

Package ‘BgeeCall’

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Type Package

Title Automatic RNA-Seq present/absent gene expression calls generation

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Description BgeeCall allows to generate present/absent gene expression calls without using an arbitrary cutoff like TPM<1.

Calls are generated based on reference intergenic sequences. These sequences are generated based on expression of all RNA-Seq libraries of each species integrated in Bgee (<https://bgee.org>).

Depends R (>= 3.6)

Imports GenomicFeatures, rhdf5, tximport, Biostrings, rtracklayer, biomaRt, jsonlite, methods, grDevices, graphics, stats, utils

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URL <https://github.com/BgeeDB/BgeeCall>

BugReports <https://github.com/BgeeDB/BgeeCall/issues>

VignetteBuilder knitr

biocViews Software, GeneExpression, RNASEq

Suggests knitr, testthat, rmarkdown, AnnotationHub, httr

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AbundanceMetadata-class

AbundanceMetadata s4 class

Description

An S4 class that is the parent class of all abundance tool Classes. It contains information needed to all abundance tools. This class can be seen as an abstract class, you should never instanciate it.

Slots

txOut Similar to tximport txOut parameter. Allows to keep abundance at transcript level if TRUE (default = FALSE)
 ignoreTxVersion logical used to remove transcript version in transcript ID if TRUE (default = FALSE)

cutoff numeric corresponding to the proportion of intergenic present divided by proportion of protein coding present (default = 0.05). In the Bgee pipeline this cutoff is fixed and its value is 0.05. Be careful when changing this parameter as it could have a huge impact on your present/absent calls.

full_transcriptome_file Name of the fasta file containing both transcriptomic and intergenic regions. This file is created by the pipeline. You should edit this slot only if you already have such a file with a different name.

tx2gene_file Name of the file containing the mapping between transcript IDs and gene IDs (See the tximport package vignette for more details). This file is created by the pipeline. You should edit this slot only if you already have such a file with a different name. This file must be stored at get_species_path()

tx2gene_file_without_version Name of the file containing the mapping between transcript IDs and gene IDs if ignoreTxVersion == TRUE (See the tximport package vignette for more details). This file is created by the pipeline. You should edit this slot only if you already have such a file with a different name. This file must be stored at get_species_path()

gene2biotype_file Name of the file containing the mapping between gene IDs and biotypes. This file is created by the pipeline. You should edit this slot only if you already have such a file with a different name.

tool_name Name of the tool that will be used to generate transcript abundance estimation. All descendants of this class have to define a value for this slot (in the prototype section)

abundance_file Name of the transcript-level abundance file. All descendants of this class have to define a value for this slot (in the prototype section)

read_size_kmer_threshold read size of the library below which transcript index is created using a smaller kmer size

transcript_id_header Name of the header of the column that contains transcript ID

count_header Name of the header of the column that contains count

abundance_header Name of the header of the column that contains abundance

eff_length_header Name of the header of the column that contains effective length

transcript_calls_file_name default name of file containing all transcript ids and calls (if calls created at transcript level)

gene_calls_file_name default name of file containing all gene ids and calls (if calls created at gene level)

transcript_cutoff_file_name default name of file containing summary of cutoff used to generate transcript expression calls (if calls created at transcript level)

gene_cutoff_file_name default name of file containing summary of cutoff used to generate gene expression calls (if calls created at gene level)

transcript_distribution_file_name default name of density plot file containing TPM distribution of all transcripts (if calls created at transcript level)

gene_distribution_file_name default name of density plot file containing TPM distribution of all genes (if calls created at gene level)

BgeeMetadata-class *BgeeMetadata S4 class*

Description

An S4 class that contains all information to retrieve intergenic regions generated by Bgee.

Slots

intergenic_release Bgee intergenic release that will be used
 all_releases list of all reference intergenic releases that can be used to generate your present/absent expression calls.
 intergenic_prefix String used to generate an intergenic release specific output directory
 custom_intergenic_path path to a local version of reference intergenic fasta file. If NULL (by default) the reference intergenic fasta file will be downloaded. If not NULL BgeeCall will merge this local reference intergenic file with the transcriptome. Except if you generated your own intergenic regions always keep it NULL.

create_kallisto_index *Create kallisto indexes.*

Description

This function creates kallisto indexes. Two indexes can be created depending on the reads size (see ‘AbundanceMetadata@read_size_kmer_threshold’ and ‘UserMetadata@reads_size’ for more information). One with default kmer value (31 nt) and one with kmer size of 15 nt. In order to generate.

Usage

```
create_kallisto_index(  
  myKallistoMetadata,  
  myBgeeMetadata,  
  myUserMetadata,  
  transcriptome_path = ""  
)
```

Arguments

myKallistoMetadata
 A Reference Class KallistoMetadata object.
myBgeeMetadata A Reference Class BgeeMetadata object.
myUserMetadata A Reference Class UserMetadata object.
transcriptome_path
 path to the transcriptome fasta file. If no path is provided the default path created using BgeeCall will be used. IMPORTANT : in BgeeCall the transcriptome used to generate present/absent calls contains both intergenic sequences downloaded from Bgee and the reference transcriptome. If this function is run to generate present/absent then ‘transcriptome_path’ has to be empty

Value

create kallisto index and save it on the hard drive

Author(s)

Julien Wollbrett.

Examples

```
## Not run:
# first a transcriptome is needed. Here it is downloaded from AnnotationHub
library(AnnotationHub)
ah <- AnnotationHub()
ah_resources <- query(ah, c('Ensembl', 'Caenorhabditis elegans', '84'))

# kallisto can not deal with S4 objects. A Path to a transcriptome file is
# required
transcriptome_object <- rtracklayer::import.2bit(ah_resources[['AH50453']])
transcriptome_path <- file.path(getwd(), 'transcriptome.fa')
Biostrings::writeXStringSet(transcriptome_object, transcriptome_path)

# initialize objects needed to create destination folder
bgee <- new('BgeeMetadata')
user <- new('UserMetadata', species_id = '6239')
kallisto <- new('KallistoMetadata')

# generate transcriptome index
create_kallisto_index(kallisto, bgee, user, transcriptome_path)

## End(Not run)
```

download_fasta_intergenic

Download fasta intergenic

Description

Check if reference intergenic fasta file has already been downloaded. If not the file is downloaded from Bgee FTP or from the community repository depending on myBgeeMetadata@intergenic_release. if myBgeeMetadata@intergenic_release == "community" then reference intergenic wil be downloaded from the Zenodo community repository. Otherwise Reference intergenic sequences will be downloaded from the official Bgee FTP. Be careful when using reference intergenic sequences generated by the community as the Bgee team do not deeply review them.

Usage

```
download_fasta_intergenic(
  myBgeeMetadata = new("BgeeMetadata"),
  myUserMetadata,
  intergenic_file
)
```

Arguments

`myBgeeMetadata` A Reference Class BgeeMetadata object (optional)
`myUserMetadata` A Reference Class UserMetadata object.
`intergenic_file`
 path where intergenic file will be saved

Value

download fasta intergenic from Bgee FTP or from the Zenodo community and save it locally

Examples

```
{
  bgee_intergenic_file <- file.path(getwd(), 'intergenic.fasta')
  userMetadata <- new('UserMetadata', species_id = '7227')
}
```

`download_kallisto` *Download binary version of kallisto.*

Description

Check your OS and download correct binary version of kallisto.

Usage

```
download_kallisto(myKallistoMetadata, myUserMetadata)
```

Arguments

`myKallistoMetadata`
 A Reference Class KallistoMetadata object.
`myUserMetadata` A Reference Class UserMetadata object.

Value

save uncompressed executable of kallisto on the hard drive

Author(s)

Julien Wollbrett.

Examples

```
{
  kallisto <- new('KallistoMetadata')
  user <- new('UserMetadata')
  download_kallisto(kallisto, user)
}
```

generate_calls_workflow
generate present/absent calls

Description

Main function running the workflow allowing to generate present/absent calls from a file, a data.frame, or objects of the classe UserMetadata (please choose only 1 out of the 3) This workflow is highly tunable by editing default values of the slots of S4 objects. For more information on how to tune the workflow please have a look at the vignette and the documentation of the classes KallistoMetadata, AbundanceMetadata, UserMetadata, and BgeeMetadata

Usage

```
generate_calls_workflow(  
  abundanceMetadata = new("KallistoMetadata"),  
  bgeeMetadata = new("BgeeMetadata"),  
  userMetadata = NULL,  
  userDataFrame = NULL,  
  userFile = NULL  
)
```

Arguments

abundanceMetadata	A Reference Class BgeeMetadata object (optional) allowing to tune your gene quantification abundance analyze
bgeeMetadata	A Reference Class BgeeMetadata object (optional) allowing to choose the version of reference intergenic sequences
userMetadata	A Reference Class UserMetadata object (optional). generate present/allows calls using objects of the UserMetadata class. Can be one object or a list of objects.
userDataFrame	a data.frame comtaining all information to generate present/absent calls. Each line of this data.frame will generate calls for one RNA-Seq library. This data.frame must contains 7 columns : - species_id : The ensembl species ID - run_ids : (optional) allows to generate calls for a subpart of all runs of the library. must be a character or a list of characters. - reads_size (optional) the size of the reads of the library (Default = 50) if the reads size is lower than 51 abundance quantification wil be run with a smaller kmer size - rnaseq_lib_path : path to RNA-Seq library directory - transcriptome_path : path to transcriptome file - annotation_path : path to annotation file - working_path : root of the directory where results will be written
userFile	path to a tsv file containing 7 columns. these columns are the same than for userDataFrame (see above). a template of this file is available at the root of the package and accessible with the command system.file('userMetadataTemplate.tsv', package = 'BgeeCall')

Value

paths to the 4 results files (see vignette for more details)

Author(s)

Julien Wollbrett

See Also

[AbundanceMetadata](#), [KallistoMetadata](#), [BgeeMetadata](#), [UserMetadata](#)

Examples

```
## Not run:
# import gene annotation and transcriptome from AnnotationHub
library(AnnotationHub)
ah <- AnnotationHub()
ah_resources <- query(ah, c('Ensembl', 'Caenorhabditis elegans', '84'))
annotation_object <- ah_resources[['AH50789']]
transcriptome_object <- rtracklayer::import.2bit(ah_resources[['AH50453']])

# instantiate BgeeCall object
# add annotation and transcriptome in the user_BgeeCall object
# it is possible to import them using an S4 object (GRanges, DNAStringSet)
# or a file (gtf, fasta) with methods setAnnotationFromFile() and
# setTranscriptomeFromFile()
user_BgeeCall <- setAnnotationFromObject(user_BgeeCall,
                                             annotation_object,
                                             'WBcel235_84')
user_BgeeCall <- setTranscriptomeFromObject(user_BgeeCall,
                                              transcriptome_object,
                                              'WBcel235')
# provide path to the directory of your RNA-Seq library
user_BgeeCall <- setRNASeqLibPath(user_BgeeCall,
                                   system.file('extdata', 'SRX099901_subset',
                                   package = 'BgeeCall'))

# run the full BgeeCall workflow
calls_output <- generate_calls_workflow(
    userMetadata = user_BgeeCall)

## End(Not run)
```

generate_presence_absence

Generate presence absence

Description

Generate presence absence calls. It correponds to the last part of the generation of the expression calls workflow. It runs the last part of the workflow generating present/absent expression calls. This function should only be used by advanced user who already manually run all previous parts of the pipeline. If you are not an advanced user it is safer to run the function “generate_calls_workflow“ that run all steps of the workflow

Usage

```
generate_presence_absence(
  myAbundanceMetadata = new("KallistoMetadata"),
  myBgeeMetadata = new("BgeeMetadata"),
  myUserMetadata
)
```

Arguments

`myAbundanceMetadata`
A descendant object of the Class `myAbundanceMetadata` (optional).

`myBgeeMetadata` A Class `BgeeMetadata` object (optional).

`myUserMetadata` A Class `UserMetadata` object.

Value

path to the 4 output files

Author(s)

Julien Wollbrett

Julien Roux

See Also

`generate_calls_workflow`

Examples

```
{
# this example reuse data present in the directory 'extdata' of the package.
user <- new('UserMetadata', working_path = system.file('extdata',
package = 'BgeeCall'), species_id = '6239', rnaseq_lib_path = system.file(
'extdata', 'SRX099901_subset', package = 'BgeeCall'),
annotation_name = 'WBcel235_84', simple_arborescence = TRUE)
calls_output <- generate_presence_absence(myUserMetadata = user)

#
```

`getIntergenicPrefix` ‘`intergenic_prefix`‘ Getter

Description

Get value of the ‘`intergenic_prefix`‘ slot

Usage

```
getIntergenicPrefix(bgeeObject)

## S4 method for signature 'BgeeMetadata'
getIntergenicPrefix(bgeeObject)
```

Arguments

`bgeeObject` The `BgeeMetadata` object

Value

the value of the ‘`intergenic_prefix`‘ slot of the object

Examples

```
{
bgee <- new("BgeeMetadata")
intergenic_prefix <- getIntergenicPrefix(bgee)
}
```

`getIntergenicRelease` ‘`intergenic_release`‘ Getter

Description

Get value of the ‘`intergenic_release`‘ slot

Usage

```
getIntergenicRelease(bgeeObject)

## S4 method for signature 'BgeeMetadata'
getIntergenicRelease(bgeeObject)
```

Arguments

`bgeeObject` The `BgeeMetadata` object

Value

the value of the ‘`intergenic_release`‘ slot of the object

Examples

```
{
bgee <- new("BgeeMetadata")
intergenic_release <- getIntergenicRelease(bgee)
}
```

```
getRunIds           'run_ids' Getter
```

Description

Get value of the ‘run_ids‘ slot

Usage

```
getRunIds(userObject)  
## S4 method for signature 'UserMetadata'  
getRunIds(userObject)
```

Arguments

userObject The UserMetadata object

Value

the value of the ‘run_ids‘ slot of the object

Examples

```
{  
  user <- new("UserMetadata")  
  run_ids <- getRunIds(user)  
}
```

```
getSimpleArborescence 'simple_arborescence' Getter
```

Description

Get value of the ‘simple_arborescence‘ slot

Usage

```
getSimpleArborescence(userObject)  
## S4 method for signature 'UserMetadata'  
getSimpleArborescence(userObject)
```

Arguments

userObject The UserMetadata object

Value

the value of the ‘simple_arborescence‘ slot of the object

Examples

```
{
  user <- new("UserMetadata")
  simple_arborescence <- getSimpleArborescence(user)
}
```

getWorkingPath	‘working_path’ Getter
----------------	-----------------------

Description

Get value of the ‘working_path’ slot

Usage

```
getWorkingPath(userObject)

## S4 method for signature 'UserMetadata'
getWorkingPath(userObject)
```

Arguments

userObject The UserMetadata object

Value

the value of the ‘working_path’ slot of the object

Examples

```
{
  user <- new("UserMetadata")
  working_path <- getWorkingPath(user)
}
```

KallistoMetadata-class	<i>KallistoMetadata S4 class</i>
------------------------	----------------------------------

Description

An S4 class that is the descendant of the AbundanceMetadata class. It contains all metadata needed to run kallisto analysis. All slots of this class have a default value. You do not need to edit them to run the package

Slots

`download_kallisto` A logical allowing to use an already installed version of kallisto or to download a version that will be used only by this package

`kallisto_windows_url` URL to the binary of kallisto for windows

`kallisto_linux_url` URL to the binary of kallisto for linux

`kallisto_osx_url` URL to the binary of kallisto for MacOS

`kallisto_windows_dir` Name of the directory where kallisto will be installed on windows

`kallisto_linux_dir` Name of the directory where kallisto will be installed on linux

`kallisto_osx_dir` Name of the directory where kallisto will be installed on Mac

`unix_kallisto_name` Name of the kallisto executable in linux and macOS

`windows_kallisto_name` Name of the kallisto executable in windows

`index_file` Name of index file generated by kallisto with default kmer size. It will be generated using the fasta file that contains both transcriptomic and intergenic regions. Do not use an index you generated outside of this package. This file is created by the pipeline. You should edit this slot only if you already have such a file with a different name. This file must be stored at `get_tool_path()`

`k15_index_file` same as `index_file`. This index is generated with smallest kmers and will be used only for libraries containing reads smaller than 50nt.

`single_end_parameters` kallisto parameters used to run a single end mapping

`pair_end_parameters` kallisto parameters used to run a pair end mapping

`overwrite_index` logical allowing to overwrite already existing index. FALSE by default. Then by default already existing index files will not be generated again.

`overwrite_quant` logical allowing to overwrite already existing abundance.txt files. FALSE by default. Then by default already existing quantification files will not be generated again.

`overwrite_calls` logical allowing to overwrite already existing present/absent calls. FALSE by default. Then by default already generated calls will not be generated again.

list_bgee_ref_intergenic_species

List species having Bgee reference intergenic sequences

Description

Return information related to species having Bgee reference intergenic sequences available for the selected Bgee intergenic release:

- `speciesId` : the NCBI species ID of the species
- `specieName` : scientific species name
- `numberOfLibraries` : number of libraries used to generate these reference intergenic sequences

If a `BgeeMetadata` object is provided this function retrieves the list of species using `BgeeMetadata@intergenic_release`. If only a ‘release’ is provided it will use it to retrieve the list of species. If none of them are provided the default Bgee reference intergenic release will be used.

Usage

```
list_bgee_ref_intergenic_species(myBgeeMetadata = NULL, release = NULL)
```

Arguments

<code>myBgeeMetadata</code>	A Reference Class BgeeMetadata object
<code>release</code>	A Bgee reference intergenic release name

Value

list all species having reference intergenic sequences available in the selected release

Author(s)

Julien Wollbrett

Examples

```
{
  bgee <- new("BgeeMetadata")
  list_bgee_ref_intergenic_species(myBgeeMetadata = bgee)
  list_bgee_ref_intergenic_species(release = '0.2')
}
```

list_community_ref_intergenic_species

List species having reference intergenic sequences created by the BgeeCall community

Description

Return information related to species having reference intergenic sequences created by the BgeeCall community - speciesId : the NCBI species ID of the species - url : url to the reference intergenic fasta file - numberOfLibraries : number of libraries used to generate these reference intergenic sequences

Usage

```
list_community_ref_intergenic_species()
```

Value

list all species having reference intergenic sequences created by the community

Author(s)

Julien Wollbrett

Examples

```
{
  list_community_ref_intergenic_species()
}
```

list_intergenic_release

List reference intergenic releases usable with the BgeeCall package

Description

Returns information on available Bgee intergenic releases, the access URL for FTP, and the date of release

Usage

```
list_intergenic_release(release = NULL)
```

Arguments

release	A character specifying a targeted release number (e.g., '0.1'). If not specified, all available releases are shown.
---------	---

Value

A data frame with information on Bgee intergenic releases available to use with the BgeeCall package.

Author(s)

Julien Wollbrett

Examples

```
{  
  list_intergenic_release()  
}
```

merge_transcriptome_and_intergenic

Merge transcriptome file provided by the user with the Bgee intergenic fasta file.

Description

This function will create a file corresponding to the concatenation of the transcriptome fasta file provided by the user and the corresponding intergenic fasta file created by Bgee.

Usage

```
merge_transcriptome_and_intergenic(  
  myKallistoMetadata,  
  myBgeeMetadata,  
  myUserMetadata  
)
```

Arguments

`myKallistoMetadata` A Reference Class KallistoMetadata object.
`myBgeeMetadata` A Reference Class BgeeMetadata object.
`myUserMetadata` A Reference Class UserMetadata object.

Value

save merged file on the hard drive

Author(s)

Julien Wollbrett.

Examples

```
{
bgee <- new('BgeeMetadata', intergenic_release = '0.1')
user <- new ('UserMetadata', species_id = '6239')
kallisto <- new('KallistoMetadata')
user <- setTranscriptomeFromFile(user, system.file("extdata",
"transcriptome.fa", package = "BgeeCall"), 'WBcel235')
merge_transcriptome_and_intergenic(kallisto, bgee, user)
}
```

run_kallisto

Run one kallisto abundance analyse

Description

Run kallisto and all preliminary steps if needed like : - creation of transcriptome with intergenic (if needed) - installation of kallisto (if needed) - index creation (if needed) - run kallisto quantification

Usage

```
run_kallisto(
  myKallistoMetadata,
  myBgeeMetadata,
  myUserMetadata,
  transcriptome_path = ""
)
```

Arguments

`myKallistoMetadata` A Reference Class KallistoMetadata object.
`myBgeeMetadata` A Reference Class BgeeMetadata object.
`myUserMetadata` A Reference Class UserMetadata object. This object has to be edited before running kallisto @seealso UserMetadata.R

transcriptome_path

path to the transcriptome fasta file. If no path is provided the default path created using BgeeCall will be used. IMPORTANT : in BgeeCall the transcriptome used to generate present/absent calls contains both intergenic sequences downloaded from Bgee and the reference transcriptome.

Value

create kallisto output files and save them on the hard drive

Author(s)

Julien Wollbrett.

Examples

```
## Not run:
# first a transcriptome is needed. Here it is downloaded from AnnotationHub
library(AnnotationHub)
ah <- AnnotationHub()
ah_resources <- query(ah, c('Ensembl', 'Caenorhabditis elegans', '84'))

# kallisto can not deal with S4 objects. Path to transcriptome file is
# required
transcriptome_object <- rtracklayer::import.2bit(ah_resources[['AH50453']])
transcriptome_path <- file.path(getwd(), 'transcriptome.fa')
Biostrings::writeXStringSet(transcriptome_object, transcriptome_path)

# initialize objects needed to create destination folder
bgee <- new('BgeeMetadata')
user <- new('UserMetadata', species_id = '6239')
user <- setRNASEqLibPath(user, system.file(
  'extdata', 'SRX099901_subset',
  package = 'BgeeCall'))
kallisto <- new('KallistoMetadata')

# generate transcriptome index
run_kallisto(kallisto, bgee, user, transcriptome_path)

## End(Not run)
```

Description

Run tximport. Will summarize abundance estimation from transcript level to gene level if ‘myAbundanceMetadata@txout == FALSE’. Otherwise keep abundance estimation at transcript level.

Usage

```
run_tximport(
  myAbundanceMetadata = new("KallistoMetadata"),
  myBgeeMetadata = new("BgeeMetadata"),
  myUserMetadata,
  abundanceFile = ""
)
```

Arguments

myAbundanceMetadata A descendant object of the Class myAbundanceMetadata.

myBgeeMetadata A Reference Class BgeeMetadata object.

myUserMetadata A Reference Class UserMetadata object.

abundanceFile (Optional) Path to the abundance file. NULL by default. If not NULL, the file located at ‘abundanceFile’ will be used to run tximport. Otherwise (Default) the path to the abundance file is deduced from attributes of classes ‘BgeeMetadata’, ‘UserMetadata’ and ‘AbundanceMetadata’

Value

a tximport object

Author(s)

Julien Wollbrett

Examples

```
{
  user <- new("UserMetadata", working_path = system.file("extdata",
    package = "BgeeCall"), species_id = "6239",
    rnaseq_lib_path = system.file("extdata",
      "SRX099901_subset", package = "BgeeCall"),
    annotation_name = "WBcel235_84", simple_arborescence = TRUE)
  abundance_file <- system.file('extdata', 'abundance.tsv', package = 'BgeeCall')
  tx_import <- run_tximport(myUserMetadata = user,
    abundanceFile = abundance_file)
}
```

setAnnotationFromFile *Set annotation_object of one UserMetadata object*

Description

Method of the class UserMetadata. Set annotation_object of one UserMetadata object by providing the path to a fasta transcriptome file.

Usage

```
setAnnotationFromFile(userObject, annotationPath, annotationName)

## S4 method for signature 'UserMetadata,character,missing'
setAnnotationFromFile(userObject, annotationPath, annotationName)

## S4 method for signature 'UserMetadata,character,character'
setAnnotationFromFile(userObject, annotationPath, annotationName)
```

Arguments

`userObject` The UserMetadata object
`annotationPath` Absolute path to the annotation file
`annotationName` (optional) Name of the annotation. Will be used to create folders.

Details

If no annotationName is provided the name of the annotation file will be used to create folders.

Value

An object of the class UserMetadata

Examples

```
{
# path to gtf annotation file
annotation_file <- system.file("extdata", "annotation.gtf", package = "BgeeCall")
user <- new("UserMetadata")
user <- setAnnotationFromFile(user, annotation_file,
                             "annotation_name")
}
```

setAnnotationFromObject

Set annotation_object of one UserMetadata object

Description

Method of the class UserMetadata. Set annotation_object of one UserMetadata object by using one GRanges object as input.

Usage

```
setAnnotationFromObject(userObject, annotationObject, annotationName)

## S4 method for signature 'UserMetadata,GRanges,character'
setAnnotationFromObject(userObject, annotationObject, annotationName = "")
```

Arguments

`userObject` The `UserMetadata` object
`annotationObject` object of the `GRanges` S4 class
`annotationName` (optional) Name of the annotation. Will be used to create folders.

Details

If no `annotationName` is provided the name of the file is used to create folders.

Value

An object of the class `UserMetadata`

Examples

```
{
  user <- new("UserMetadata")
  annotation_object <- rtracklayer::import(system.file("extdata",
  "annotation.gtf", package = "BgeeCall"))
  user <- setAnnotationFromObject(user, annotation_object,
  "annotation_name")
}
```

`setIntergenicRelease` ‘`intergenic_release`’ Setter

Description

Set value of the ‘`intergenic_release`’ slot

Usage

```
setIntergenicRelease(bgeeObject, intergenicRelease)

## S4 method for signature 'BgeeMetadata,character'
setIntergenicRelease(bgeeObject, intergenicRelease)
```

Arguments

`bgeeObject` The `BgeeMetadata` object
`intergenicRelease` character corresponding to the ‘`intergenic_release`’

Value

An object of the class `BgeeMetadata` with new ‘`intergenic_release`’ value

Examples

```
{
  bgee <- new("BgeeMetadata")
  bgee <- setIntergenicRelease(bgee, "0.1")
}
```

`setOutputDir` ‘*output_dir*‘ *Setter*

Description

Set value of the ‘*output_dir*‘ slot

Usage

```
setOutputDir(userObject, outputDir)

## S4 method for signature 'UserMetadata,character'
setOutputDir(userObject, outputDir)
```

Arguments

<code>userObject</code>	The <code>UserMetadata</code> object
<code>outputDir</code>	path to the directory wanted as ‘ <i>output_dir</i> ‘

Value

An object of the class `UserMetadata` with new ‘*output_dir*‘ value

Examples

```
{
  user <- new("UserMetadata")
  user <- setOutputDir(user, getwd())
}
```

`setRNASEqLibPath` ‘*rnaseq_lib_path*‘ *Setter*

Description

Set value of the ‘*rnaseq_lib_path*‘ slot

Usage

```
setRNASEqLibPath(userObject, rnaSeqLibPath)

## S4 method for signature 'UserMetadata,character'
setRNASEqLibPath(userObject, rnaSeqLibPath)
```

Arguments

`userObject` The `UserMetadata` object
`rnaSeqLibPath` path to the directory wanted as ‘rnaseq_lib_path’

Value

An object of the class `UserMetadata` with new ‘rnaseq_lib_path’ value

Examples

```
{
  user <- new("UserMetadata")
  user <- setRNASEqLibPath(user, getwd())
}
```

`setRunIds` ‘run_ids’ Setter

Description

Method of the class `UserMetadata`. Set `run_ids` of one `UserMetadata` object by providing the id of all wanted runs

Usage

```
setRunIds(userObject, runIds)

## S4 method for signature 'UserMetadata,character'
setRunIds(userObject, runIds)
```

Arguments

`userObject` The `UserMetadata` object
`runIds` id of all wanted runs

Value

An object of the class `UserMetadata`

Examples

```
{
  user <- new("UserMetadata")
  user <- setRunIds(user, c("RUN_1", "RUN_2"))
}
```

`setSimpleArborescence` ‘*simple_arborescence*‘ *Setter*

Description

Set value of the ‘simple_arborescence’ slot

Usage

```
setSimpleArborescence(userObject, simpleArborescence)
```

```
## S4 method for signature 'UserMetadata,logical'
setSimpleArborescence(userObject, simpleArborescence)
```

Arguments

<code>userObject</code>	The UserMetadata object
<code>simpleArborescence</code>	boolean defining if output files will be created a simple arborescence (TRUE) or not (FALSE)

Value

An object of the class UserMetadata with new ‘simple_arborescence‘ value

Examples

```
{
  user <- new("UserMetadata")
  user <- setSimpleArborescence(user, FALSE)
}
```

`setTranscriptomeFromFile`

Set transcriptome_object of one UserMetadata object

Description

Method of the class UserMetadata. Set transcriptome_object of one UserMetadata object by providing the path to a fasta transcriptome file.

Usage

```
setTranscriptomeFromFile(userObject, transcriptomePath, transcriptomeName)
```

```
## S4 method for signature 'UserMetadata,character,missing'
setTranscriptomeFromFile(userObject, transcriptomePath, transcriptomeName)
```

```
## S4 method for signature 'UserMetadata,character,character'
setTranscriptomeFromFile(userObject, transcriptomePath, transcriptomeName)
```

Arguments

userObject The UserMetadata object
transcriptomePath Absolute path to the transcriptome file
transcriptomeName (optional) Name of the transcriptome. Will be used to create folders.

Details

If no transcriptomeName is provided the name of the transcriptome file will be used to create folders.

Value

An object of the class UserMetadata

Examples

```
{
  transcriptome_path <- system.file("extdata", "transcriptome.fa", package = "BgeeCall")
  user <- new("UserMetadata")
  user <- setTranscriptomeFromFile(user, transcriptome_path,
                                    "transcriptome_name")
}
```

setTranscriptomeFromObject

Set transcriptome_object of one UserMetadata object

Description

Method of the class UserMetadata. Set transcriptome_object of one UserMetadata object by using one DNAStringSet object as input.

Usage

```
setTranscriptomeFromObject(userObject, transcriptomeObject, transcriptomeName)

## S4 method for signature 'UserMetadata,DNAStringSet,character'
setTranscriptomeFromObject(userObject, transcriptomeObject, transcriptomeName)
```

Arguments

userObject UserMetadata object
transcriptomeObject Object of the DNAStringSet S4 class
transcriptomeName Name of the transcriptome. Will be used to create transcriptome folders.

Details

Please use a DNAStringSet object as input. This class is defined in the Biostrings package

Value

an object of UserMetadata

Examples

```
{
  user <- new("UserMetadata")
  transcriptome_object <- Biostrings::readDNAStringSet(
    system.file("extdata", "transcriptome.fa", package = "BgeeCall"))
  user <- setTranscriptomeFromObject(user,
    transcriptome_object,
    "transcriptome_name")
}
```

`setWorkingPath` ‘`working_path`‘ Setter

Description

Set value of the ‘`working_path`‘ slot

Usage

```
setWorkingPath(userObject, workingPath)

## S4 method for signature 'UserMetadata,character'
setWorkingPath(userObject, workingPath)
```

Arguments

<code>userObject</code>	The <code>UserMetadata</code> object
<code>workingPath</code>	path to the directory wanted as ‘ <code>working_path</code> ‘

Value

An object of the class `UserMetadata` with new ‘`working_path`‘ value

Examples

```
{
  user <- new("UserMetadata")
  user <- setWorkingPath(user, getwd())
}
```

UserMetadata-class *UserMetadata S4 class*

Description

An S4 class containing all metadata that have to be provided by the user. It is mandatory to edit ‘species_id’, ‘rnaseq_lib_path’, ‘transcriptome_path’, ‘annotation_name’, ‘annotation_object’ and potentially ‘run_ids’ before using the package.

Slots

species_id The NCBI Taxon Id of the species

run_ids A vector of character. Has to be provided only if a subset of runs present in UserMetadata@rnaseq_lib_path has to be run. If empty, all fastq files present in the rnaseq_lib_path will be considered as technical replicates and merged to run one transcript expression estimation analyse.

reads_size The size of the reads. If smaller than ‘KallistoMetadata@read_size_kmer_threshold’, an index with a kmer size of 15 bp will be used.

rnaseq_lib_path Path to the directory of the RNA-Seq library that contains fastq files. The extension of the fastq files name must be .fq, .fastq, .fq.gz, or .fastq.gz

transcriptome_name Name of the transcriptome used to generate arborescence of output repositories.

transcriptome_object Object containing transcriptome

annotation_name Name of the annotation used to generate arborescence of output repositories.

annotation_object Object containing annotations from GTF or GFF file

working_path Working directory. By default the working directory is defined with the ‘getwd()’ function.

simple_arborescence logical allowing to create a simple arborescence of directory. If ‘TRUE’ (default), all results will be on the same directory (working_path/intergenic_release/all_results/libraryId). Use ‘FALSE’ if you plan to generate expression calls for the same library using different transcriptomes or gene annotations, otherwise you will overwrite previous results. When ‘FALSE’ the path to result folder looks like : working_path/intergenic_release/speciesId/kallisto/transcriptome_name/annotation_name

output_dir (optional) Allows to manually define your output directory. By default the path to output directory is created automatically from the working_path (working_path/intergenic_release/all_results/libraryId)

verbose logical allowing to use the verbose mode. TRUE by default.

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