames = TRUE, background = NA_integer_)
disjoinAssay(x, simplifyDisjoin, i = 1, withDimnames = TRUE,
background = NA_integer_)
qreduceAssay(x, query, simplifyReduce, i = 1, withDimnames = TRUE,
background = NA_integer_)
```

Arguments

х	A RaggedExperiment object
i	integer(1) or character(1) name of assay to be transformed.
withDimnames	logical(1) include dimnames on the returned matrix. When there are no explict rownames, these are manufactured with as.character(rowRanges(x)); rownames are always manufactured for compactAssay() and disjoinAssay().
background	A value (default NA) for the returned matrix after *Assay operations
simplifyDisjoin	
	A function / functional operating on a *List, where the elements of the list are all within-sample assay values from ranges overlapping each disjoint range. For instance, to use the simplifyDisjoin=mean of overlapping ranges, where ranges are characterized by integer-valued scores, the entries are calculated as

query

GRanges providing regions over which reduction is to occur.

simplifyReduce A function / functional accepting arguments score, range, and qrange:

- score A *List, where each list element corresponds to a cell in the matrix to be returned by qreduceAssay. Vector elements correspond to ranges overlapping query. The *List objects support many vectorized mathematical operations, so simplifyReduce can be implemented efficiently.
- range A GRangesList instance, 'parallel' to score. Each element of the list corresponds to a cell in the matrix to be returned by qreduceAssay. Each range in the element corresponds to the range for which the score element applies.
- qrange A GRanges instance with the same length as score, providing the query range window to which the corresponding scores apply.

Value

sparseAssay(): A matrix() with dimensions dim(x). Elements contain the assay value for the *i*th range and *j*th sample.

compactAssay(): Samples with identical range are placed in the same row. Non-disjoint ranges are NOT collapsed.

disjoinAssay(): A matrix with number of rows equal to number of disjoint ranges across all samples. Elements of the matrix are summarized by applying simplifyDisjoin() to assay values of overlapping ranges

qreduceAssay(): A matrix() with dimensions $length(query) \times ncol(x)$. Elements contain assay values for the ith query range and jth sample, summarized according to the function simplifyReduce.

Examples

```
x <- RaggedExperiment(GRangesList(</pre>
    GRanges(c("A:1-3", "A:4-5", "A:10-15"), score=1:3),
    GRanges(c("A:4-5", "B:1-3"), score=4:5)
))
query <- GRanges(c("A:1-2", "A:4-5", "B:1-5"))</pre>
weightedmean <- function(scores, ranges, qranges)</pre>
    ## weighted average score per query range
    sum(scores * width(ranges)) / sum(width(ranges))
qreduceAssay(x, query, weightedmean)
## Not run:
    ## Extended example: non-silent mutations, summarized by genic
    ## region
    suppressPackageStartupMessages({
        library(TxDb.Hsapiens.UCSC.hg19.knownGene)
        library(org.Hs.eg.db)
        library(GenomeInfoDb)
        library(MultiAssayExperiment)
    })
    ## TCGA Multi-assay experiment to RaggedExperiment
    url <- "http://s3.amazonaws.com/multiassayexperiments/accMAE0.rds"</pre>
    ## download.file(url, fl <- tempfile())</pre>
    ## fl <- "accMAE0.rds"</pre>
    mae <- readRDS(f1)[, , c("RNASeq2GeneNorm", "CNASNP", "Mutations")]</pre>
    ## genomic coordinates
    gn <- genes(TxDb.Hsapiens.UCSC.hg19.knownGene)</pre>
    gn <- keepStandardChromosomes(granges(gn), pruning.mode="coarse")</pre>
    seqlevelsStyle(gn) <- "NCBI"</pre>
    ## reduce mutations, marking any genomic range with non-silent
    ## mutation as FALSE
    nonsilent <- function(scores, ranges, granges)</pre>
        any(scores != "Silent")
    re <- as(mae[["Mutations"]], "RaggedExperiment")</pre>
    mutations <- qreduceAssay(re, gn, nonsilent, "Variant_Classification")</pre>
```

4

```
## reduce copy number
re <- as(mae[["CNASNP"]], "RaggedExperiment")
cn <- qreduceAssay(re, gn, weightedmean, "Segment_Mean")
## End(Not run)</pre>
```

RaggedExperiment-class

RaggedExperiment objects

Description

The RaggedExperiment class is a container for storing range-based data, including but not limited to copy number data, and mutation data. It can store a collection of GRanges objects, as it is derived from the GenomicRangesList.

Usage

```
RaggedExperiment(..., colData = DataFrame())
## S4 method for signature 'RaggedExperiment'
rowRanges(x, ...)
## S4 method for signature 'RaggedExperiment'
mcols(x, use.names = FALSE, ...)
## S4 replacement method for signature 'RaggedExperiment'
mcols(x, ...) <- value</pre>
## S4 method for signature 'RaggedExperiment'
rowData(x, ...)
## S4 replacement method for signature 'RaggedExperiment'
rowData(x, ...) <- value</pre>
## S4 method for signature 'RaggedExperiment'
dim(x)
## S4 method for signature 'RaggedExperiment'
dimnames(x)
## S4 replacement method for signature 'RaggedExperiment,list'
dimnames(x) <- value
## S4 method for signature 'RaggedExperiment'
length(x)
## S4 method for signature 'RaggedExperiment'
colData(x, ...)
```

```
## S4 replacement method for signature 'RaggedExperiment,DataFrame'
colData(x) <- value</pre>
## S4 method for signature 'RaggedExperiment,missing'
assay(x, i, ...)
## S4 method for signature 'RaggedExperiment,ANY'
assay(x, i, ..., withDimnames = TRUE)
## S4 method for signature 'RaggedExperiment'
assays(x, ..., withDimnames = TRUE)
## S4 method for signature 'RaggedExperiment'
assayNames(x, ...)
## S4 method for signature 'RaggedExperiment'
show(object)
## S4 method for signature 'RaggedExperiment,ANY,ANY,ANY'
x[i, j, ..., drop = TRUE]
## S4 method for signature 'RaggedExperiment,Vector'
overlapsAny(query, subject, maxgap = 0L,
 minoverlap = 1L, type = c("any", "start", "end", "within", "equal"), ...)
## S4 method for signature 'RaggedExperiment,Vector'
subsetByOverlaps(x, ranges, maxgap = -1L,
 minoverlap = 0L, type = c("any", "start", "end", "within", "equal"),
 invert = FALSE, ...)
```

Arguments

	Constructor: GRanges, list of GRanges, or GRangesList OR assay: Additional arguments for assay. See details for more information.
colData	A DataFrame describing samples. Length of rowRanges must equal the number of rows in colData
х	A RaggedExperiment object.
use.names	(logical default FALSE) whether to propagate rownames from the object to rownames of metadata ${\tt DataFrame}$
value	• dimnames: A list of dimension names
	• mcols: A DataFrame representing the assays
i	logical(1), integer(1), or character(1) indicating the assay to be reported. For [, i can be any supported Vector object, e.g., GRanges.
withDimnames	logical (default TRUE) whether to use dimension names in the resulting object
object	A RaggedExperiment object.
j	integer(), character(), or logical() index selecting columns from RaggedExperiment
drop	logical (default TRUE) whether to drop empty samples
query	A RaggedExperiment instance.

6

subject	Each of them can be an IntegerRanges (e.g. IRanges, Views) or IntegerRanges- List (e.g. IRangesList, ViewsList) derivative. In addition, if subject or ranges is an IntegerRanges object, query or x can be an integer vector to be converted to length-one ranges.
	If query (or x) is an IntegerRangesList object, then subject (or ranges) must also be an IntegerRangesList object.
	If both arguments are list-like objects with names, each list element from the 2nd argument is paired with the list element from the 1st argument with the matching name, if any. Otherwise, list elements are paired by position. The overlap is then computed between the pairs as described below.
	If subject is omitted, query is queried against itself. In this case, and only this case, the drop.self and drop.redundant arguments are allowed. By default, the result will contain hits for each range against itself, and if there is a hit from A to B, there is also a hit for B to A. If drop.self is TRUE, all self matches are dropped. If drop.redundant is TRUE, only one of A->B and B->A is returned.
maxgap	A single integer $>= -1$.
	If type is set to "any", maxgap is interpreted as the maximum <i>gap</i> that is allowed between 2 ranges for the ranges to be considered as overlapping. The <i>gap</i> between 2 ranges is the number of positions that separate them. The <i>gap</i> between 2 adjacent ranges is 0. By convention when one range has its start or end strictly inside the other (i.e. non-disjoint ranges), the <i>gap</i> is considered to be -1.
	If type is set to anything else, maxgap has a special meaning that depends on the particular type. See type below for more information.
minoverlap	A single non-negative integer.
	Only ranges with a minimum of minoverlap overlapping positions are considered to be overlapping.
	When type is "any", at least one of maxgap and minoverlap must be set to its default value.
type	By default, any overlap is accepted. By specifying the type parameter, one can select for specific types of overlap. The types correspond to operations in Allen's Interval Algebra (see references). If type is start or end, the intervals are required to have matching starts or ends, respectively. Specifying equal as the type returns the intersection of the start and end matches. If type is within, the query interval must be wholly contained within the subject interval. Note that all matches must additionally satisfy the minoverlap constraint described above.
	The maxgap parameter has special meaning with the special overlap types. For start, end, and equal, it specifies the maximum difference in the starts, ends or both, respectively. For within, it is the maximum amount by which the subject may be wider than the query. If maxgap is set to -1 (the default), it's replaced internally by 0.
ranges	Each of them can be an IntegerRanges (e.g. IRanges, Views) or IntegerRanges- List (e.g. IRangesList, ViewsList) derivative. In addition, if subject or ranges is an IntegerRanges object, query or x can be an integer vector to be converted to length-one ranges.
	If query (or x) is an IntegerRangesList object, then subject (or ranges) must also be an IntegerRangesList object.
	If both arguments are list-like objects with names, each list element from the 2nd argument is paired with the list element from the 1st argument with the matching

	name, if any. Otherwise, list elements are paired by position. The overlap is then computed between the pairs as described below.
	If subject is omitted, query is queried against itself. In this case, and only this
	case, the drop.self and drop.redundant arguments are allowed. By default,
	the result will contain hits for each range against itself, and if there is a hit from
	A to B, there is also a hit for B to A. If drop.self is TRUE, all self matches are
	dropped. If drop.redundant is TRUE, only one of A->B and B->A is returned.
invert	If TRUE, keep only the ranges in x that do <i>not</i> overlap ranges.

Value

constructor returns a RaggedExperiment object

'rowRanges' returns a GRanges object summarizing ranges corresponding to assay() rows.

'mcols' returns a DataFrame object of the metadata columns

'assays' returns a SimpleList

'overlapsAny' returns a logical vector of length equal to the number of rows in the query; TRUE when the copy number region overlaps the subject.

'subsetByOverlaps' returns a RaggedExperiment containing only copy number regions overlapping subject.

Methods (by generic)

- rowRanges: rowRanges accessor
- mcols: get the metadata columns of the ranges, rectangular representation of the 'assays'
- mcols<-: set the metadata columns of the ranges corresponding to the assays
- rowData: get the rowData or metadata for the ranges
- rowData<-: set the rowData or metadata for the ranges
- dim: get dimensions (number of sample-specific row ranges by number of samples)
- dimnames: get row (sample-specific) range names and sample names
- dimnames<-: set row (sample-specific) range names and sample names
- · length: get the length of row vectors in the object, similar to SummarizedExperiment
- colData: get column data
- colData<-: change the colData
- · assay: assay missing method uses first metadata column
- assay: assay numeric method.
- assays: assays
- assayNames: names in each assay
- · show: show method
- [: Subset a RaggedExperiment object
- overlapsAny: Determine whether copy number ranges defined by query overlap ranges of subject.
- subsetByOverlaps: Subset the RaggedExperiment to contain only copy number ranges overlapping ranges of subject.

RaggedExperiment-class

Constructors

RaggedExperiment(..., colData=DataFrame()): Creates a RaggedExperiment object using multiple GRanges objects or a list of GRanges objects. Additional column data may be provided as a DataFrame object.

Accessors

In the following, 'x' represents a RaggedExperiment object:

rowRanges(x):

Get the ranged data. Value is a GenomicRanges object.

assays(x):

Get the assays. Value is a SimpleList.

assay(x, i):

An alternative to assays(x)[[i]] to get the *i*th (default first) assay element.

mcols(x), mcols(x) <- value:</pre>

Get or set the metadata columns. For RaggedExperiment, the columns correspond to the assay *i*th elements.

rowData(x), rowData(x) <- value:</pre>

Get or set the row data. Value is a DataFrame object. Also corresponds to the mcols data.

Note for advanced users and developers. Both mcols and rowData setters may reduce the size of the internal RaggedExperiment data representation. Particularly after subsetting, the internal row index is modified and such setter operations will use the index to subset the data and reduce the "rows" of the internal data representation.

Subsetting

x[i, j]: Get ranges or elements (i and j, respectively) with optional metadata columns where i or j can be missing, an NA-free logical, numeric, or character vector.

Coercion

In the following, 'object' represents a RaggedExperiment object:

as(object, "GRangesList"):

Creates a GRangesList object from a RaggedExperiment.

as(from, "RaggedExperiment"):

Creates a RaggedExperiment object from a GRangesList, or GRanges object.

Examples

```
## Create an empty RaggedExperiment instance
re0 <- RaggedExperiment()
re0
## Create a couple of GRanges objects with row ranges names
sample1 <- GRanges(
    c(a = "chr1:1-10:-", b = "chr1:11-18:+"),
    score = 1:2)
sample2 <- GRanges(
    c(c = "chr2:1-10:-", d = "chr2:11-18:+"),</pre>
```

```
score = 3:4)
## Include column data
colDat <- DataFrame(id = 1:2)</pre>
## Create a RaggedExperiment object from a couple of GRanges
re1 <- RaggedExperiment(sample1=sample1, sample2=sample2, colData = colDat)</pre>
re1
## With list of GRanges
lgr <- list(sample1 = sample1, sample2 = sample2)</pre>
## Create a RaggedExperiment from a list of GRanges
re2 <- RaggedExperiment(lgr, colData = colDat)</pre>
grl <- GRangesList(sample1 = sample1, sample2 = sample2)</pre>
## Create a RaggedExperiment from a GRangesList
re3 <- RaggedExperiment(grl, colData = colDat)</pre>
## Subset a RaggedExperiment
assay(re3[c(1, 3),])
subsetByOverlaps(re3, GRanges("chr1:1-5")) # by ranges
```

sparseSummarizedExperiment

Create SummarizedExperiment representations by transforming ragged assays to rectangular form.

Description

These methods transform RaggedExperiment objects to similar SummarizedExperiment objects. They do so by transforming assay data to more rectangular representations, following the rules outlined for similarly names transformations sparseAssay(), compactAssay(), disjoinAssay(), and qreduceAssay(). Because of the complexity of the transformation, ti usually only makes sense transform RaggedExperiment objects with a single assay; this is currently enforced at time of coercion.

Usage

```
sparseSummarizedExperiment(x, i = 1, withDimnames = TRUE)
compactSummarizedExperiment(x, i = 1L, withDimnames = TRUE)
disjoinSummarizedExperiment(x, simplify, i = 1L, withDimnames = TRUE)
qreduceSummarizedExperiment(x, query, simplify, i = 1L, withDimnames = TRUE)
```

Arguments

Х	RaggedExperiment
i	integer(1), character(1), or logical() selecting the assay to be transformed.

withDimnames	logical(1) default TRUE. propagate dimnames to SummarizedExperiment.
simplify	function of 1 (for disjoinSummarizedExperiment) or 3 (for qreduceSummarizedExperiment) arguments, used to transform assays. See assay-functions.
query	GRanges provding regions over which reduction is to occur.

Value

All functions return RangedSummarizedExperiment.

sparseSummarizedExperiment has rowRanges() identical to the row ranges of x, and assay()
data as sparseAssay(). This is very space-inefficient representation of ragged data.

compactSummarizedExperiment has rowRanges() identical to the row ranges of x, and assay() data as compactAssay(). This is space-inefficient representation of ragged data when samples are primarily composed of different ranges.

disjoinSummarizedExperiment has rowRanges() identical to the disjoint row ranges of x, disjoint(rowRanges(x)),
and assay() data as disjoinAssay().

qreduceSummarizedExperiment has rowRanges() identical to query, and assay() data as qreduceAssay().

Examples

```
x <- RaggedExperiment(GRangesList(</pre>
    GRanges(c("A:1-5", "A:4-6", "A:10-15"), score=1:3),
GRanges(c("A:1-5", "B:1-3"), score=4:5)
))
## sparseSummarizedExperiment
sse <- sparseSummarizedExperiment(x)</pre>
assay(sse)
rowRanges(sse)
## compactSummarizedExperiment
cse <- compactSummarizedExperiment(x)</pre>
assay(cse)
rowRanges(cse)
## disjoinSummarizedExperiment
disjoinAssay(x, lengths)
dse <- disjoinSummarizedExperiment(x, lengths)</pre>
assay(dse)
rowRanges(dse)
## qreduceSummarizedExperiment
x <- RaggedExperiment(GRangesList(</pre>
    GRanges(c("A:1-3", "A:4-5", "A:10-15"), score=1:3),
    GRanges(c("A:4-5", "B:1-3"), score=4:5)
))
query <- GRanges(c("A:1-2", "A:4-5", "B:1-5"))</pre>
weightedmean <- function(scores, ranges, granges)</pre>
    ## weighted average score per query range
```

```
sum(scores * width(ranges)) / sum(width(ranges))
```

11

```
qreduceAssay(x, query, weightedmean)
qse <- qreduceSummarizedExperiment(x, query, weightedmean)
assay(qse)
rowRanges(qse)</pre>
```

Index

```
[,RaggedExperiment,ANY,ANY,ANY-method
        (RaggedExperiment-class), 5
assay, RaggedExperiment, ANY-method
        (RaggedExperiment-class), 5
assay,RaggedExperiment,missing-method
        (RaggedExperiment-class), 5
assay-functions, 2
assayNames,RaggedExperiment-method
        (RaggedExperiment-class), 5
assays, RaggedExperiment-method
        (RaggedExperiment-class), 5
coerce, GRangesList, RaggedExperiment-method
        (RaggedExperiment-class), 5
coerce,RaggedExperiment,GRangesList-method
        (RaggedExperiment-class), 5
colData,RaggedExperiment-method
        (RaggedExperiment-class), 5
colData<-,RaggedExperiment,DataFrame-method</pre>
        (RaggedExperiment-class), 5
compactAssay (assay-functions), 2
compactSummarizedExperiment
        (sparseSummarizedExperiment),
        10
```

```
DataFrame, 6, 8, 9

dim,RaggedExperiment-method

(RaggedExperiment-class), 5

dimnames,RaggedExperiment-method

(RaggedExperiment-class), 5

dimnames<-,RaggedExperiment,list-method

(RaggedExperiment-class), 5

disjoinAssay (assay-functions), 2

disjoinSummarizedExperiment

(sparseSummarizedExperiment),

10
```

GRanges, 8, 9
GRangesList, 2, 9
IntegerRanges, 7
IntegerRangesList, 7
IRanges, 7
IRangesList, 7

length,RaggedExperiment-method
 (RaggedExperiment-class), 5
mcols,RaggedExperiment-method
 (RaggedExperiment-class), 5
mcols<-,RaggedExperiment-method
 (RaggedExperiment-class), 5</pre>

```
RaggedExperiment, 2
RaggedExperiment
        (RaggedExperiment-class), 5
RaggedExperiment-class, 5
RaggedExperiment-package, 2
rowData,RaggedExperiment-method
        (RaggedExperiment-class), 5
rowData<-,RaggedExperiment-method</pre>
        (RaggedExperiment-class), 5
rowRanges, 2
rowRanges,RaggedExperiment-method
        (RaggedExperiment-class), 5
show,RaggedExperiment-method
        (RaggedExperiment-class), 5
SimpleList, 8, 9
sparseAssay (assay-functions), 2
sparseSummarizedExperiment, 10
subsetByOverlaps,RaggedExperiment,Vector-method
        (RaggedExperiment-class), 5
SummarizedExperiment, 2, 8
```

Views,7 ViewsList,7