Package 'mSigHdp'

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Burnin

Run the Gibbs sampling burnin (one thread)

Description

Run the Gibbs sampling burnin (one thread)

Usage

```
Burnin(
  hdp.state,
  seedNumber = 1,
  burnin = 5000,
  cpiter = 3,
  burnin.verbosity = 0,
  burnin.multiplier = 2,
  checkpoint = TRUE
)
```

Arguments

hdp.state An hdpState-class object or a list representation of an hdpState-class

object.

seedNumber Set the random seed to this value.

burnin The number of burn-in iterations in one batch. The total number of burnin iter-

ations is burnin * burnin.multiplier.

cpiter The number of iterations of concentration parameter sampling to perform after

each main Gibbs-sample iteration. (See Teh et al., "Hierarchical Dirichlet Processes", Journal of the American Statistical Association 2006;101(476):1566-

1581 (https://doi.org/10.1198/016214506000000302).)

burnin.verbosity

Verbosity of message statements.

burnin.multiplier

Run burnin.multiplier rounds of burnin iterations. If checkpoint is TRUE, save the burnin chain (see parameter checkpoint.) The diagnostic plot diagnostics.likelihood.pdf can help determine if the chain is stationary. The burnin can be continued from a checkpoint file with ExtendBurnin

(see argument checkpoint).

 $\verb| checkpoint| If \verb| TRUE|, create| a checkpoint| file called mSigHdp.burnin.checkpoint|.seedNumber.Rdata|$

in the current working directory.

Value

A list with 2 elements:

hdplist A list representation of an hdpState-class object.

likelihood A numeric vector with the likelihood at each iteration. This is the same type as returned from link [hdp] {hdp_burnin} in package hdpx.

CombineChainsAndExtractSigs

Extract signatures etc. from multiple Gibbs sample chains

Description

Extract signatures etc. from multiple Gibbs sample chains

Usage

```
CombineChainsAndExtractSigs(
 clean.chlist,
  input.catalog,
  verbose = FALSE,
 high.confidence.prop = 0.9,
 hc.cutoff = 0.1
)
```

Arguments

clean.chlist A list of hdpSampleChain-class S4 objects, each with the information from one Gibbs sampling chain. See hdpSampleChain-class in package hdpx.

input.catalog

Input spectra catalog as a matrix.

verbose If TRUE then message progress information.

high.confidence.prop

Raw clusters of mutations found in >= high.confidence.prop proportion of posterior samples are signatures with high confidence.

hc.cutoff

The cutoff of height of the hierarchical clustering dendrogram used in combining raw clusters of mutations into aggregated clusters.

Value

Invisibly, a list with the following elements:

signature The extracted signature profiles as a matrix; rows are mutation types, columns are signatures with high confidence.

signature.post.samp.number A data frame with two columns. The first column corresponds to each signature in signature and the second columns contains the number of posterior samples that found the raw clusters contributing to the signature.

signature.cdc A numeric data frame. Columns correspond to signatures as in signature. Rows correspond to either biological samples or to parent and grandparent Dirichlet processes.

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exposureProbs The inferred exposures as a matrix of mutation probabilities; rows are signatures, columns are samples (e.g. tumors). This is similar to signature.cdc, but every column was normalized to sum to 1.

low.confidence.signature The profiles of signatures extracted with low confidence as a matrix; rows are mutation types, columns are signatures with < high.confidence.prop of posterior samples.

low.confidence.post.samp.number Analogous to signature.post.samp.number, except that this one is for signatures in low.confidence.signature.

low.confidence.cdc Analogous to signature.cdc, except that columns in this matrix correspond to columns in low.confidence.signature.

extracted.retval A list object returned from extract_components in package hdpx.

ExtendBurnin

Extend burnin iterations generated from Burnin

Description

Extend burnin iterations generated from Burnin

Usage

```
ExtendBurnin(
  previous.burnin.output,
  burnin = 5000,
  cpiter = 3,
  burnin.verbosity = 0,
  seedNumber = NULL
)
```

Arguments

previous.burnin.output

Output from Burnin or the file path of a checkpoint file written by Burnin.

burnin The number of burnin iterations to perform.

cpiter The number of iterations of concentration parameter sampling to perform after

each main Gibbs-sample iteration. (See Teh et al., "Hierarchical Dirichlet Processes", Journal of the American Statistical Association 2006;101(476):1566-

1581 (https://doi.org/10.1198/016214506000000302).)

burnin.verbosity

Number that controls whether progress messages are printed.

seedNumber A random seed for reproducible results.

Value

The same type of object as returned from Burnin.

The envisioned application is extending burnins from burnin checkpoints.

```
GibbsSamplingAfterBurnin
```

Start Gibbs sampling on one chain after burnin

Description

This function might be used to start Gibbs sampling after ExtendBurnin.

Usage

```
GibbsSamplingAfterBurnin(
  burnin.output,
  post.n,
  post.space,
  post.cpiter = 3,
  post.verbosity = 0,
  seedNumber = NULL
)
```

Arguments

burnin.output

A path to burnin checkpoint Rdata or to an S4 object from Burnin.

post.n The number of posterior samples to collect.

post.space The number of iterations between collected samples.

post.cpiter The number of iterations of concentration parameter sampling to perform after

each main Gibbs-sample iteration. (See Teh et al., "Hierarchical Dirichlet Processes", Journal of the American Statistical Association 2006;101(476):1566-

1581 (https://doi.org/10.1198/016214506000000302).)

post.verbosity

Verbosity of debugging statements. No need to change unless for testing or

debugging.

seedNumber A random seed that ensures reproducible results.

Value

An hdpSampleChain S4 object with the salient information from each posterior sample. See hdpSampleChain-class in package hdpx.

RunHdpxParallel Extract (discover) mutational signatures from a matrix of mutational spectra

Description

Please see the vignette for an example.

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Usage

```
RunHdpxParallel(
  input.catalog,
  seedNumber = 123,
 K.quess,
 multi.types = FALSE,
  verbose = FALSE,
 burnin = 1000,
 burnin.multiplier = 10,
  post.n = 200,
 post.space = 100,
 post.cpiter = 3,
 post.verbosity = 0,
  CPU.cores = 20,
  num.child.process = 20,
 high.confidence.prop = 0.9,
 hc.cutoff = 0.1,
  overwrite = TRUE,
  out.dir = NULL,
  gamma.alpha = 1,
  gamma.beta = 20,
  checkpoint = TRUE
```

Arguments

input.catalog

Input spectra catalog as a matrix or in ICAMS format.

seedNumber

A random seed that ensures ensures reproducible results.

K.quess

Suggested initial value of the number of raw clusters. Usually, the number of raw clusters is roughly twice the number of extracted signatures. Passed to hdpx::dp_activate as argument initcc.

multi.types

A logical scalar or a character vector.

If FALSE, The HDP analysis will regard all input spectra as one tumor type, and the HDP structure will have one parent node for all tumors.

If TRUE, Sample IDs in input.catalog must have the form sample_type::sample_id. If a character vector, then its length must be ncol (input.catalog), and

each value is the sample type of the corresponding column in input.catalog, e.g. c(rep("Type-A", 23), rep("Type-B", 10)) for 23 Type-A sam-

ples and 10 Type-B samples.

If not FALSE, HDP will have one parent node for each sample type and one grandparent node.

verbose

If TRUE then message progress information.

burnin

The number of burn-in iterations in one batch. The total number of burnin iterations is burnin * burnin.multiplier.

burnin.multiplier

Run burnin.multiplier rounds of burnin iterations. If checkpoint is TRUE, save the burnin chain (see parameter checkpoint.) The diagnostic plot diagnostics.likelihood.pdf can help determine if the chain is stationary. The burnin can be continued from a checkpoint file with ExtendBurnin (see argument checkpoint).

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post.n The number of posterior samples to collect. The number of iterations between collected samples. post.space post.cpiter The number of iterations of concentration parameter samplings to perform after each iteration. post.verbosity Verbosity of debugging statements. No need to change except for development purposes. CPU.cores Number of CPUs to use; this should be no more than num.child.process. num.child.process Number of posterior sampling chains; can set to 1 for testing. We recommend 20 for real data analysis high.confidence.prop Raw clusters of mutations found in >= high.confidence.prop proportion of posterior samples are signatures with high confidence. The cutoff of height of the hierarchical clustering dendrogram used in combining hc.cutoff raw clusters of mutations into aggregated clusters. If TRUE overwrite out .dir if it exists, otherwise raise an error. overwrite out.dir If not NULL then a character string specifying a directory that will be created for the output, including csv files and plots (pdfs) of extracted signatures and their exposures. If NULL no directory will be created and no files will be generated. Shape parameter of the gamma distribution prior for the Dirichlet process congamma.alpha

centration parameters α_0 and all α_j in Figure B.1 of

 https://www.repository.cam.ac.uk/bitstream/handle/1810/275454/Roberts-2018-PhD.pdf

gamma.beta

Inverse scale parameter (rate parameter) of the gamma distribution prior for the Dirichlet process concentration parameters: β_0 and all β_i in Figure B.1 of

 https://www.repository.cam.ac.uk/bitstream/handle/1810/275454/Roberts-2018-PhD.pdf

We recommend gamma.alpha = 1 and gamma.beta = 20 for single-base-substitution signature extraction; gamma.alpha = 1 and gamma.beta = 50 for doublet-basesubstitution and indel signature extraction

checkpoint If TRUE, then

- Checkpoint each final Gibbs sample chain to the current working directory, in a file called mSigHdp.sample.checkpoint.x.Rdata, where x depends on seedNumber.
- Periodically checkpoint the burnin state to the current working directory, in files called mSigHdp.burnin.checkpoint.x.Rdata, where x depends on the seedNumber.

Value

Invisibly, a list with the following elements:

signature The extracted signature profiles as a matrix; rows are mutation types, columns are signatures with high confidence.

signature.post.samp.number A data frame with two columns. The first column corresponds to each signature in signature and the second columns contains the number of posterior samples that found the raw clusters contributing to the signature.

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signature.cdc A numeric data frame. Columns correspond to signatures as in signature. Rows correspond to either biological samples or to parent and grandparent Dirichlet processes.

- **exposureProbs** The inferred exposures as a matrix of mutation probabilities; rows are signatures, columns are samples (e.g. tumors). This is similar to signature.cdc, but every column was normalized to sum to 1.
- **low.confidence.signature** The profiles of signatures extracted with low confidence as a matrix; rows are mutation types, columns are signatures with < high.confidence.prop of posterior samples.
- low.confidence.post.samp.number Analogous to signature.post.samp.number, except
 that this one is for signatures in low.confidence.signature.
- **low.confidence.cdc** Analogous to signature.cdc, except that columns in this matrix correspond to columns in low.confidence.signature.
- extracted.retval A list object returned from extract_components in package hdpx.

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