Package 'mSigHdp'

May 16, 2020

Title Mutational signature extraction using hdp

Version 0.0.0.9005

Description Calls hdp for mutational signature analysis, with performance issues in hdp:::stirling() corrected.

License GPL-3

Encoding UTF-8

LazyData true

Language en-US

biocViews

Imports hdp,
hdpx,
SynSigGen,
SynSigEval

Roxygen list(markdown = TRUE)

Depends R (>= 3.5)

RoxygenNote 7.1.0

Remotes github::nicolaroberts/hdp,
github::steverozen/hdpx,
github::steverozen/SynSigGen,
github::WuyangFF95/SynSigEval

Suggests testthat,
ICAMS,
hdp,
hdpx,
SynSigEval,
SynSigGen,
utils

R topics documented:

`RunAndEvalHdp4` *Run and evaluate hdp*

Description

Run and evaluate hdp

Usage

```
RunAndEvalHdp4(
  input.catalog,
  ground.truth.exposure.file,
  ground.truth.sig.file = NULL,
  ground.truth.sig.catalog = NULL,
  out.dir,
  CPU.cores = 1,
  seedNumber = 1,
  K.guess,
  multi.types = FALSE,
  remove.noise = FALSE,
  test.only = 0,
  overwrite = FALSE,
  verbose = TRUE,
  num.posterior = 4,
  post.burnin = 4000,
  post.n = 50,
  post.space = 50,
  post.cpiter = 3,
  post.verbosity = 0,
  cos.merge = 0.9,
  min.sample = 1
)
```

Arguments

`input.catalog`	Either a character string, in which case this is the path to a file containing a spectra catalog in `ICAMS` format, or an `ICAMS` catalog.
`ground.truth.exposure.file`	Path to file with ground truth exposures.
`ground.truth.sig.file`	Path to file with ground truth signatures.
`ground.truth.sig.catalog`	`ICAMS` catalog with signatures used to construct the ground truth spectra. Specify only one of `ground.truth.sig.file.path` or `ground.truth.sig.catalog`.
`out.dir`	Directory that will be created for the output; if `overwrite` is `FALSE` then abort if `out.dir` already exits.
`CPU.cores`	Number of CPUs to use in running `hdp_posterior`; this is used to parallize running the posterior sampling chains, so there is no point in making this larger than `num.posterior`.

seedNumber	An integer that is used to generate separate random seeds for each call to dp_activate, and each call of hdp_posterior; please see the code on how this is done. But repeated calls with same value of seedNumber and other inputs should produce the same results.
K.guess	Suggested initial value of the number of signatures, passed to dp_activate as initcc.
multi.types	A logical scalar or a character vector. If FALSE, hdp will regard all input spectra as one tumor type. If TRUE, hdp will infer tumor types based on the string before "::" in their names. e.g. tumor type for "SA.Syn.Ovary-AdenoCA::S.500" would be "SA.Syn.Ovary-AdenoCA" If multi.types is a character vector, then it should be of the same length as the number of columns in input.catalog, and each value is the name of the tumor type of the corresponding column in input.catalog, e.g. c("SA.Syn.Ovary-AdenoCA","SA.Syn.O
remove.noise	Deprecated; ignored
test.only	If > 0, only analyze the first test.only columns in input.catalog.
overwrite	If TRUE overwrite out.dir if it exists, otherwise raise an error.
verbose	If TRUE then message progress information.
num.posterior	Number of posterior sampling chains; can set to 1 for testing.
post.burnin	Pass to hdp_posterior burnin.
post.n	Pass to hdp_posterior n.
post.space	Pass to hdp_posterior space.
post.cpiter	Pass to hdp_posterior cpiter.
post.verbosity	Pass to hdp_posterior verbosity.
cos.merge	The cosine similarity threshhold for merging raw clusters from the posterior sampling chains into "components" i.e. signatures; passed to hdp_extract_components.
min.sample	A "component" (i.e. signature) must have at least this many samples; passed to hdp_extract_components.

RunAndEvalHdp5 *Run and evaluate hdp*

Description

Run and evaluate hdp

Usage

```
RunAndEvalHdp5(
  input.catalog.file,
  ground.truth.exposure.file,
  ground.truth.sig.file = NULL,
  ground.truth.sig.catalog = NULL,
  out.dir,
  CPU.cores = 1,
  seedNumber = 1,
```

```
  K.guess,
  multi.types = FALSE,
  remove.noise = FALSE,
  test.only = 0,
  overwrite = FALSE,
  verbose = TRUE,
  num.posterior = 4,
  post.burnin = 4000,
  post.n = 50,
  post.space = 50,
  post.cpiter = 3,
  post.verbosity = 0,
  cos.merge = 0.9,
  min.sample = 1
)
```

Arguments

input.catalog.file	File containing a spectra catalog in ICAMS format.
ground.truth.exposure.file	Path to file with ground truth exposures.
ground.truth.sig.file	Path to file with ground truth signatures.
ground.truth.sig.catalog	ICAMS catalog with signatures used to construct the ground truth spectra. Specify only one of ground.truth.sig.file.path or ground.truth.sig.catalog.
out.dir	Directory that will be created for the output; if overwrite is FALSE then abort if out.dir already exits.
CPU.cores	Number of CPUs to use in running hdp_posterior; this is used to parallize running the posterior sampling chains, so there is no point in making this larger than num.posterior.
seedNumber	An integer that is used to generate separate random seeds for each call to dp_activate, and each call of hdp_posterior; please see the code on how this is done. But repeated calls with same value of seedNumber and other inputs should produce the same results.
K.guess	Suggested initial value of the number of signatures, passed to dp_activate as initcc.
multi.types	A logical scalar or a character vector. If FALSE, hdp will regard all input spectra as one tumor type. If TRUE, hdp will infer tumor types based on the string before "::" in their names. e.g. tumor type for "SA.Syn.Ovary-AdenoCA::S.500" would be "SA.Syn.Ovary-AdenoCA" If multi.types is a character vector, then it should be of the same length as the number of columns in input.catalog, and each value is the name of the tumor type of the corresponding column in input.catalog, e.g. c("SA.Syn.Ovary-AdenoCA","SA.Syn.O
remove.noise	Deprecated; ignored
test.only	If > 0, only analyze the first test.only columns in input.catalog.file.
overwrite	If TRUE overwrite out.dir if it exists, otherwise raise an error.

verbose	If TRUE then message progress information.
num.posterior	Number of posterior sampling chains; can set to 1 for testing.
post.burnin	Pass to hdp_posterior burnin.
post.n	Pass to hdp_posterior n.
post.space	Pass to hdp_posterior space.
post.cpiter	Pass to hdp_posterior cpiter.
post.verbosity	Pass to hdp_posterior verbosity.
cos.merge	The cosine similarity threshhold for merging raw clusters from the posterior sampling chains into "components" i.e. signatures; passed to hdp_extract_components.
min.sample	A "component" (i.e. signature) must have at least this many samples; passed to hdp_extract_components.

Runhdp4 *Run hdp extraction and attribution on a spectra catalog file using hdpx*

Description

Run hdp extraction and attribution on a spectra catalog file using hdpx

Usage

```
Runhdp4(
  input.catalog,
  out.dir,
  CPU.cores = 1,
  seedNumber = 1,
  K.guess,
  multi.types = FALSE,
  remove.noise = FALSE,
  test.only = 0,
  overwrite = FALSE,
  verbose = TRUE,
  num.posterior = 4,
  post.burnin = 4000,
  post.n = 50,
  post.space = 50,
  post.cpiter = 3,
  post.verbosity = 0,
  cos.merge = 0.9,
  min.sample = 1,
  plot.extracted.sig = FALSE
)
```

Arguments

input.catalog	Either a character string, in which case this is the path to a file containing a spectra catalog in ICAMS format, or an ICAMS catalog.
out.dir	Directory that will be created for the output; if overwrite is FALSE then abort if out.dir already exits.

CPU.cores	Number of CPUs to use in running hdp_posterior; this is used to parallize running the posterior sampling chains, so there is no point in making this larger than num.posterior.
seedNumber	An integer that is used to generate separate random seeds for each call to dp_activate, and each call of hdp_posterior; please see the code on how this is done. But repeated calls with same value of seedNumber and other inputs should produce the same results.
K.guess	Suggested initial value of the number of signatures, passed to dp_activate as initcc.
multi.types	A logical scalar or a character vector. If FALSE, hdp will regard all input spectra as one tumor type. If TRUE, hdp will infer tumor types based on the string before "::" in their names. e.g. tumor type for "SA.Syn.Ovary-AdenoCA::S.500" would be "SA.Syn.Ovary-AdenoCA" If multi.types is a character vector, then it should be of the same length as the number of columns in input.catalog, and each value is the name of the tumor type of the corresponding column in input.catalog, e.g. c("SA.Syn.Ovary-AdenoCA","SA.Syn.O
remove.noise	Deprecated; ignored
test.only	If > 0, only analyze the first test.only columns in input.catalog.
overwrite	If TRUE overwrite out.dir if it exists, otherwise raise an error.
verbose	If TRUE then message progress information.
num.posterior	Number of posterior sampling chains; can set to 1 for testing.
post.burnin	Pass to hdp_posterior burnin.
post.n	Pass to hdp_posterior n.
post.space	Pass to hdp_posterior space.
post.cpiter	Pass to hdp_posterior cpiter.
post.verbosity	Pass to hdp_posterior verbosity.
cos.merge	The cosine similarity threshhold for merging raw clusters from the posterior sampling chains into "components" i.e. signatures; passed to hdp_extract_components.
min.sample	A "component" (i.e. signature) must have at least this many samples; passed to hdp_extract_components.
plot.extracted.sig	If TRUE then plot the extracted signatures.

Details

Creates several files in out.dir. These are: call.and.session.info.txt, hdp.diagnostics.pdf, Runhdp4.retval.Rdata, extracted.signatures.csv, extracted.signature.pdf (optional), inferred.exposures.csv.

Value

The same list as returned by RunhdpInternal4.

Runhdp5 *Run hdp extraction and attribution on a spectra catalog file using the original hdp package.*

Description

Run hdp extraction and attribution on a spectra catalog file using the original hdp package.

Usage

```
Runhdp5(
  input.catalog.file,
  out.dir,
  CPU.cores = 1,
  seedNumber = 1,
  K.guess,
  multi.types = FALSE,
  remove.noise = FALSE,
  test.only = 0,
  overwrite = FALSE,
  verbose = TRUE,
  num.posterior = 4,
  post.burnin = 4000,
  post.n = 50,
  post.space = 50,
  post.cpiter = 3,
  post.verbosity = 0,
  cos.merge = 0.9,
  min.sample = 1
)
```

Arguments

input.catalog.file	File containing a spectra catalog in ICAMS format.
out.dir	Directory that will be created for the output; if overwrite is FALSE then abort if out.dir already exits.
CPU.cores	Number of CPUs to use in running hdp_posterior; this is used to parallize running the posterior sampling chains, so there is no point in making this larger than num.posterior.
seedNumber	An integer that is used to generate separate random seeds for each call to dp_activate, and each call of hdp_posterior; please see the code on how this is done. But repeated calls with same value of seedNumber and other inputs should produce the same results.
K.guess	Suggested initial value of the number of signatures, passed to dp_activate as initcc.
multi.types	A logical scalar or a character vector. If FALSE, hdp will regard all input spectra as one tumor type.

	If TRUE, hdp will infer tumor types based on the string before "::" in their names. e.g. tumor type for "SA.Syn.Ovary-AdenoCA::S.500" would be "SA.Syn.Ovary-AdenoCA" If multi.types is a character vector, then it should be of the same length as the number of columns in input.catalog, and each value is the name of the tumor type of the corresponding column in input.catalog, e.g. c("SA.Syn.Ovary-AdenoCA","SA.Syn.O
remove.noise	Deprecated; ignored
test.only	If > 0, only analyze the first test.only columns in input.catalog.file.
overwrite	If TRUE overwrite out.dir if it exists, otherwise raise an error.
verbose	If TRUE then message progress information.
num.posterior	Number of posterior sampling chains; can set to 1 for testing.
post.burnin	Pass to hdp_posterior burnin.
post.n	Pass to hdp_posterior n.
post.space	Pass to hdp_posterior space.
post.cpiter	Pass to hdp_posterior cpiter.
post.verbosity	Pass to hdp_posterior verbosity.
cos.merge	The cosine similarity threshhold for merging raw clusters from the posterior sampling chains into "components" i.e. signatures; passed to hdp_extract_components.
min.sample	A "component" (i.e. signature) must have at least this many samples; passed to hdp_extract_components.

Details

Creates several files in out.dir. These are: TODO(Steve): list the files

Value

The same list as returned by RunhdpInternal5.

RunhdpInternal4 *Run hdp extraction and attribution on a spectra catalog file*

Description

Run hdp extraction and attribution on a spectra catalog file

Usage

```
RunhdpInternal4(
  input.catalog,
  CPU.cores = 1,
  seedNumber = 1,
  K.guess,
  multi.types = FALSE,
  verbose = TRUE,
  num.posterior = 4,
  post.burnin = 4000,
  post.n = 50,
```

```
  post.space = 50,
  post.cpiter = 3,
  post.verbosity = 0,
  cos.merge = 0.9,
  min.sample = 1
)
```

Arguments

input.catalog	Input spectra catalog as a matrix or in ICAMS format.
CPU.cores	Number of CPUs to use in running hdp_posterior; this is used to parallize running the posterior sampling chains, so there is no point in making this larger than num.posterior.
seedNumber	An integer that is used to generate separate random seeds for each call to dp_activate, and each call of hdp_posterior; please see the code on how this is done. But repeated calls with same value of seedNumber and other inputs should produce the same results.
K.guess	Suggested initial value of the number of signatures, passed to dp_activate as initcc.
multi.types	A logical scalar or a character vector. If FALSE, hdp will regard all input spectra as one tumor type. If TRUE, hdp will infer tumor types based on the string before "::" in their names. e.g. tumor type for "SA.Syn.Ovary-AdenoCA::S.500" would be "SA.Syn.Ovary-AdenoCA" If multi.types is a character vector, then it should be of the same length as the number of columns in input.catalog, and each value is the name of the tumor type of the corresponding column in input.catalog, e.g. c("SA.Syn.Ovary-AdenoCA","SA.Syn.O
verbose	If TRUE then message progress information.
num.posterior	Number of posterior sampling chains; can set to 1 for testing.
post.burnin	Pass to hdp_posterior burnin.
post.n	Pass to hdp_posterior n.
post.space	Pass to hdp_posterior space.
post.cpiter	Pass to hdp_posterior cpiter.
post.verbosity	Pass to hdp_posterior verbosity.
cos.merge	The cosine similarity threshhold for merging raw clusters from the posterior sampling chains into "components" i.e. signatures; passed to hdp_extract_components.
min.sample	A "component" (i.e. signature) must have at least this many samples; passed to hdp_extract_components.

Value

A list with the following elements:

signature The extracted signature profiles as a matrix; rows are mutation types, columns are samples (e.g. tumors).

exposure The inferred exposures as a matrix of mutation counts; rows are signatures, columns are samples (e.g. tumors).

exposure.p exposure converted to proportions.

multi.chains A hdpSampleMulti-class object. This object has the method chains which returns a list of hdpSampleChain-class objects. Each of these sample chains objects has a method final_hdpState (actually the methods seems to be just hdp) that returns the hdpState from which it was generated.

RunhdpInternal5 *Run hdp extraction and attribution on a spectra catalog file*

Description

Run hdp extraction and attribution on a spectra catalog file

Usage

```
RunhdpInternal5(
  input.catalog,
  CPU.cores = 1,
  seedNumber = 1,
  K.guess,
  multi.types = FALSE,
  verbose = TRUE,
  num.posterior = 4,
  post.burnin = 4000,
  post.n = 50,
  post.space = 50,
  post.cpiter = 3,
  post.verbosity = 0,
  cos.merge = 0.9,
  min.sample = 1
)
```

Arguments

input.catalog	Input spectra catalog as a matrix or in ICAMS format.
CPU.cores	Number of CPUs to use in running hdp_posterior; this is used to parallize running the posterior sampling chains, so there is no point in making this larger than num.posterior.
seedNumber	An integer that is used to generate separate random seeds for each call to dp_activate, and each call of hdp_posterior; please see the code on how this is done. But repeated calls with same value of seedNumber and other inputs should produce the same results.
K.guess	Suggested initial value of the number of signatures, passed to dp_activate as initcc.
multi.types	A logical scalar or a character vector. If FALSE, hdp will regard all input spectra as one tumor type. If TRUE, hdp will infer tumor types based on the string before "::" in their names. e.g. tumor type for "SA.Syn.Ovary-AdenoCA::S.500" would be "SA.Syn.Ovary-AdenoCA" If multi.types is a character vector, then it should be of the same length as the number of columns in input.catalog, and each value is the name of the tumor type of the corresponding column in input.catalog, e.g. c("SA.Syn.Ovary-AdenoCA","SA.Syn.O

verbose	If TRUE then message progress information.
num.posterior	Number of posterior sampling chains; can set to 1 for testing.
post.burnin	Pass to hdp_posterior burnin.
post.n	Pass to hdp_posterior n.
post.space	Pass to hdp_posterior space.
post.cpiter	Pass to hdp_posterior cpiter.
post.verbosity	Pass to hdp_posterior verbosity.
cos.merge	The cosine similarity threshhold for merging raw clusters from the posterior sampling chains into "components" i.e. signatures; passed to hdp_extract_components.
min.sample	A "component" (i.e. signature) must have at least this many samples; passed to hdp_extract_components.

Value

A list with the following elements:

signature The extracted signature profiles as a matrix; rows are mutation types, columns are samples (e.g. tumors).

exposure The inferred exposures as a matrix of mutation counts; rows are signatures, columns are samples (e.g. tumors).

exposure.p exposure converted to proportions.

multi.chains A hdpSampleMulti-class object. This object has the method chains which returns a list of hdpSampleChain-class objects. Each of these sample chains objects has a method final_hdpState (actually the methods seems to be just hdp) that returns the hdpState from which it was generated.

---

xmake.s *Return a function to calculate the unsigned Stirling numbers of the first kind*

Description

Return a function to calculate the unsigned Stirling numbers of the first kind

Usage

```
xmake.s()
```

Value

A function to caluculate a vector of unsigned Stirling numbers, $s(n,k), k = 1...n$, each divided by the maxium Stirling number in the series. The returned function is a closure with state that includes a list of all the unsigned Stirling number series $<=$ the argument, n,

i.e. $[s(1,1)], [s(2,1), s(2,2)], ..., [s(n,1), ..., s(n,n)]$. Memory usage could be substantial, but the stored state does not include the many trailing zeros in the vectors. For this to work within the hdp (https://github.com/nicolaroberts/hdp) package the function returned *must* be called stir.closure.

Index