Package 'msaR'

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Title Multiple Sequence Alignment for R Shiny
Version 0.6.0
Description Visualizes multiple sequence alignments dynamically within the Shiny web application framework.
Depends R (>= $3.2.2$)
License BSL-1.0
Imports ape, htmlwidgets
Suggests Biostrings, knitr, rmarkdown, testthat (>= 3.0.0)
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.check_lengths

check_lengths

Description

check_lengths

Usage

```
.check_lengths(seqlist)
```

Arguments

seqlist

optimistic loading can lead to issues. All MSAs should have identical length. if not theres a problem.

as.fasta

as.fasta

Description

functionality to convert objects to a fasta string. Currently this can handle character objects which are interpreted as filenames or several of the popular means of storing sequence data: DNAbin, DNAStringSet, AAStringSet, RNAStringSet, BStringSet, DNAMultipleAlignment, RNAMultipleAlignment, or AAMultipleAlignment.

Usage

```
as.fasta(seqs)
```

Arguments

seqs

(Required.) the sequence/alignment to be displayed. A character vector, DNAbin, DNAStringSet, AAStringSet, or RNAStringSet.

Value

A character string in fasta format.

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Examples

```
seqfile <- system.file("sequences","AHBA.aln",package="msaR")
as.fasta(seqfile)
help("as.fasta")

## Not run:
if (requireNamespace("Biostrings")) {
   seqs <- readDNAStringSet(seqfile)
   as.fasta(seqs)
}

## End(Not run)</pre>
```

as.sequences

as.sequences

Description

functionality to convert sequence objects into R lists that can be serialized to JS as JSON. Currently, this can handle character objects which are interpreted as filenames or several of the popular means of storing sequence data: DNAbin, DNAStringSet, AAStringSet, RNAStringSet, BStringSet, DNAMultipleAlignment, RNAMultipleAlignment, or AAMultipleAlignment.

Usage

```
as.sequences(seqs)
```

Arguments

seqs

(Required.) the sequence/alignment to be displayed. A character vector, DNAbin, DNAStringSet, AAStringSet, or RNAStringSet.

Value

A list of named lists where each sublist has name, id, and seq members.

Examples

```
seqfile <- system.file("sequences","AHBA.aln",package="msaR")
as.sequences(seqfile)
help("as.sequences")

## Not run:
if (requireNamespace("Biostrings")) {
   seqs <- readDNAStringSet(seqfile)
   as.sequences(seqs)
}

## End(Not run)</pre>
```

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msaR msaR

Description

Dynamic Multiple Sequence Alignments in R and Shiny

Usage

```
msaR(
 msa,
 menu = TRUE,
 width = NULL,
 height = NULL,
  rowheight = 15,
  alignmentHeight = 225,
 overviewbox = TRUE,
  seqlogo = TRUE,
  colorscheme = "nucleotide",
  conservation = FALSE,
 markers = TRUE,
 metacell = FALSE,
 leftheader = TRUE,
  labels = TRUE,
  labelname = TRUE,
  labelid = FALSE,
  labelNameLength = 100,
 overviewboxWidth = "auto",
  overviewboxHeight = "fixed"
)
```

Arguments

msa	File or BioString Object representing a multiple Sequence Alignment.
menu	Optional. Default TRUE. Determines whether to include the interactive menu.
width	Optional. Default NULL. The width of the html widget element.
height	Optional. Default NULL. The height of the html widget element.
rowheight Optional. Default 20. Height of a row in the MSA.	
alignmentHeigh	t
	Optional. Default 225. Height of the MSA.
overviewbox	optional. Default TRUE. Include the overview box?
seqlogo	optional. Default TRUE. Include the seqlogo?
colorscheme	optional. Default "nucleotide". The color scheme to use. Can be one of the following: "buried", "cinema", "clustal", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "cinema", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "cinema", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "cinema", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "cinema", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "clustal2", "helix", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "hydro", "hydro", "lesk", "mae", "nucleotide", "purine", "stransfer of the following: "buried", "hydro", "hydr
conservation	optional. Default TRUE. Include the conservation widget?

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markers optional. Default TRUE. Include the alignment markers? These are the numbers

along the top that

metacell optional. Default FALSE. Include the per-sequence metadata.

leftheader optional. Default TRUE. Include the header information.

labels optional. Default TRUE. Include all of the sequence information msa Labels.

labelname optional. Default TRUE. Include sequence name? labelid optional. Default FALSE. Include the labelid?

labelNameLength

optional. Default 100. Width of the Label Names.

overviewboxWidth

optional. Default. "auto". Can also be "fixed"

overviewboxHeight

optional. Default. "fixed". Can also be an integer value.

Examples

```
seqfile <- system.file("sequences","AHBA.aln",package="msaR")
msaR(seqfile)</pre>
```

msaROutput

Widget output function for use in Shiny

Description

Widget output function for use in Shiny

Usage

```
msaROutput(outputId, width = "100%", height = "100%")
```

Arguments

outputId output id width width height height

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	renderMsaR	Widget render function for use in Shiny
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Description

Widget render function for use in Shiny

Usage

```
renderMsaR(expr, env = parent.frame(), quoted = FALSE)
```

Arguments

```
expr expr env env quoted quoted
```

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