

Package: metabinR (via r-universe)

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Type Package

Title Abundance and Compositional Based Binning of Metagenomes

Version 2.0.0

biocViews Classification, Clustering, Microbiome, Sequencing, Software

Description Provide functions for performing abundance and compositional based binning on metagenomic samples, directly from FASTA or FASTQ files. Functions are implemented in Java and called via rJava. Parallel implementation that operates directly on input FASTA/FASTQ files for fast execution. Inputs may be file paths or Biostrings/ShortRead sequence objects; results are returned as a MetabinResult S4 object wrapping cluster assignments, algorithm parameters, and input metadata.

License GPL-3

Encoding UTF-8

Language en-US

LazyData false

Depends R (>= 4.4), methods

Imports BiocGenerics, BiocParallel, Biostrings, checkmate, cli, rJava, S4Vectors, ShortRead, utils

SystemRequirements Java (>= 11)

RoxygenNote 7.3.3

URL <https://github.com/gkanogiannis/metabinR>

BugReports <https://github.com/gkanogiannis/metabinR/issues>

Suggests BiocStyle, covr, cvms, data.table, dplyr, ggplot2, gridExtra, knitr, R.utils, rmarkdown, sabre, spelling, testthat (>= 3.0.0)

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abundance_based_binning

Abundance based binning on metagenomic samples

Description

This function performs abundance based binning on metagenomic samples, directly from FASTA or FASTQ files, by long kmer analysis ($k > 8$). See [doi:10.1186/s1285901611863](https://doi.org/10.1186/s1285901611863) for more details.

Usage

```
abundance_based_binning(
  ...,
  eMin = 1,
  eMax = 0,
  kMerSizeAB = 10,
  numOfClustersAB = 3,
  outputAB = "AB.cluster",
  keepQuality = FALSE,
  dryRun = FALSE,
  gzip = FALSE,
  numOfThreads = BiocParallel::bpworkers()
)
```

Arguments

... Input sequences. Either character paths to FASTA/FASTQ files (uncompressed or gzip compressed), a [DNAStrngSet](#) / [QualityScaledDNAStrngSet](#), or a [ShortReadQ](#) object. Non-file inputs are staged to a temporary FASTA/FASTQ file for the Java backend.

eMin	Exclude kmers of less or equal count.
eMax	Exclude kmers of more or equal count.
kMerSizeAB	kmer length for Abundance based Binning.
numOfClustersAB	Number of Clusters for Abundance based Binning.
outputAB	Output Abundance based Binning Clusters files location and prefix.
keepQuality	Keep fastq qualities on the output files. (will produce .fastq)
dryRun	Don't write any output files.
gzip	Gzip output files.
numOfThreads	Number of threads to use. Defaults to <code>bpworkers()</code> .

Value

A `MetabinResult` object. Its assignments slot is a `DataFrame` with `numOfClustersAB + 2` columns:

- `read_id`: read identifier from fasta header
- `AB`: read was assigned to this AB cluster index
- `AB.n`: read to cluster AB.n distance

For backwards-compatible data.frame output use `as.data.frame(result)`.

Author(s)

Anestis Gkanogiannis, <anestis@gkanogiannis.com>

References

<https://github.com/gkanogiannis/metabinR>

Examples

```
res <- abundance_based_binning(  
  system.file("extdata", "reads.metagenome.fasta.gz", package = "metabinR"),  
  dryRun = TRUE, kMerSizeAB = 8  
)  
res  
head(as.data.frame(res))
```

```
as.data.frame, MetabinResult-method
```

Convert a [MetabinResult](#) to a data.frame

Description

Preserves backwards compatibility with the data.frame return of metabinR <= 1.x.

Usage

```
## S4 method for signature 'MetabinResult'
as.data.frame(x, row.names = NULL, optional = FALSE, ...)
```

Arguments

```
x          A MetabinResult.
row.names, optional, ...
           Unused; retained for S3 signature compatibility.
```

Value

A base data.frame of the assignments.

```
composition_based_binning
```

Composition based binning on metagenomic samples

Description

This function performs composition based binning on metagenomic samples, directly from FASTA or FASTQ files, by short kmer analysis ($k < 8$). See [doi:10.1186/s1285901611863](https://doi.org/10.1186/s1285901611863) for more details.

Usage

```
composition_based_binning(
  ...,
  kMerSizeCB = 4,
  numOfClustersCB = 5,
  outputCB = "CB.cluster",
  keepQuality = FALSE,
  dryRun = FALSE,
  gzip = FALSE,
  numOfThreads = BiocParallel::bpworkers()
)
```

Arguments

...	Input sequences. Either character paths to FASTA/FASTQ files (uncompressed or gzip compressed), a DNAStrngSet / QualityScaledDNAStrngSet , or a ShortReadQ object. Non-file inputs are staged to a temporary FASTA/FASTQ file for the Java backend.
kMerSizeCB	kmer length for Composition based Binning.
numOfClustersCB	Number of Clusters for Composition based Binning.
outputCB	Output Composition based Binning Clusters files location and prefix.
keepQuality	Keep fastq qualities on the output files. (will produce .fastq)
dryRun	Don't write any output files.
gzip	Gzip output files.
numOfThreads	Number of threads to use. Defaults to bpworkers() .

Value

A [MetabinResult](#) object. Its assignments slot is a [DataFrame](#) with numofClustersCB + 2 columns:

- read_id : read identifier from fasta header
- CB : read was assigned to this CB cluster index
- CB.n : read to cluster CB.n distance

For backwards-compatible data.frame output use `as.data.frame(result)`.

Author(s)

Anestis Gkanogiannis, <anestis@gkanogiannis.com>

References

<https://github.com/gkanogiannis/metabinR>

Examples

```
res <- composition_based_binning(
  system.file("extdata", "reads.metagenome.fasta.gz", package = "metabinR"),
  dryRun = TRUE, kMerSizeCB = 2
)
res
```

hierarchical_binning *Hierarchical (ABxCB) binning on metagenomic samples*

Description

This function performs hierarchical binning on metagenomic samples, directly from FASTA or FASTQ files. First it analyzes sequences by long kmer analysis ($k > 8$), as in [abundance_based_binning](#). Then for each AB bin, it guesses the number of composition bins in it and performs composition based binning by short kmer analysis ($k < 8$), as in [composition_based_binning](#). See [doi:10.1186/s1285901611863](https://doi.org/10.1186/s1285901611863) for more details.

Usage

```
hierarchical_binning(
  ...,
  eMin = 1,
  eMax = 0,
  kMerSizeAB = 10,
  kMerSizeCB = 4,
  genomeSize = 3e+06,
  numOfClustersAB = 3,
  outputC = "ABxCB.cluster",
  keepQuality = FALSE,
  dryRun = FALSE,
  gzip = FALSE,
  numOfThreads = BiocParallel::bpworkers()
)
```

Arguments

...	Input sequences. Either character paths to FASTA/FASTQ files (uncompressed or gzip compressed), a DNAStrngSet / QualityScaledDNAStrngSet , or a ShortReadQ object. Non-file inputs are staged to a temporary FASTA/FASTQ file for the Java backend.
eMin	Exclude kmers of less or equal count.
eMax	Exclude kmers of more or equal count.
kMerSizeAB	kmer length for Abundance based Binning.
kMerSizeCB	kmer length for Composition based Binning.
genomeSize	Average genome size of taxa in the metagenome data.
numOfClustersAB	Number of Clusters for Abundance based Binning.
outputC	Output Hierarchical Binning (ABxCB) Clusters files location and prefix.
keepQuality	Keep fastq qualities on the output files. (will produce .fastq)
dryRun	Don't write any output files.
gzip	Gzip output files.
numOfThreads	Number of threads to use. Defaults to bpworkers() .

Value

A `MetabinResult` object. Its assignments slot is a `DataFrame`:

- `read_id` : read identifier from fasta header
- `ABxCB` : read was assigned to this ABxCB cluster index
- `ABxCB.n` : read to cluster ABxCB.n distance

For backwards-compatible `data.frame` output use `as.data.frame(result)`.

Author(s)

Anestis Gkanogiannis, <anestis@gkanogiannis.com>

References

<https://github.com/gkanogiannis/metabinR>

Examples

```
res <- hierarchical_binning(  
  system.file("extdata", "reads.metagenome.fasta.gz", package = "metabinR"),  
  dryRun = TRUE, kMerSizeAB = 4, kMerSizeCB = 2  
)  
res
```

`metabinR_jvm_options` *JVM options used when metabinR loads*

Description

Returns the vector of JVM flags passed to `.jpackage` on package load. Set `options(metabinR.jvm.flags = c(...))` before loading the package to override; set `options(java.parameters = ...)` to prepend heap-size flags (e.g. `"-Xmx4g"`) in the usual rJava way.

Usage

```
metabinR_jvm_options()
```

Value

A character vector of JVM flags.

Examples

```
metabinR_jvm_options()
```

MetabinResult-accessors

Accessors for [MetabinResult](#)

Description

Accessors for [MetabinResult](#)

Usage

```
assignments(x)

nClusters(x)

parameters(x)

algorithm(x)

## S4 method for signature 'MetabinResult'
assignments(x)

## S4 method for signature 'MetabinResult'
parameters(x)

## S4 method for signature 'MetabinResult'
algorithm(x)

## S4 method for signature 'MetabinResult'
nClusters(x)
```

Arguments

x A [MetabinResult](#) object.

Value

`assignments()` returns a [DataFrame](#) with the per-read cluster assignments and distances. `nClusters()` returns an integer scalar: the number of clusters inferred by the algorithm. `parameters()` returns the list of arguments passed to the binning function. `algorithm()` returns the algorithm tag ("AB", "CB", or "ABxCB").

Examples

```
res <- abundance_based_binning(
  system.file("extdata", "reads.metagenome.fasta.gz", package = "metabinR"),
  dryRun = TRUE, kMerSizeAB = 4, numOfClustersAB = 2
)
assignments(res)
```

```
nClusters(res)
algorithm(res)
parameters(res)$kMerSizeAB
```

MetabinResult-class *MetabinResult: binning result container*

Description

An S4 class returned by [`abundance_based_binning()`], [`composition_based_binning()`] and [`hierarchical_binning()`].

Usage

```
## S4 method for signature 'MetabinResult'
show(object)
```

Arguments

object A [MetabinResult](#).

Value

Objects of this class are returned by the binning functions. Use [assignments](#), [nClusters](#), [parameters](#), [algorithm](#), or `as.data.frame()` to access the results.

Slots

`assignments` A [DataFrame](#) of cluster assignments. The first column is `read_id`; subsequent columns are algorithm-specific (see the corresponding binning function).

`parameters` Named list of the parameters passed to the algorithm.

`inputs` Character vector of input file paths that were processed.

`algorithm` Character scalar: one of "AB", "CB", "ABxCB".

Examples

```
res <- abundance_based_binning(
  system.file("extdata", "reads.metagenome.fasta.gz", package = "metabinR"),
  dryRun = TRUE, kMerSizeAB = 4, numOfClustersAB = 2
)
res
is(res, "MetabinResult")
```

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