

# Package ‘nonparametricSummaryPSM’

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**Type** Package

**Title** Nonparametric methods to find optimal weights to combine posterior similarity matrices

**Version** 0.1.0

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**Description** The purpose of this code is to compute summary PSMs (posterior similarity matrices) from a set of multiple PSMs obtained for instance by means of subsampling. This implements the Dirichlet process and Pitman-Yor process based methods for combining PSMs proposed in Strauss et al. (2019). GPpseudoClust: deconvolution of shared pseudo-trajectories at single-cell resolution.

**License** GPL-3.0

**Encoding** UTF-8

**LazyData** true

**RoxygenNote** 6.1.1

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checkConvergence	<i>checkConvergence</i>
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## Description

checkConvergence

## Usage

checkConvergence(PSMs)

**Arguments**

PSMs                      3-dimensional array of PSMs, for each  $j$  PSMs[,j] is the PSM of subsampled chain  $j$ ,  $j = 1, \dots, m$

**Value**

results: list of length  $m-1$  of results of processPSM function applied to the following subsets of PSMs: 1,2; 1,2,3; 1,2,3,4; ... 1,2,...,m

coph: vector of length  $m-1$  of cophenetic correlation coefficients measuring how well the summary PSM obtained using the PY+PEAR method on the first  $k$  submatrices ( $k=1, \dots, l$ ) is described by a hierarchical clustering tree for a summary PSM based on the first  $l$  subsampled chains

coph\_DP: as coph, but for summary PSMs obtained by the DPM+PEAR method

distPY: vector of length  $m-2$  of Frobenius (Euclidean) norm of distances between consecutive summary PSMs, that is summary PSMs obtained from chains 1,2,...,k,k+1 and 1,2,...,k (PY+PEAR)

distDP: as distPY, but for DPM+PEAR method

**Author(s)**

Magdalena Strauss

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computeSumClustPEAR      *computeSumClustPear*

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**Description**

Compute summary PSMs (posterior similarity matrices) from a set of multiple PSMs obtained for instance by means of subsampling. This implements the Dirichlet process and Pitman-Yor process based methods for combining PSMs proposed in Strauss et al. Unravelling shared pseudo-trajectories at single-cell resolution. Internal function

**Usage**

```
computeSumClustPEAR(PSM, maxCl = 10)
```

**Arguments**

PSM                      posterior similarity matrix

**Value**

Summary clustering computed using the PEAR criterion (Fritsch and Ickstadt, 2009, using the mclust package (Fritsch, 2012))

**Author(s)**

Magdalena Strauss

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`computeWeightsSumClust`*computeWeightsSumClust*

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**Description**

Internal function

**Usage**

`computeWeightsSumClust(allocs)`

**Value**

PSM and summary clustering obtained from Pitman-Yor process with `allocs` as input, weights used to compute the summary PSM from the individual PSMs

**Author(s)**

Magdalena Strauss

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`computeWeightsSumClustDPM`*computeWeightsSumClustDPM*

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**Description**

Internal function

**Usage**

`computeWeightsSumClustDPM(allocs)`

**Value**

PSM and summary clustering obtained from Dirichlet process with `allocs` as input, weights used to compute the summary PSM from the individual PSMs

**Author(s)**

Magdalena Strauss

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processPSMs

*processPSMs*


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### Description

processPSMs

### Usage

processPSMs(PSMs)

### Arguments

PSMs                      3-dimensional array of PSMs, for each j PSMs[,j] is the PSM of subsampled chain j

### Value

weightedPSM: weighted summary PSM obtained using a Pitman-Yor process mixture model with variable selection

sumClustPEAR: final summary clustering obtained from weightedPSM using the PEAR criterion

weightedPSM\_DP weighted summary PSM obtained using a Dirichlet process mixture model with variable selection

sumClustPEAR\_DP: final summary clustering obtained from weightedPSM\_DP using the PEAR criterion

weights: weights which were used for the computation of the summary PSM (Pitman-Yor based model)

weights\_DP: weights which were used for the computation of the summary PSM (Dirichlet based model)

### Author(s)

Magdalena Strauss

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RhatConcentration

*RhatConcentration*


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### Description

RhatConcentration

### Usage

RhatConcentration(concentrationSamples, nChainsTest)

**Arguments**

`concentrationSamples`:  
nIterations x nChains matrix of nIterations samples of the concentration parameter alpha for each of nChains subsampled chains

`nChainsTest`: number of subsamples for which we want to test if the number of sufficient for convergence

**Value**

GR-statistics across subsampled chains (see Strauss et al. 2019) for groups of

**Author(s)**

Magdalena Strauss

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