Package 'SynSigEval'

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Type Package

Title Evaluate Results of Mutational Signature Analysis Software on Synthetic Spectra Created by Package SynSigGen

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Author Steven G. Rozen, Yang Wu

Maintainer Steven G. Rozen <steverozen@gmail.com>

Description
Examine and evaluate the output of mutational signature analysis computational approaches.

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LazyData true

Imports data.table,
ICAMS,
ICAMSxtra (>= 0.0.2),
stats,
devtools,
utils,
rlang,
graphics,
grDevices,
ggplot2,
ggpubr,
ggbeeswarm,
gtools

Remotes github::steverozen/ICAMSxtra@*release

Depends R (>= 3.5)

RoxygenNote 7.1.1

Suggests testthat,
knitr,
lsa,
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DelayedArray,
BiocManager,
ggpubr

R topics documented:

`CopyBestSignatureAnalyzerResult`

Find the SignatureAnalyzer results directory with the best results and make a copy of it as `sa.results.dir/best.run/`

Description

Find the SignatureAnalyzer results directory with the best results and make a copy of it as `sa.results.dir/best.run/`

Usage

```
CopyBestSignatureAnalyzerResult(
  sa.results.dir,
  verbose = FALSE,
  overwrite = FALSE
)
```

Arguments

sa.results.dir	See BestSignatureAnalyzerResult
verbose	See BestSignatureAnalyzerResult
overwrite	If TRUE overwrite existing "best.run"

Value

The path of the best directory that was copied as a string, with the list directories examined as the attribute run.directories.

CreateEMuOutput *Prepare input file for EMu from a EMu formatted catalog file.*

Description

Prepare input file for EMu from a EMu formatted catalog file.

Usage

```
CreateEMuOutput(
  catalog,
  out.dir = paste0(dirname(catalog), "/ExtrAttr/EMu.results"),
  overwrite = FALSE
)
```

Arguments

catalog	a catalog in ICAMS format. It can be a .csv file, or a matrix or data.frame. Usually, it refers to "ground.truth.syn.catalog.csv".
out.dir	Directory that will be created for the output; abort if it already exists. Usually, the out.dir will be a EMu.results folder directly under the folder storing catalog.
overwrite	If TRUE, overwrite existing output

Details

Creates folder named EMu.results containing catalogs in EMu-formatted catalogs: Rows are signatures; the first column is the name of the mutation type, while the remaining columns are samples (tumors). These EMu-formatted catalogs will the input when running EMu program later on compiled binary.

Value

invisible(catalog), original catalog in EMu format

CreatehelmsmanOutput *Prepare input file for helmsman from a helmsman formatted catalog file.*

Description

Prepare input file for helmsman from a helmsman formatted catalog file.

Usage

```
CreatehelmsmanOutput(
  catalog,
  out.dir = paste0(dirname(catalog), "/ExtrAttr/helmsman.results"),
  overwrite = FALSE
)
```

Arguments

catalog	a catalog in ICAMS format. It can be a .csv file, or a matrix or data.frame. Usually, it refers to "ground.truth.syn.catalog.csv".
out.dir	Directory that will be created for the output; abort if it already exists. Usually, the out.dir will be a helmsman.results folder directly under the folder storing catalog.
overwrite	If TRUE, overwrite existing output

Details

Creates folder named helmsman.results containing catalogs in helmsman-formatted catalogs: Rows are signatures; the first column is the name of the mutation type, while the remaining columns are samples (tumors). These helmsman-formatted catalogs will the input when running helmsman program later on Python platform.

Value

invisible(catMatrix), original catalog in helmsman format

CreateMultiModalMuSigOutput *Prepare input file for MultiModalMuSig from a MultiModalMuSig formatted catalog file.*

Description

Prepare input file for MultiModalMuSig from a MultiModalMuSig formatted catalog file.

Usage

```
CreateMultiModalMuSigOutput(
  catalog,
  out.dir = paste0(dirname(catalog), "/ExtrAttr/MultiModalMuSig.results"),
  overwrite = FALSE
)
```

Arguments

catalog	a catalog in ICAMS format. It can be a .csv file, or a matrix or data.frame. Usually, it refers to "ground.truth.syn.catalog.csv".
out.dir	Directory that will be created for the output; abort if it already exists. Usually, the out.dir will be a MultiModalMuSig.results folder directly under the folder storing catalog.
overwrite	If TRUE, overwrite existing output

Details

Creates folder named MultiModalMuSig.results containing catalogs in MultiModalMuSig-formatted catalogs: Rows are signatures; the first column is the name of the mutation type, while the remaining columns are samples (tumors). These MM-formatted catalogs will the input when running MultiModalMuSig program later on Julia platform.

Value

invisible(catMatrix), original catalog in MultiModalMuSig format

---

helmsmanCatalog2ICAMS *Read Catalog files or matrices in helmsman format.*

Description

Read Catalog files or matrices in helmsman format.

Usage

```
helmsmanCatalog2ICAMS(
  cat,
  region = "unknown",
  catalog.type = "counts.signature"
)
```

Arguments

cat	Input catalog, can be a tab-delimited text file in helmsman format, or a matrix/data.frame object.
region	Catalog region. Can be a specific genomic or exomic region, or "unknown". Default: "unknown"
catalog.type	Is the catalog a signature catalog, or a spectrum catalog? Default: "counts.signature"

Value

a catalog matrix in ICAMS format.

ICAMSCatalog2EMu *Convert Catalogs from ICAMS format to EMu format*

Description

Convert Catalogs from ICAMS format to EMu format

Usage

```
ICAMSCatalog2EMu(catalog)
```

Arguments

catalog	A catalog matrix in ICAMS format. (SNS only!)

Value

a matrix without any dimnames, but the values are the transposition of the values in `catalog`.

ICAMSCatalog2helmsman *Convert Catalogs from ICAMS format to helmsman format*

Description

Convert Catalogs from ICAMS format to helmsman format

Usage

```
ICAMSCatalog2helmsman(catalog, type = "spectra")
```

Arguments

catalog	A catalog matrix in ICAMS format. (SNS only!)
type	Whether it is a spectra catalog ("spectra") or a signature catalog ("signature").

Value

a catalog matrix in helmsman format.

ICAMSCatalog2MM *Convert Catalogs from ICAMS format to MM format*

Description

Convert Catalogs from ICAMS format to MM format

Usage

```
ICAMSCatalog2MM(catalog)
```

Arguments

catalog	A catalog matrix in ICAMS format. (SNS/DNS/ID)

Value

a catalog matrix in MultiModalMuSig format.

MMCatalog2ICAMS *Convert Catalogs (File or Matrix) from MM format to ICAMS format*

Description

Convert Catalogs (File or Matrix) from MM format to ICAMS format

Usage

```
MMCatalog2ICAMS(cat, region = "unknown", catalog.type = "counts.signature")
```

Arguments

cat	Input catalog, can be a tab-delimited file or matrix in MultiModalMuSig format.
region	Catalog region. Can be a specific genomic or exomic region, or "unknown". Default: "unknown"
catalog.type	Is the catalog a signature catalog, or a spectrum catalog? Default: "counts.signature"

Value

a catalog matrix in ICAMS format.

`PlotCatCOMPOSITE` *Plot the a SignatureAnalyzer COMPOSITE signature or catalog into separate pdfs*

Description

Plot the a SignatureAnalyzer COMPOSITE signature or catalog into separate pdfs

Usage

```
PlotCatCOMPOSITE(catalog, filename.header, type, id = colnames(catalog))
```

Arguments

`catalog`	Catalog or signature matrix
`filename.header`	Contain path and the beginning part of the file name. The name of the pdf files will be: `filename.header`.SNS.96.pdf `filename.header`.SNS.1536.pdf `filename.header`.DNS.78.pdf `filename.header`.ID.83.pdf
`type`	See `PlotCatalogToPdf`.
`id`	A vector containing the identifiers of the samples or signatures in `catalog`.

`ReadAndAnalyzeExposures` *Assess how well inferred exposures match input exposures We assume that in many cases attribution programs will be run outside of R on file inputs and will generate fill outputs.*

Description

Assess how well inferred exposures match input exposures

We assume that in many cases attribution programs will be run outside of R on file inputs and will generate fill outputs.

Usage

```
ReadAndAnalyzeExposures(
  extracted.sigs,
  ground.truth.sigs,
  inferred.exp.path,
  ground.truth.exposures
)
```

Arguments

`extracted.sigs` Path to file containing the extracted signature profiles.

`ground.truth.sigs`
File containing signature profiles from which the synthetic data were generated.

`inferred.exp.path`
File containing mutation counts (exposures) of synthetic tumors which are inferred to extracted or input signatures.

`ground.truth.exposures`
File containing the exposures from which the synthetic catalogs were generated. This file is used to restrict assessment of signature exposures to only those signatures in `ground.truth.sigs` that were actually represented in the exposures.

Details

Generates output files by calling `MatchSigsAndRelabel`

Value

A `data.frame` recording:

`Ground.truth.exposure`: sum of ground truth exposures of all tumors to all ground-truth signatures.

`Inferred.exposure`: sum of inferred exposures of all tumors to all ground-truth signatures. Here, inferred exposure of a tumor to a ground-truth signature equals to the sum of the exposures of this tumor to all extracted signatures which are most similar to a ground-truth signature. If there is no extracted signature resembling an ground-truth signature, the inferred exposure of this ground-truth signature will be `0`.

`Absolute.difference`: sum of absolute difference between ground-truth exposure and inferred exposure of all tumors to all ground-truth signatures.

`ReadAndAnalyzeSigs` *Assess how well extracted signatures match input signatures We assume that in many cases extraction programs will be run outside of R on file inputs and will generate fill outputs.*

Description

Assess how well extracted signatures match input signatures

We assume that in many cases extraction programs will be run outside of R on file inputs and will generate fill outputs.

Usage

```
ReadAndAnalyzeSigs(extracted.sigs, ground.truth.sigs, ground.truth.exposures)
```

Arguments

`extracted.sigs`	Path to file containing the extracted signature profiles.
`ground.truth.sigs`	File containing signature profiles from which the synthetic data were generated.
`ground.truth.exposures`	File containing the exposures from which the synthetic catalogs were generated. This file is used to restrict assessment to only those signatures in `ground.truth.sigs` that were actually represented in the exposures.

Details

Generates output files by calling `MatchSigsAndRelabel`

Value

See `MatchSigsAndRelabel`

`ReadEMuCatalog` *Read Catalog files in EMu format.*

Description

Read Catalog files in EMu format.

Usage

```
ReadEMuCatalog(
  cat,
  mutTypes,
  sigOrSampleNames,
  region = "unknown",
  catalog.type = "counts.signature"
)
```

Arguments

`cat`	A tab-delimited catalog text file in EMu format; or a EMu formatted matrix or data.frame.
`mutTypes`	Types of mutations. They are usually from an `ICAMS:::catalog.row.header` object.
`sigOrSampleNames`	If input file is a counts signature file (`catalog.type == "counts.signature"`), signature names should be provided. If input file is a counts spectra file (`catalog.type == "counts"`), names of samples should be provided.
`region`	Catalog region. Can be a specific genomic or exomic region, or "unknown". Default: "unknown"
`catalog.type`	Is the catalog a signature catalog, or a spectrum catalog? Default: "counts"

Value

a catalog matrix in ICAMS format.

`ReadEMuExposureFile` *Read Exposure files in EMu format.*

Description

Read Exposure files in EMu format.

Usage

```
ReadEMuExposureFile(exposureFile, sigNames, sampleNames)
```

Arguments

`exposureFile`	Exposure file generated by EMu. Usually, it is called "W_components.txt".
`sigNames`	Names of signatures. These will be served as the rownames of the exposure matrix.
`sampleNames`	Names of samples in exposure file.
	Return ICAMS/SynSigEval formatted exposure matrix.

`ReadExposureMM` *Read Catalog files in MM format*

Description

Read Catalog files in MM format

Usage

```
ReadExposureMM(exposureFile)
```

Arguments

`exposureFile`	Input exposure file, can be a tab-delimited text file in MultiModalMuSig format.

Value

a exposure matrix in ICAMS format.

ReadhelmsmanExposure *Read Exposure files in helmsman format.*

Description

Read Exposure files in helmsman format.

Usage

```
ReadhelmsmanExposure(exposure, check.names = TRUE)
```

Arguments

exposure	Exposure file generated by helmsman. Usually, it is called "W_components.txt".
check.names	logical. If TRUE then the names of the variables in the data frame are checked to ensure that they are syntactically valid variable names. If necessary they are adjusted (by make.names) so that they are, and also to ensure that there are no duplicates.
	Return ICAMS/SynSigEval formatted exposure matrix.

ReadSigProfilerExposure *Read a file containing exposures attributed by SigProfiler/Python*

Description

Read a file containing exposures attributed by SigProfiler/Python

Usage

```
ReadSigProfilerExposure(file)
```

Arguments

file	The name of the file to read.

Value

The corresponding signature matrix in standard internal representation.

ReadSigProfilerSigDBS78

Read a file containing DBS78 signatures extracted by SigProfiler/Python

Description

Read a file containing DBS78 signatures extracted by SigProfiler/Python

Usage

```
ReadSigProfilerSigDBS78(file)
```

Arguments

file	The name of the file to read.

Value

The corresponding signature matrix in standard internal representation.

ReadSigProfilerSigSBS96

Read a file containing SBS96 signatures extracted by SigProfiler/Python

Description

Read a file containing SBS96 signatures extracted by SigProfiler/Python

Usage

```
ReadSigProfilerSigSBS96(file)
```

Arguments

file	The name of the file to read.

Value

The corresponding signature matrix in standard internal representation.

`SignatureAnalyzerSummarizeSBS1SBS5`

Summarize all sub-directories of SignatureAnalyzer results on the correlated SBS1 / SBS5.

Description

This is special-purpose function to summarize results from one in-silico experiment that examines how well signatures can be extracted from synthetic tumors with correlated SBS1 and SBS5.

Usage

```
SignatureAnalyzerSummarizeSBS1SBS5(top.level.dir, overwrite = FALSE)
```

Arguments

`top.level.dir`	Path to top level directory.
`overwrite`	If TRUE overwrite existing directories and files.

`SignatureAnalyzerSummarizeTopLevel`

Summarize all subdirectories of SignatureAnalyzer results on a major dataset.

Description

This function depends on a particular directory structure: see argument `top.level.dir`. This function finds the best of multiple SignatureAnalyzer extraction runs and summarizes the comparison of the best run with the ground truth.

Usage

```
SignatureAnalyzerSummarizeTopLevel(top.level.dir, overwrite = FALSE)
```

Arguments

`top.level.dir`	Path to top level directory, which must contain the following subdirectories: • `sa.sa.96/sa.results/` • `sp.sp/sa.results/` • `sa.sa.COMPOSITE/sa.results/` • `sp.sa.COMPOSITE/sa.results/` Each of the directories must contain additional subdirectories, one for each SignatureAnalyzer run, names `sa.run.<n>`, where <n> is an integer (string of digits).
`overwrite`	If TRUE overwrite existing summary files.

SplitCatCOMPOSITE *Split COMPOSITE (SNS1536+DBS78+ID83) catalogs in ICAMS format into 3 individual catalogs.*

Description

Split COMPOSITE (SNS1536+DBS78+ID83) catalogs in ICAMS format into 3 individual catalogs.

Usage

```
SplitCatCOMPOSITE(catalog)
```

Arguments

catalog	Input catalog, can be a .csv file or matrix in ICAMS COMPOSITE format.

Value

a list, containing 3 catalog matrices in MultiModalMuSig format. Each matrix contains SNS1536, DBS78 and ID83 information, respectively.

SummarizeMultiRuns *Assess/evaluate multiple summarized runs for one dataset from one computational approach.*

Description

Summarize results from each computational approach in `tool.dir/run.names` (generated by running a computational approach), combine them into `tool.dir`.

Usage

```
SummarizeMultiRuns(datasetName, toolName, tool.dir, run.names)
```

Arguments

datasetName	Name of the dataset. (e.g. "S.0.1.Rsq.0.1"). Usually, it is has the same name as `basename(top.dir)`.
toolName	Name of computational approach. (e.g. "SigProExtractor")
tool.dir	Fourth level path from the `top.dir`. Expected to have multiple runs with different names (e.g. "seed.1") That is, `top.dir/sp.sp/ExtrAttr/sa.results/`. or `top.dir/sa.sa.96/Attr/deconstructSigs.results/` Here, `top.dir` refers to a top-level directory which contains the full information of a synthetic dataset. (e.g. `syn.2.7a.7b.abst.v8`) This code depends on a conventional directory structure documented elsewhere. However there should be a directory within the `tool.dir` which stores the software output.
run.names	A character vector records the list of `run.dir`, or fifth level directories from the dataset top-level folder. E.g., c("seed.1","seed.691")

Details

Also writes multiple files into folder `tool.dir`.

Value

A list contain values of measures measures in multiple runs:

- $averCosSim Cosine similarity
- $truePos True Positives(TP): Ground-truth signatures which are active in the spectra, and extracted.
- $falseNeg False Negatives(FN): Ground-truth signatures not extracted.
- $falsePos False Positives(FP): Signatures wrongly extracted, not resembling any ground-truth signatures.
- $TPR True positive rate (TPR, Sensitivity): TP / (TP + FN)
- $PPV Positive predictive value (PPV): TP / (FP + TP)
- $cosSim Average cosine similarity to each of the ground-truth signatures.
- $AggManhattanDist Scaled Manhattan distance between ground-truth and inferred exposures to each of the ground-truth signatures.

This list also contains `mean` and `sd`, and other statistics of these measures in

- $fivenum
- $fivenumMD
- $meanSD
- $meanSDMD

SummarizeMultiToolsMultiDatasets

Summarize results for multiple datasets, by different computational approaches.

Description

Summarize results of mutational signature extraction and exposure inferrence by multiple computational approaches on multiple datasets. Before running this function, make sure the summary file for each single data set `third.level.dir/multiTools.Rda` exists.

Usage

```
SummarizeMultiToolsMultiDatasets(
  dataset.dirs,
  second.third.level.dirname,
  out.dir,
  overwrite = FALSE
)
```

Arguments

dataset.dirs	Paths of top-level dataset directories trees you want to investigate. E.g. "./S.0.1.Rsq.0.1"
second.third.level.dirname	Name of the second.level.dir (e.g. "sp.sp") and the third.level.dir (e.g. "ExtrAttr") to be investigated. Examples are: "sp.sp/ExtrAttr", "sa.sa.96/Attr" Note: multiTools.RDa are expected to be exist under dataset.dirs/second.third.level.dirnam
out.dir	Path of the output directory.
overwrite	Whether to overwrite the contents in out.dir if it already exists. (Default: FALSE)

Details

multiTools.Rda is generated by SummarizeMultiToolsOneDataset).

SummarizeMultiToolsOneDataset

Combine results for a single dataset, from different computational approaches.

Description

Summarize results from each computational approach in third.level.dir/tool.dirnames (generated by SummarizeMultiRuns), combine them into third.level.dir.

Usage

```
SummarizeMultiToolsOneDataset(
  third.level.dir,
  toolNames,
  tool.dirnames,
  datasetGroup,
  datasetGroupName,
  datasetSubGroup = NULL,
  datasetSubGroupName = NULL
)
```

Arguments

third.level.dir	Third level path distinguishing de novo extraction + attribution packages from attribution-only packages. Examples: top.dir/sp.sp/ExtrAttr/ top.dir/sa.sa/Attr/
toolNames	Names of computational approach. (e.g. "SigProExtractor")
tool.dirnames	Third level path from the top.dir. Expected to have summarized results generated by SummarizeMultiRuns. (multiRun.RDa, ManhattanDist.csv, meanSD.csv, meanSD.Manhattan.dist.csv) Examples: "signeR.results" (Under third.level.dir "ExtrAttr") "deconstructSigs.results" (Under third.level.dir "Attr") Here, top.dir refers to a top-level directory which contains the full information of a synthetic dataset. (e.g. syn.2.7a.7b.abst.v8) This code depends on a conventional directory structure documented elsewhere. However there should be a directory within the tool.names which stores the software output.

datasetGroup
: Numeric or character vector specifying the groups each dataset belong to. E.g. For SBS1-SBS5 correlated datasets, we can consider slope as the group: c("slope=0.1","slope=0.5","s Default: "Default"

datasetGroupName
: Meaning or label of all datasetGroup. E.g. For SBS1-SBS5 correlated datasets, we can consider "SBS1:SBS5 mutation count ratio" as the label of the datasetGroup slope.

datasetSubGroup
: Optional. Numeric or character vector differentiating datasets within each group. E.g. For SBS1-SBS5 correlated datasets, we can consider Pearson's R^2 as the subgroup: c("Rsq=0.1","Rsq=0.2","Rsq=0.3","Rsq=0.6") Default: Names of datasets, which are basename(dataset.dirs)

datasetSubGroupName
: Optional. Meaning or label of all datasetSubGroup. E.g. For SBS1-SBS5 correlated datasets, we can consider "Pearson's R squared" as the label of the datasetSubGroup Pearson's R^2.

Details

This function generates multiTools.RDa under third.level.dir

Value

A list contain c(mean,sd) of multiple runs: Cosine similarity True Positives(TP): Ground-truth signatures which are active in the spectra, and extracted. False Negatives(FN): Ground-truth signatures not extracted. False Positives(FP): Signatures wrongly extracted, not resembling any ground-truth signatures. True positive rate (TPR, Sensitivity): TP / (TP + FN) Positive predictive value (PPV, Precision): TP / (FP + TP)

SummarizeOneToolMultiDatasets

Combine results for multiple datasets, from one computational approaches.

Description

Summarize results from each computational approach in third.level.dir/tool.dirnames (generated by SummarizeMultiRuns), combine them into third.level.dir.

Usage

```
SummarizeOneToolMultiDatasets(
  dataset.dirs,
  datasetGroup,
  datasetGroupName,
  datasetSubGroup = NULL,
  datasetSubGroupName = NULL,
  toolName,
  tool.dirname,
  out.dir,
  overwrite = FALSE
)
```

Arguments

dataset.dirs	Paths of top-level dataset directories trees you want to investigate. E.g. "./S.0.1.Rsq.0.1"
datasetGroup	Numeric or character vector specifying the group each dataset belong to. E.g. For SBS1-SBS5 correlated datasets, we can consider slope (SBS1:SBS5 count ratio) as the group: c(0.1,0.5,1,2,5,10) Default: "Default"
datasetGroupName	Meaning or label of all datasetGroup. E.g. For SBS1-SBS5 correlated datasets, we can consider "SBS1:SBS5 mutation count ratio" as the label of the datasetGroup slope.
datasetSubGroup	Numeric or character vector differentiating datasets within each group. E.g. For SBS1-SBS5 correlated datasets, we can consider Pearson's R^2 as the subgroup: c(0.1,0.2,0.3,0.6) Default: Names of datasets, which are basename(dataset.dirs)
datasetSubGroupName	Meaning or label of all datasetSubGroup. E.g. For SBS1-SBS5 correlated datasets, we can consider "Pearson's R squared" as the label of the datasetSubGroup Pearson's R^2.
toolName	Name of computational approach to be investigated (e.g. "SigProExtractor")
tool.dirname	Name of the second.level.dir (e.g. "sp.sp"), third.level.dir (e.g. "ExtrAttr") and tool.dir (e.g. "SigProExtractor.results") to be investigated. One example: "sp.sp/ExtrAttr/SigProExtractor.results" Note: this function expects the summary generated by SummarizeSigOneSubdir under dataset.dirs/tool.dirname
out.dir	Path of the output directory.
overwrite	Whether to overwrite the contents in out.dir if it already exists. (Default: FALSE)

SummarizeSigOneAttrSubdir

Assess/evaluate results from packages which can ONLY do exposure attribution.

Description

Packages including but not limited to: deconstructSigs, YAPSA.

Usage

```
SummarizeSigOneAttrSubdir(
  run.dir,
  ground.truth.exposure.dir = paste0(run.dir, "/../../../"),
  overwrite = FALSE
)
```

Arguments

run.dir	Lowest level path to results, e.g. <top.dir>/sa.sa.96/Attr/YAPSA.results/seed.1/ Here, <top.dir> refers to a top-level directory which contains the full information of a synthetic dataset. (e.g. syn.2.7a.7b.abst.v8) This code depends on a conventional directory structure documented elsewhere. For packages which can do both extraction and attribution, we expect two files, ground.truth.signatures.csv and inferred.exposures.csv are in the folder.
ground.truth.exposure.dir	Folder which stores ground-truth exposures. It defaults to be sub.dir, i.e. run.dir/../../
overwrite	If TRUE overwrite existing directories and files.

Details

Here, we excluded SignatureEstimation. Although it is also a package with only attribution, but it has two attribution algorithms. Therefore the naming of the results are slightly different from the other two packages.

SummarizeSigOneExtrAttrSubdir

Assess/evaluate results from packages which can do BOTH extraction and attribution, excluding SigProfiler-Python and SignatureAnalyzer.

Description

Packages including but not limited to: hdp, MutationalPatterns, sigfit, signeR, SomaticSignatures.

Usage

```
SummarizeSigOneExtrAttrSubdir(
  run.dir,
  ground.truth.exposure.dir = paste0(run.dir, "/../../../"),
  overwrite = FALSE
)
```

Arguments

run.dir	Lowest level path to result of a run. E.g. <top.dir>/sa.sa.96/ExtrAttr/SomaticSignatures.res Here, <top.dir> refers to a top-level directory which contains the full information of a synthetic dataset. (e.g. syn.2.7a.7b.abst.v8) This code depends on a conventional directory structure documented elsewhere. For packages which can do both extraction and attribution, we expect two files, extracted.signatures.csv and inferred.exposures.csv are in the folder.
ground.truth.exposure.dir	Folder which stores ground-truth exposures. It defaults to be sub.dir, i.e. run.dir/../../../
overwrite	If TRUE overwrite existing directories and files.

SummarizeSigOnehelmsmanSubdir

Assess/evaluate results from SigProfiler-python (a.k.a. SigProExtractor) Assessment is restricted to v0.0.5.43, because different version has different folder structure.

Description

Assess/evaluate results from SigProfiler-python (a.k.a. SigProExtractor) Assessment is restricted to v0.0.5.43, because different version has different folder structure.

Usage

```
SummarizeSigOnehelmsmanSubdir(
  run.dir,
  ground.truth.exposure.dir = paste0(run.dir, "/../../../"),
  overwrite = FALSE,
  hierarchy = FALSE
)
```

Arguments

run.dir	Lowest level path to results, e.g. <top.dir>/sa.sa.96/ExtrAttr/SigProExtractor.results/see Here, <top.dir> refers to a top-level directory which contains the full information of a synthetic dataset. (e.g. syn.2.7a.7b.abst.v8) This code depends on a conventional directory structure documented elsewhere. However there should be a directory <run.dir>/SBS96 which stores SigProfiler results.
ground.truth.exposure.dir	Folder which stores ground-truth exposures. Usually, it refers to sub.dir, i.e. run.dir/../../../
overwrite	If TRUE overwrite existing directories and files.
hierarchy	Whether the user have enabled hierarchy = True when running SigProExtractor. specifying True or False into SigProExtractor will cause the program to generate different folder structure. (Default: FALSE)

SummarizeSigOneSigProExtractorSubdir

Assess/evaluate results from SigProExtractor SigProfiler-python de novo extraction and attribution package. Assessment is restricted to v0.0.5.43+, because different version has different folder structure.

Description

Assess/evaluate results from SigProExtractor SigProfiler-python de novo extraction and attribution package. Assessment is restricted to v0.0.5.43+, because different version has different folder structure.

Usage

```
SummarizeSigOneSigProExtractorSubdir(
  run.dir,
  ground.truth.exposure.dir = paste0(run.dir, "/../../../"),
  overwrite = FALSE,
  hierarchy = FALSE
)
```

Arguments

run.dir	Lowest level path to results, e.g. <top.dir>/sa.sa.96/ExtrAttr/SigProExtractor.results/see Here, <top.dir> refers to a top-level directory which contains the full information of a synthetic dataset. (e.g. syn.2.7a.7b.abst.v8) This code depends on a conventional directory structure documented elsewhere. However there should be a directory <run.dir>/SBS96 which stores SigProfiler results.
ground.truth.exposure.dir	TODO(Wu Yang): Fix this File name which stores ground-truth exposures; defaults to "ground.truth.syn.exposures.csv". This file can be found in the sub.dir, i.e. <run.dir>/../../../
overwrite	If TRUE overwrite existing directories and files.
hierarchy	Whether the user have enabled hierarchy = True when running SigProExtractor. specifying True or False into SigProExtractor will cause the program to generate different folder structure. (Default: FALSE)

SummarizeSigOneSigProSSSubdir

Assess/evaluate results from sigproSS (a.k.a. SigProfiler Python attribution package)

Description

Assess/evaluate results from sigproSS (a.k.a. SigProfiler Python attribution package)

Usage

```
SummarizeSigOneSigProSSSubdir(
  run.dir,
  ground.truth.exposure.dir = paste0(run.dir, "/../../../"),
  overwrite = FALSE
)
```

Arguments

run.dir	Lowest level path to results, e.g. <top.dir>/sa.sa.96/ExtrAttr/SigProExtractor.results/see Here, <top.dir> refers to a top-level directory which contains the full information of a synthetic dataset. (e.g. syn.2.7a.7b.abst.v8) This code depends on a conventional directory structure documented elsewhere. However there should be a directory <run.dir>/SBS96 which stores SigProfiler results.

ground.truth.exposure.dir
: TODO(Wu Yang): Fix this File name which stores ground-truth exposures; defaults to "ground.truth.syn.exposures.csv". This file can be found in the sub.dir, i.e. <run.dir>/../../../

overwrite
: If TRUE overwrite existing directories and files.

SummarizeSigProExtractor *Summarize SigProfiler results in the sa.sa.96 and/or sp.sp subdirectories.*

Description

Summarize SigProfiler results in the sa.sa.96 and/or sp.sp subdirectories.

Usage

```
SummarizeSigProExtractor(
  top.dir,
  sub.dir = c("sa.sa.96", "sp.sp"),
  overwrite = FALSE
)
```

Arguments

top.dir
: The top directory of a conventional data structure containing at least one of the subdirectories: sa.sa.96/sp.results and sp.sp/sp.results; see further documentation elsewhere.

sub.dir
: The subdirectory under top.dir, and containing a folder named sp.results. By default, it contains both c("sa.sa","sp.sp"). But you should specify sub.dir = "sp.sp" for top.dir with only the sp.sp subdirectory (as is the case for the correlated SBS1-and-SBS5-containing data sets).

overwrite
: whether to overwrite the existing run.dir/summary folder? If chosen to be FALSE and there is an existing summary folder, an error will be raised.

Details

Results are put in standardized subdirectories of top.dir.

SynSigEval *SynSigEval*

Description

Assess the performance of two steps in mutational signature analysis:

- signature extraction
- exposure inferrence (a.k.a. signature attribution)

by computational approaches, using catalogs of synthetic mutational spectra created by package SynSigGen.

Input

SynSigEval requires the input data listed below:

1. E, matrix of synthetic exposures (signatures x samples)
2. S, mutational signature profiles (mutation type x signature)
3. synthetic.spectra, synthetic mutational spectra with known ground-truth mutational signature profiles (S) and exposures (synthetic.exposures). It can be created from SynSigGen.
4. T, signatures extracted by SignatureAnalzer, SigProfiler, or other computational approaches on synthetic.spectra. For attribution-only approaches, T=S.
5. F, exposures inferred by computational approaches on synthetic.spectra.

Folder structure for SynSigEval v0.2

Summary function will fit to the new 5-level folder structure:

First Level - top.level.dir: dataset folder (e.g. "S.0.1.Rsq.0.1", "syn.pancreas"). All spectra datasets under any top.level.dir have the same exposure.

Second Level - ground.truth.exposure.dir: spectra folder: (e.g. "sp.sp", "sa.sa.96"). All spectra datasets under any second.level.dir have the same signature and the same exposure counts.

Third Level - third.level.dir: It can be ("Attr") for storing results of packages which can only do exposure attribution of known signatures ("Attr"); it can also be ("ExtrAttr"), folder to store results of software packages which can do de-novo extraction and following attribution.

Fourth Level - tool.dir: The results of a software package (e.g. "SigProExtractor.results","SignatureEstimation.QP.resu
Under this level, tool.dir may contain multiple run.dir, each is a run of the software package using a specific number of seed.

Fifth level - run.dir: contains results from a run of the software package using a specific number of seed. (e.g. "seed.1")

Summarize results

1. Summarize results in fifth-level run.dir:
   Relevant functions are:
   - SummarizeSigProExtractor
   - SignatureAnalyzerSummarizeTopLevel
   - SignatureAnalyzerSummarizeSBS1SBS5
   - SummarizeSigOneExtrAttrSubdir
   - SummarizeSigOneAttrSubdir
   - SummarizeSigOnehelmsmanSubdir
   - SummarizeSigOneSigProSSSubdir
2. Summarize results of multiple runs by a computational approach on one spectra data set:
   SummarizeMultiRuns
3. Summarize results of multiple computational approaches on one spectra data set:
   SummarizeMultiToolsOneDataset
4. Summarize results of multiple computational approaches on multiple spectra data sets:
   SummarizeMultiToolsMultiDatasets

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