Package 'fabMix'

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Type Package

Title Overfitting Parsimonious Bayesian Mixtures of Factor Analyzers with an Unknown Number of Components

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Description Model-based clustering of multivariate continuous data using overfitting Bayesian mixtures of factor analyzers (Papastamoulis, 2018 CSDA). Suitable prior assumptions ensure that asymptotically the extra components will have zero posterior weight (empty), therefore, the inference on the number of clusters is based on the ``alive" components. The number of factors is considered fixed, and the optimal one can be estimated using information criteria. Eight parameterizations are available, namely the ``UUU", ``UCU", ``UCC", ``CUU", ``CUC", ``CCC", ``CCU" and ``UUC" models (see McNicholas et al, 2008 Stat Comp). Identifiability issues related to label switching are dealt by post-processing the simulated output with the ECR algorithm (Papastamoulis 2010 JCGS, 2016 JSS). Missing values are currently allowed for the UUU and UCU model parameterizations.

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Imports Rcpp (>= 0.12.17), MASS, doParallel, foreach, label.switching, mvtnorm, doRNG, RColorBrewer, corrplot, mclust

LinkingTo Rcpp, RcppArmadillo

NeedsCompilation yes

R topics documented:

`fabMix-package` *Overfitting Parsimonious Bayesian Mixtures of Factor Analyzers with an Unknown Number of Components*

Description

Model-based clustering of multivariate continuous data using overfitting Bayesian mixtures of factor analyzers (Papastamoulis, 2018 CSDA). Suitable prior assumptions ensure that asymptotically the extra components will have zero posterior weight (empty), therefore, the inference on the number of clusters is based on the "alive" components. The number of factors is considered fixed, and the optimal one can be estimated using information criteria. Eight parameterizations are available, namely the "UUU", "UCU", "UCC", "CUU", "CUC", "CCC", "CCU" and "UUC" models (see McNicholas et al, 2008 Stat Comp). Identifiability issues related to label switching are dealt by post-processing the simulated output with the ECR algorithm (Papastamoulis 2010 JCGS, 2016 JSS). Missing values are currently allowed for the UUU and UCU model parameterizations.

Author(s)

Panagiotis Papastamoulis

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References

Fokoue, E. and Titterington, D.M. (2003). Mixtures of Factor Analysers: Bayesian Estimation and Inference by Stochastic Simulation. Machine Learing, 50(1): 73-94.

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van Havre, Z., White, N., Rousseau, J. and Mengersen, K. (2015). Overfitting Bayesian Mixture Models with an Unknown Number of Components. PLOS ONE, 10(7): 1-27.

Papastamoulis, P. (2016). `label.switching`: An R Package for Dealing with the Label Switching Problem in MCMC Outputs. Journal of Statistical Software, 69(1), 1-24.

Papastamoulis, P. (2018). Overfitting Bayesian mixtures of factor analyzers with an unknown number of components. Computational Statistics and Data Analysis, 124: 220-234. DOI: 10.1016/j.csda.2018.03.007.

See Also

fabMix

Examples

```
# simulate a synthetic dataset along the lines of the paper:
n = 1000              # sample size
p = 40                # number of variables
q = 4                 # number of factors
K = 10                # number of clusters
sINV_diag = 1/((1:p)) # diagonal of inverse variance of errors
set.seed(10)
syntheticDataset <- simData(K.true = K, n = n, q = q, p = p, sINV_values = sINV_diag )

## Not run:

# define parameters
Kmax <- 20    # number of overfitted mixture components
nChains <- 8  # number of parallel chains
dN <- 1
# Dirichlet prior of mixture weights per chain.
#   The target chain corresponds to the first entry.
dirPriorAlphas <- c(1, 1 + dN * (2:nChains - 1))/Kmax
outputFolder <- "fabMixExample"
# Run algorithm
fabMix( dirPriorAlphas = dirPriorAlphas,
        rawData = syntheticDataset$data,
        outDir = outputFolder, Kmax = Kmax, mCycles = 1200,
        burnCycles = 200, q = q)

# Compute information criteria:
getStuffForDIC(x_data = syntheticDataset$data, outputFolder = outputFolder, q = q)

# Deal with label switching:
dealWithLabelSwitching(x_data = syntheticDataset$data,
        outputFolder = outputFolder, q = q,
        compute_regularized_expression = TRUE, Km = Kmax)

## End(Not run)
```

---

complete.log.likelihood

Complete log-likelihood function

Description

Complete log-likelihood function

Usage

```
complete.log.likelihood(x_data, w, mu, Lambda, SigmaINV, z)
```

Arguments

x_data	Data
w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
SigmaINV	Precision matrix (inverse covariance)
z	Allocation vector of the data to the mixture components

Value

complete log-likelihood value

Author(s)

Panagiotis Papastamoulis

---

complete.log.likelihood_q0

Complete log-likelihood function for $q = 0$

Description

Complete log-likelihood function

Usage

```
complete.log.likelihood_q0(x_data, w, mu, SigmaINV, z)
```

Arguments

x_data	Data
w	Mixture weights
mu	Marginal means
SigmaINV	Precision matrix (inverse covariance)
z	Allocation vector of the data to the mixture components

Value

complete log-likelihood value

Author(s)

Panagiotis Papastamoulis

complete.log.likelihood_q0_sameSigma

Complete log-likelihood function for $q = 0$

Description

Complete log-likelihood function

Usage

```
complete.log.likelihood_q0_sameSigma(x_data, w, mu, SigmaINV, z)
```

Arguments

x_data	Data
w	Mixture weights
mu	Marginal means
SigmaINV	Precision matrix (inverse covariance)
z	Allocation vector of the data to the mixture components

Value

complete log-likelihood value

Author(s)

Panagiotis Papastamoulis

complete.log.likelihood_Sj

Complete log-likelihood function

Description

Complete log-likelihood function

Usage

```
complete.log.likelihood_Sj(x_data, w, mu, Lambda, SigmaINV, z)
```

Arguments

x_data	Data
w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
SigmaINV	Precision matrix (inverse covariance) per component
z	Allocation vector of the data to the mixture components

Value

complete log-likelihood value

Author(s)

Panagiotis Papastamoulis

---

compute_A_B_G_D_and_simulate_mu_Lambda

Computation and simulations

Description

This function simulates μ and Λ.

Usage

```
compute_A_B_G_D_and_simulate_mu_Lambda(SigmaINV,
suff_statistics, OmegaINV, K, priorConst1, T_INV, v_r)
```

Arguments

SigmaINV	Precision matrix Σ^{-1}
suff_statistics	Sufficient statistics
OmegaINV	Prior parameter: Ω^{-1}
K	Number of overfitting mixture components
priorConst1	Prior constant: $T^{-1}\xi$
T_INV	Prior parameter: $T^{-1}\xi$
v_r	This vector is used to set to zero the upper right $(q-1)\times(q-1)$ diagonal block of factor loadings for identifiability purposes.

Value

A list containing A, B, Γ, Δ and a draw from the conditional distributions of μ and Λ.

Author(s)

Panagiotis Papastamoulis

`compute_A_B_G_D_and_simulate_mu_Lambda_CCU`
Computation and simulations for CCU

Description

This function simulates μ and Λ for the CCU model.

Usage

```
compute_A_B_G_D_and_simulate_mu_Lambda_CCU(SigmaINV,
suff_statistics, OmegaINV, K, priorConst1, T_INV, v_r)
```

Arguments

`SigmaINV`	Precision matrix Σ^{-1}
`suff_statistics`	Sufficient statistics
`OmegaINV`	Prior parameter: Ω^{-1}
`K`	Number of overfitting mixture components
`priorConst1`	Prior constant: $T^{-1}\xi$
`T_INV`	Prior parameter: $T^{-1}\xi$
`v_r`	This vector is used to set to zero the upper right $(q-1)\times(q-1)$ diagonal block of factor loadings for identifiability purposes.

Value

A list containing A, B, Γ, Δ and a draw from the conditional distributions of μ and Λ.

Author(s)

Panagiotis Papastamoulis

`compute_A_B_G_D_and_simulate_mu_Lambda_CUU` *Computation and simulations for CUU*

Description

This function simulates μ and Λ for the CUU model.

Usage

```
compute_A_B_G_D_and_simulate_mu_Lambda_CUU(SigmaINV,
suff_statistics, OmegaINV, K, priorConst1, T_INV, v_r)
```

Arguments

`SigmaINV`	Precision matrix Σ^{-1}
`suff_statistics`	Sufficient statistics
`OmegaINV`	Prior parameter: Ω^{-1}
`K`	Number of overfitting mixture components
`priorConst1`	Prior constant: $T^{-1}\xi$
`T_INV`	Prior parameter: $T^{-1}\xi$
`v_r`	This vector is used to set to zero the upper right $(q-1)\times(q-1)$ diagonal block of factor loadings for identifiability purposes.

Value

A list containing A, B, Γ, Δ and a draw from the conditional distributions of μ and Λ.

Author(s)

Panagiotis Papastamoulis

`compute_A_B_G_D_and_simulate_mu_Lambda_q0` *Computation and simulations for $q = 0$.*

Description

This function simulates μ.

Usage

```
compute_A_B_G_D_and_simulate_mu_Lambda_q0(SigmaINV,
suff_statistics, K, priorConst1, T_INV, v_r)
```

Arguments

SigmaINV	Precision matrix Σ^{-1}
suff_statistics	Sufficient statistics
K	Number of overfitting mixture components
priorConst1	Prior constant: $T^{-1}\xi$
T_INV	Prior parameter: $T^{-1}\xi$
v_r	This vector is used to set to zero the upper right $(q-1)\times(q-1)$ diagonal block of factor loadings for identifiability purposes.

Value

A list containing A, B, Γ, Δ and a draw from the conditional distributions of μ and Λ.

Author(s)

Panagiotis Papastamoulis

---

compute_A_B_G_D_and_simulate_mu_Lambda_q0_sameSigma

Computation and simulations for $q = 0$.

Description

This function simulates μ.

Usage

```
compute_A_B_G_D_and_simulate_mu_Lambda_q0_sameSigma(SigmaINV,
suff_statistics, K, priorConst1, T_INV, v_r)
```

Arguments

SigmaINV	Precision matrix Σ^{-1}
suff_statistics	Sufficient statistics
K	Number of overfitting mixture components
priorConst1	Prior constant: $T^{-1}\xi$
T_INV	Prior parameter: $T^{-1}\xi$
v_r	This vector is used to set to zero the upper right $(q-1)\times(q-1)$ diagonal block of factor loadings for identifiability purposes.

Value

A list containing A, B, Γ, Δ and a draw from the conditional distributions of μ and Λ.

Author(s)

Panagiotis Papastamoulis

`compute_A_B_G_D_and_simulate_mu_Lambda_Sj`
Computation and simulations

Description

This function simulates μ and Λ.

Usage

```
compute_A_B_G_D_and_simulate_mu_Lambda_Sj(SigmaINV,
suff_statistics, OmegaINV, K, priorConst1, T_INV, v_r)
```

Arguments

`SigmaINV`	Precision matrix Σ^{-1} per component
`suff_statistics`	Sufficient statistics
`OmegaINV`	Prior parameter: Ω^{-1}
`K`	Number of overfitting mixture components
`priorConst1`	Prior constant: $T^{-1}\xi$
`T_INV`	Prior parameter: $T^{-1}\xi$
`v_r`	This vector is used to set to zero the upper right $(q-1)\times(q-1)$ diagonal block of factor loadings for identifiability purposes.

Value

A list containing A, B, Γ, Δ and a draw from the conditional distributions of μ and Λ.

Author(s)

Panagiotis Papastamoulis

`compute_sufficient_statistics` *Compute sufficient statistics*

Description

Compute sufficient statistics given y and z.

Usage

```
compute_sufficient_statistics(y, z, K, x_data)
```

Arguments

`y`	Matrix of factors
`z`	Allocation vector
`K`	Number of components
`x_data`	Data

Value

A list with six entries of sufficient statistics.

Author(s)

Panagiotis Papastamoulis

`compute_sufficient_statistics_given_mu` *Compute sufficient statistics given mu*

Description

Compute sufficient statistics given y and z.

Usage

```
compute_sufficient_statistics_given_mu(y, z, K, x_data,mu)
```

Arguments

`y`	Matrix of factors
`z`	Allocation vector
`K`	Number of components
`x_data`	Data
`mu`	Means per component

Value

A list with six entries of sufficient statistics.

Author(s)

Panagiotis Papastamoulis

`compute_sufficient_statistics_q0`
Compute sufficient statistics for $q = 0$

Description

Compute sufficient statistics given z.

Usage

```
compute_sufficient_statistics_q0(z, K, x_data)
```

Arguments

`z`	Allocation vector
`K`	Number of components
`x_data`	Data

Value

A list with six entries of sufficient statistics.

Author(s)

Panagiotis Papastamoulis

`dealWithLabelSwitching` *Apply label switching algorithms*

Description

This functions is a wrapper for the `label.switching` package and applies the `ECR` and `ECR.ITERATIVE.1` algorithms. The model may have the same variance of error terms per cluster or not.

Usage

```
dealWithLabelSwitching(sameSigma, x_data, outputFolder, q, burn,
z.true, compute_regularized_expression, Km)
```

Arguments

`sameSigma`	Logical value indicating whether the parameterization with the same error variance per cluster is used.
`x_data`	Data
`outputFolder`	Name of the folder where the `fabMix` function has saved its output
`q`	Number of factors
`burn`	Discard observations as burn-in period (optional).
`z.true`	An (optional) vector of cluster assignments which is considered as the grountruth clustering of the data. Useful for direct comparisons against the real parameter values in simulated data.
`compute_regularized_expression`	Logical. Should regularized expression be computed?
`Km`	Number of components in the overfitted mixture model.

Value

The following files are produced in the output folder:

Author(s)

Panagiotis Papastamoulis

References

Papastamoulis, P. (2016). `label.switching`: An R Package for Dealing with the Label Switching Problem in MCMC Outputs. Journal of Statistical Software, 69(1), 1-24.

fabMix *Main function*

Description

This function runs parallel chains under a prior tempering scheme of the Dirichlet prior distribution of mixture weights.

Usage

```
fabMix(model, dirPriorAlphas, rawData, outDir, Kmax, mCycles,
burnCycles, g, h, alpha_sigma, beta_sigma, q, normalize,
thinning, zStart, nIterPerCycle, gibbs_z,
warm_up_overfitting, warm_up, overfittingInitialization,
progressGraphs, gwar)
```

Arguments

model	Any subset of "UUU" "CUU" "UCU" "CCU" "UCC" "UUC" "CUC", "CCC" indicating the fitted models.
dirPriorAlphas	The prior Dirichlet parameters for each chain.
rawData	The observed data as an $n \times p$ matrix. Clustering is performed on the rows of the matrix.
outDir	Name of the output folder.
Kmax	Number of components in the overfitted mixture. Default: 20.
mCycles	Number of MCMC cycles. Each cycle consists of nIterPerCycle MCMC iterations. At the end of each cycle a swap of the state of two randomly chosen adjacent chains is attempted.
burnCycles	Number of cycles that will be discarded as burn-in period.
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
q	A vector containing the number of factors to be fitted.
normalize	Should the observed data be normalized? Default value: TRUE.
thinning	Optional integer denoting the thinning of the keeped MCMC cycles.
zStart	Optional starting value for the allocation vector.
nIterPerCycle	Number of iteration per MCMC cycle. Default value: 10.
gibbs_z	Select the gibbs sampling scheme for updating latent allocations of mixture model. Default value: 1.
warm_up_overfitting	Number of iterations for the overfitting initialization scheme. Default value: 100.

warm_up Number of iterations that will be used to initialize the models before starting proposing switchings. Default value: 500.

overfittingInitialization
Logical value indicating whether the chains are initialized via the overfitting initialization scheme. Default: TRUE.

progressGraphs Logical value indicating whether to plot successive states of the chains while the sampler runs. Default: FALSE.

gwar Initialization parameter. Default: 0.05.

Value

An object of class fabMix.object, that is, a list consisting of the following entries:

bic Bayesian Information Criterion per model and number of factors.

class The estimated single best clustering of the observations according to the selected model.

n_Clusters_per_model
The most probable number of clusters (number of non-empty components of the overfitted mixture) per model and number of factors.

posterior_probability
The posterior probability of the estimated allocations according to the selected model.

covariance_matrix
The estimated posterior mean of the covariance matrix per cluster according to the selected model.

mu The estimated posterior mean of the mean per cluster according to the selected model.

weights The estimated posterior mean of the mixing proportions according to the selected model.

selected_model
Data frame containing the parameterization, number of clusters and factors of the selected model.

mcmc A list containing the MCMC draws for the parameters of the selected models. All component-specific parameters have been reordered according to the ECR algorithm in order to undo the label switching problem.

data The observed data.

Note

It is recommended to always use: normalize = TRUE (default). Tuning of dirPriorAlphas may be necessary to achieve reasonable acceptance rates of chain swaps. Note that the output is reordered in order to deal with the label switching problem, according to the ECR algorithm applied by dealWithLabelSwitching function.

Author(s)

Panagiotis Papastamoulis

References

Papastamoulis, P. (2018). Overfitting Bayesian mixtures of factor analyzers with an unknown number of components. Computational Statistics and Data Analysis, 124: 220-234. DOI: 10.1016/j.csda.2018.03.007.

See Also

dealWithLabelSwitching

Examples

```
## Not run:
# simulate a synthetic dataset along the lines of the paper:
n = 1000              # sample size
p = 40                # number of variables
q = 4                 # number of factors
K = 10                # number of clusters
sINV_diag = 1/((1:p)) # diagonal of inverse variance of errors
set.seed(10)
syntheticDataset <- simData(K.true = K, n = n, q = q, p = p, sINV_values = sINV_diag )

# define parameters
Kmax <- 20    # number of overfitted mixture components
nChains <- 8  # number of parallel chains
dN <- 1
# Dirichlet prior of mixture weights per chain.
#   The target chain corresponds to the first entry.
dirPriorAlphas <- c(1, 1 + dN * (2:nChains - 1))/Kmax
outputFolder <- "fabMixExample"
# Run algorithm
fabMix( dirPriorAlphas = dirPriorAlphas,
        rawData = syntheticDataset$data,
        outDir = outputFolder, Kmax = Kmax, mCycles = 1200,
        burnCycles = 200, q = q)

## End(Not run)
```

---

fabMix_CxC *Main function of the package for CUC, CCC models*

Description

This function runs parallel chains under a prior tempering scheme of the Dirichlet prior distribution of mixture weights.

Usage

```
fabMix_CxC(sameSigma, dirPriorAlphas, rawData, outDir, Kmax, mCycles,
burnCycles, g, h, alpha_sigma, beta_sigma, q, normalize,
thinning, zStart, nIterPerCycle, gibbs_z,
warm_up_overfitting, warm_up, overfittingInitialization,
progressGraphs, gwar, cccStart)
```

Arguments

sameSigma	Logical value denoting the parameterization of the error variance per component. If TRUE, the parameterization CCU is fitted. Otherwise, the parameterization CUU is fitted.
dirPriorAlphas	The prior Dirichlet parameters for each chain.
rawData	The observed data as an $n \times p$ matrix. Clustering is performed on the rows of the matrix.
outDir	Name of the output folder.
Kmax	Number of components in the overfitted mixture. Default: 20.
mCycles	Number of MCMC cycles. Each cycle consists of nIterPerCycle MCMC iterations. At the end of each cycle a swap of the state of two randomly chosen adjacent chains is attempted.
burnCycles	Number of cycles that will be discarded as burn-in period.
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
q	Number of factors q, where $1 \leq q \leq L$. An error is thrown if the Ledermann bound (L) is exceeded.
normalize	Should the observed data be normalized? Default value: TRUE.
thinning	Optional integer denoting the thinning of the keeped MCMC cycles.
zStart	Optional starting value for the allocation vector.
nIterPerCycle	Number of iteration per MCMC cycle. Default value: 10.
gibbs_z	Select the gibbs sampling scheme for updating latent allocations of mixture model. Default value: 1.
warm_up_overfitting	Number of iterations for the overfitting initialization scheme. Default value: 100.
warm_up	Number of iterations that will be used to initialize the models before starting proposing switchings. Default value: 500.
overfittingInitialization	Logical value indicating whether the chains are initialized via the overfitting initialization scheme. Default: TRUE.
progressGraphs	Logical value indicating whether to plot successive states of the chains while the sampler runs. Default: FALSE.
gwar	Initialization parameter. Default: 0.05.
cccStart	Initialization from the CCC model.

Value

List of files written to `outDir`

Note

It is recommended to always use: `normalize = TRUE` (default). Tuning of `dirPriorAlphas` may be necessary to achieve reasonable acceptance rates of chain swaps. Also note that the output is not identifiable due to label switching and the user has to subsequently call the `dealWithLabelSwitching` function.

Author(s)

Panagiotis Papastamoulis

References

Papastamoulis, P. (2017). Overfitting Bayesian mixtures of factor analyzers with an unknown number of components. arXiv:1701.04605 [stat.ME]

See Also

`dealWithLabelSwitching`

Examples

```
## Not run:
# simulate a synthetic dataset along the lines of the paper:
n = 1000              # sample size
p = 40                # number of variables
q = 4                 # number of factors
K = 10                # number of clusters
sINV_diag = 1/((1:p)) # diagonal of inverse variance of errors
set.seed(10)
syntheticDataset <- simData(K.true = K, n = n, q = q, p = p, sINV_values = sINV_diag )

# define parameters
Kmax <- 20    # number of overfitted mixture components
nChains <- 8  # number of parallel chains
dN <- 1
# Dirichlet prior of mixture weights per chain.
#   The target chain corresponds to the first entry.
dirPriorAlphas <- c(1, 1 + dN * (2:nChains - 1))/Kmax
outputFolder <- "fabMixExample"
# Run algorithm
fabMix_CxC( dirPriorAlphas = dirPriorAlphas,
        rawData = syntheticDataset$data,
        outDir = outputFolder, Kmax = Kmax, mCycles = 1200,
        burnCycles = 200, q = q)

# Compute information criteria:
#getStuffForDIC(x_data = syntheticDataset$data, outputFolder = outputFolder, q = q)
```

```
# Deal with label switching:
#dealWithLabelSwitching(x_data = syntheticDataset$data,
#        outputFolder = outputFolder, q = q,
#        compute_regularized_expression = TRUE, Km = Kmax)

## End(Not run)
```

`fabMix_CxU` *Main function of the package for CCU, CUU models*

Description

This function runs parallel chains under a prior tempering scheme of the Dirichlet prior distribution of mixture weights.

Usage

```
fabMix_CxU(sameSigma, dirPriorAlphas, rawData, outDir, Kmax, mCycles,
burnCycles, g, h, alpha_sigma, beta_sigma, q, normalize,
thinning, zStart, nIterPerCycle, gibbs_z,
warm_up_overfitting, warm_up, overfittingInitialization,
progressGraphs, gwar)
```

Arguments

`sameSigma`	Logical value denoting the parameterization of the error variance per component. If `TRUE`, the parameterization CCU is fitted. Otherwise, the parameterization CUU is fitted.
`dirPriorAlphas`	The prior Dirichlet parameters for each chain.
`rawData`	The observed data as an $n \times p$ matrix. Clustering is performed on the rows of the matrix.
`outDir`	Name of the output folder.
`Kmax`	Number of components in the overfitted mixture. Default: 20.
`mCycles`	Number of MCMC cycles. Each cycle consists of `nIterPerCycle` MCMC iterations. At the end of each cycle a swap of the state of two randomly chosen adjacent chains is attempted.
`burnCycles`	Number of cycles that will be discarded as burn-in period.
`g`	Prior parameter g. Default value: $g = 2$.
`h`	Prior parameter h. Default value: $h = 1$.
`alpha_sigma`	Prior parameter α. Default value: $\alpha = 2$.
`beta_sigma`	Prior parameter β. Default value: $\beta = 1$.

q	Number of factors q, where $1 \leq q \leq L$. An error is thrown if the Ledermann bound (L) is exceeded.
normalize	Should the observed data be normalized? Default value: TRUE.
thinning	Optional integer denoting the thinning of the keeped MCMC cycles.
zStart	Optional starting value for the allocation vector.
nIterPerCycle	Number of iteration per MCMC cycle. Default value: 10.
gibbs_z	Select the gibbs sampling scheme for updating latent allocations of mixture model. Default value: 1.
warm_up_overfitting	Number of iterations for the overfitting initialization scheme. Default value: 100.
warm_up	Number of iterations that will be used to initialize the models before starting proposing switchings. Default value: 500.
overfittingInitialization	Logical value indicating whether the chains are initialized via the overfitting initialization scheme. Default: TRUE.
progressGraphs	Logical value indicating whether to plot successive states of the chains while the sampler runs. Default: FALSE.
gwar	Initialization parameter. Default: 0.05.

Value

List of files written to outDir

Note

It is recommended to always use: normalize = TRUE (default). Tuning of dirPriorAlphas may be necessary to achieve reasonable acceptance rates of chain swaps. Also note that the output is not identifiable due to label switching and the user has to subsequently call the dealWithLabelSwitching function.

Author(s)

Panagiotis Papastamoulis

References

Papastamoulis, P. (2017). Overfitting Bayesian mixtures of factor analyzers with an unknown number of components. arXiv:1701.04605 [stat.ME]

See Also

dealWithLabelSwitching

Examples

```
## Not run:
# simulate a synthetic dataset along the lines of the paper:
n = 1000              # sample size
p = 40                # number of variables
q = 4                 # number of factors
K = 10                # number of clusters
sINV_diag = 1/((1:p)) # diagonal of inverse variance of errors
set.seed(10)
syntheticDataset <- simData(K.true = K, n = n, q = q, p = p, sINV_values = sINV_diag )

# define parameters
Kmax <- 20    # number of overfitted mixture components
nChains <- 8  # number of parallel chains
dN <- 1
# Dirichlet prior of mixture weights per chain.
#   The target chain corresponds to the first entry.
dirPriorAlphas <- c(1, 1 + dN * (2:nChains - 1))/Kmax
outputFolder <- "fabMixExample"
# Run algorithm
fabMix_CxU( dirPriorAlphas = dirPriorAlphas,
        rawData = syntheticDataset$data,
        outDir = outputFolder, Kmax = Kmax, mCycles = 1200,
        burnCycles = 200, q = q)

# Compute information criteria:
#getStuffForDIC(x_data = syntheticDataset$data, outputFolder = outputFolder, q = q)

# Deal with label switching:
#dealWithLabelSwitching(x_data = syntheticDataset$data,
#        outputFolder = outputFolder, q = q,
#        compute_regularized_expression = TRUE, Km = Kmax)

## End(Not run)
```

`fabMix_missing_values` *Main function for the case of missing values*

Description

This function runs parallel chains under a prior tempering scheme of the Dirichlet prior distribution of mixture weights. Missing values are simulated from their full conditional posterior distribution.

Usage

```
fabMix_missing_values(sameSigma, dirPriorAlphas, rawData, outDir, Kmax, mCycles,
burnCycles, g, h, alpha_sigma, beta_sigma, q, normalize,
thinning, zStart, nIterPerCycle, gibbs_z, warm_up,
progressGraphs, gwar)
```

Arguments

sameSigma	Logical value denoting the parameterization of the error variance per component. If TRUE, the parameterization $\Sigma_1 = \ldots = \Sigma_K$ is fitted.
dirPriorAlphas	The prior Dirichlet parameters for each chain.
rawData	The observed data as an $n \times p$ matrix. Clustering is performed on the rows of the matrix.
outDir	Name of the output folder.
Kmax	Number of components in the overfitted mixture. Default: 20.
mCycles	Number of MCMC cycles. Each cycle consists of nIterPerCycle MCMC iterations. At the end of each cycle a swap of the state of two randomly chosen adjacent chains is attempted.
burnCycles	Number of cycles that will be discarded as burn-in period.
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
q	Number of factors q, where $1 \leq q \leq L$. An error is thrown if the Ledermann bound (L) is exceeded.
normalize	Should the observed data be normalized? Default value: TRUE.
thinning	Optional integer denoting the thinning of the keeped MCMC cycles.
zStart	Optional starting value for the allocation vector.
nIterPerCycle	Number of iteration per MCMC cycle. Default value: 10.
gibbs_z	Select the gibbs sampling scheme for updating latent allocations of mixture model. Default value: 1.
warm_up	NUmber of iterations that will be used to initialize the models before starting proposing switchings. Default value: 500.
progressGraphs	Logical value indicating whether to plot successive states of the chains while the sampler runs. Default: FALSE.
gwar	Initialization parameter. Default: 0.05.

Value

List of files written to outDir

Note

It is recommended to always use: normalize = TRUE (default). Tuning of dirPriorAlphas may be necessary to achieve reasonable acceptance rates of chain swaps. Also note that the output is not identifiable due to label switching and the user has to subsequently call the dealWithLabelSwitching function.

Author(s)

Panagiotis Papastamoulis

References

Papastamoulis, P. (2017). Overfitting Bayesian mixtures of factor analyzers with an unknown number of components. arXiv:1701.04605 [stat.ME]

See Also

dealWithLabelSwitching

Examples

```
## Not run:
# simulate a synthetic dataset along the lines of the paper:
n = 1000              # sample size
p = 40                # number of variables
q = 4                 # number of factors
K = 10                # number of clusters
sINV_diag = 1/((1:p)) # diagonal of inverse variance of errors
set.seed(10)
syntheticDataset <- simData(K.true = K, n = n, q = q, p = p, sINV_values = sINV_diag )

# define parameters
Kmax <- 20    # number of overfitted mixture components
nChains <- 8  # number of parallel chains
dN <- 1
# Dirichlet prior of mixture weights per chain.
#   The target chain corresponds to the first entry.
dirPriorAlphas <- c(1, 1 + dN * (2:nChains - 1))/Kmax
outputFolder <- "fabMixExample"
# Run algorithm
fabMix( dirPriorAlphas = dirPriorAlphas,
        rawData = syntheticDataset$data,
        outDir = outputFolder, Kmax = Kmax, mCycles = 1200,
        burnCycles = 200, q = q)

# Compute information criteria:
getStuffForDIC(x_data = syntheticDataset$data, outputFolder = outputFolder, q = q)

# Deal with label switching:
dealWithLabelSwitching(x_data = syntheticDataset$data,
        outputFolder = outputFolder, q = q,
        compute_regularized_expression = TRUE, Km = Kmax)

## End(Not run)
```

---

fabMix_UxC *Main function of the package for UUC, UCC models*

Description

This function runs parallel chains under a prior tempering scheme of the Dirichlet prior distribution of mixture weights.

Usage

```
fabMix_UxC(sameSigma, dirPriorAlphas, rawData, outDir, Kmax, mCycles,
burnCycles, g, h, alpha_sigma, beta_sigma, q, normalize,
thinning, zStart, nIterPerCycle, gibbs_z,
warm_up_overfitting, warm_up, overfittingInitialization,
progressGraphs, gwar)
```

Arguments

`sameSigma`	Logical value denoting the parameterization of the error variance per component. If `TRUE`, the parameterization CCU is fitted. Otherwise, the parameterization CUU is fitted.
`dirPriorAlphas`	The prior Dirichlet parameters for each chain.
`rawData`	The observed data as an $n \times p$ matrix. Clustering is performed on the rows of the matrix.
`outDir`	Name of the output folder.
`Kmax`	Number of components in the overfitted mixture. Default: 20.
`mCycles`	Number of MCMC cycles. Each cycle consists of `nIterPerCycle` MCMC iterations. At the end of each cycle a swap of the state of two randomly chosen adjacent chains is attempted.
`burnCycles`	Number of cycles that will be discarded as burn-in period.
`g`	Prior parameter g. Default value: $g = 2$.
`h`	Prior parameter h. Default value: $h = 1$.
`alpha_sigma`	Prior parameter α. Default value: $\alpha = 2$.
`beta_sigma`	Prior parameter β. Default value: $\beta = 1$.
`q`	Number of factors q, where $1 \leq q \leq L$. An error is thrown if the Ledermann bound (L) is exceeded.
`normalize`	Should the observed data be normalized? Default value: TRUE.
`thinning`	Optional integer denoting the thinning of the keeped MCMC cycles.
`zStart`	Optional starting value for the allocation vector.
`nIterPerCycle`	Number of iteration per MCMC cycle. Default value: 10.
`gibbs_z`	Select the gibbs sampling scheme for updating latent allocations of mixture model. Default value: 1.
`warm_up_overfitting`	Number of iterations for the overfitting initialization scheme. Default value: 100.
`warm_up`	Number of iterations that will be used to initialize the models before starting proposing switchings. Default value: 500.

`overfittingInitialization`
: Logical value indicating whether the chains are initialized via the overfitting initialization scheme. Default: TRUE.

`progressGraphs`
: Logical value indicating whether to plot successive states of the chains while the sampler runs. Default: FALSE.

`gwar`
: Initialization parameter. Default: 0.05.

Value

List of files written to `outDir`

Note

It is recommended to always use: `normalize = TRUE` (default). Tuning of `dirPriorAlphas` may be necessary to achieve reasonable acceptance rates of chain swaps. Also note that the output is not identifiable due to label switching and the user has to subsequently call the `dealWithLabelSwitching` function.

Author(s)

Panagiotis Papastamoulis

References

Papastamoulis, P. (2017). Overfitting Bayesian mixtures of factor analyzers with an unknown number of components. arXiv:1701.04605 [stat.ME]

See Also

`dealWithLabelSwitching`

Examples

```
## Not run:
# simulate a synthetic dataset along the lines of the paper:
n = 1000              # sample size
p = 40                # number of variables
q = 4                 # number of factors
K = 10                # number of clusters
sINV_diag = 1/((1:p)) # diagonal of inverse variance of errors
set.seed(10)
syntheticDataset <- simData(K.true = K, n = n, q = q, p = p, sINV_values = sINV_diag )

# define parameters
Kmax <- 20    # number of overfitted mixture components
nChains <- 8  # number of parallel chains
dN <- 1
# Dirichlet prior of mixture weights per chain.
#   The target chain corresponds to the first entry.
dirPriorAlphas <- c(1, 1 + dN * (2:nChains - 1))/Kmax
outputFolder <- "fabMixExample"
```

```
# Run algorithm
fabMix_UxC( dirPriorAlphas = dirPriorAlphas,
        rawData = syntheticDataset$data,
        outDir = outputFolder, Kmax = Kmax, mCycles = 1200,
        burnCycles = 200, q = q)

# Compute information criteria:
#getStuffForDIC(x_data = syntheticDataset$data, outputFolder = outputFolder, q = q)

# Deal with label switching:
#dealWithLabelSwitching(x_data = syntheticDataset$data,
#        outputFolder = outputFolder, q = q,
#        compute_regularized_expression = TRUE, Km = Kmax)

## End(Not run)
```

`fabMix_UxU` *Fit UxU mixtures*

Description

This function runs parallel chains under a prior tempering scheme of the Dirichlet prior distribution of mixture weights.

Usage

```
fabMix_UxU(sameSigma, dirPriorAlphas, rawData, outDir, Kmax, mCycles,
burnCycles, g, h, alpha_sigma, beta_sigma, q, normalize,
thinning, zStart, nIterPerCycle, gibbs_z,
warm_up_overfitting, warm_up, overfittingInitialization,
progressGraphs, gwar)
```

Arguments

`sameSigma`	Logical value denoting the parameterization of the error variance per component. If `TRUE`, the parameterization $\Sigma_1 = \ldots = \Sigma_K$ is fitted.
`dirPriorAlphas`	The prior Dirichlet parameters for each chain.
`rawData`	The observed data as an $n \times p$ matrix. Clustering is performed on the rows of the matrix.
`outDir`	Name of the output folder.
`Kmax`	Number of components in the overfitted mixture. Default: 20.
`mCycles`	Number of MCMC cycles. Each cycle consists of `nIterPerCycle` MCMC iterations. At the end of each cycle a swap of the state of two randomly chosen adjacent chains is attempted.
`burnCycles`	Number of cycles that will be discarded as burn-in period.

g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
q	Number of factors q, where $1 \leq q \leq L$. An error is thrown if the Ledermann bound (L) is exceeded.
normalize	Should the observed data be normalized? Default value: TRUE.
thinning	Optional integer denoting the thinning of the keeped MCMC cycles.
zStart	Optional starting value for the allocation vector.
nIterPerCycle	Number of iteration per MCMC cycle. Default value: 10.
gibbs_z	Select the gibbs sampling scheme for updating latent allocations of mixture model. Default value: 1.
warm_up_overfitting	Number of iterations for the overfitting initialization scheme. Default value: 100.
warm_up	Number of iterations that will be used to initialize the models before starting proposing switchings. Default value: 500.
overfittingInitialization	Logical value indicating whether the chains are initialized via the overfitting initialization scheme. Default: TRUE.
progressGraphs	Logical value indicating whether to plot successive states of the chains while the sampler runs. Default: FALSE.
gwar	Initialization parameter. Default: 0.05.

Value

List of files written to `outDir`

Note

It is recommended to always use: `normalize = TRUE` (default). Tuning of `dirPriorAlphas` may be necessary to achieve reasonable acceptance rates of chain swaps. Also note that the output is not identifiable due to label switching and the user has to subsequently call the `dealWithLabelSwitching` function.

Author(s)

Panagiotis Papastamoulis

References

Papastamoulis, P. (2017). Overfitting Bayesian mixtures of factor analyzers with an unknown number of components. arXiv:1701.04605 [stat.ME]

See Also

dealWithLabelSwitching

Examples

```
## Not run:
# simulate a synthetic dataset along the lines of the paper:
n = 1000              # sample size
p = 40                # number of variables
q = 4                 # number of factors
K = 10                # number of clusters
sINV_diag = 1/((1:p)) # diagonal of inverse variance of errors
set.seed(10)
syntheticDataset <- simData(K.true = K, n = n, q = q, p = p, sINV_values = sINV_diag )

# define parameters
Kmax <- 20    # number of overfitted mixture components
nChains <- 8  # number of parallel chains
dN <- 1
# Dirichlet prior of mixture weights per chain.
#   The target chain corresponds to the first entry.
dirPriorAlphas <- c(1, 1 + dN * (2:nChains - 1))/Kmax
outputFolder <- "fabMixExample"
# Run algorithm
fabMix_UxU( dirPriorAlphas = dirPriorAlphas,
        rawData = syntheticDataset$data,
        outDir = outputFolder, Kmax = Kmax, mCycles = 1200,
        burnCycles = 200, q = q)

# Compute information criteria:
getStuffForDIC(x_data = syntheticDataset$data, outputFolder = outputFolder, q = q)

# Deal with label switching:
dealWithLabelSwitching(x_data = syntheticDataset$data,
        outputFolder = outputFolder, q = q,
        compute_regularized_expression = TRUE, Km = Kmax)

## End(Not run)
```

---

getStuffForDIC *Compute information criteria*

Description

This function computes four information criteria for a given run of the fabMix algorithm, namely: AIC, BIC, DIC and DIC_2. Given various runs with different number of factors, the selected model corresponds to the one with the smalled value of the selected criterion.

Usage

```
getStuffForDIC(sameSigma, sameLambda, isotropic, x_data, outputFolder, q, burn,
Km, normalize, discardLower)
```

Arguments

sameSigma	Logical value indicating whether the parameterization with the same variance of errors per component is used. Default: TRUE.
sameLambda	Logical value indicating whether the parameterization with same loadings per component is used. Default: FALSE.
isotropic	Logical value indicating whether the parameterization with isotropic error variance per component is used. Default: FALSE.
x_data	Observed data.
outputFolder	Name of the folder where the fabMix function has saved its output.
q	Number of factors. Note that this should coincide with the number of factors in the fabMix run.
burn	Discard observations as burn-in period (optional).
Km	Number of components in the overfitted mixture model. Note that this should coincide with the same entry in the fabMix run.
normalize	Should the observed data be normalized? Note that this should coincide with the same entry in the fabMix run. Default value: TRUE.
discardLower	Discard draws with log-likelihood values lower than the specific quantile. This applied only for the DIC computation.

Details

If necessary, more details than the description above

Value

The information criteria are saved to the informationCriteria_map_model.txt file in the code-outputFolder.

Note

It is well known that DIC tends to overfit, so it advised to compare models with different number of factors using AIC or BIC.

Author(s)

Panagiotis Papastamoulis

log_dirichlet_pdf *Log-density function of the Dirichlet distribution*

Description

Log-density function of the Dirichlet distribution

Usage

```
log_dirichlet_pdf(alpha, weights)
```

Arguments

alpha	Parameter vector
weights	Vector of weights

Value

Log-density of the $D(alpha_1, \ldots, \alpha_k)$ evaluated at $w_1, \ldots, w_k$.

Author(s)

Panagiotis Papastamoulis

myDirichlet *Simulate from the Dirichlet distribution*

Description

Generate a random draw from the Dirichlet distribution $D(\alpha_1, \ldots, \alpha_k)$.

Usage

```
myDirichlet(alpha)
```

Arguments

alpha	Parameter vector

Value

Simulated vector

Author(s)

Panagiotis Papastamoulis

observed.log.likelihood0 *Log-likelihood of the mixture model*

Description

Log-likelihood of the mixture model evaluated only at the alive components.

Usage

```
observed.log.likelihood0(x_data, w, mu, Lambda, Sigma, z)
```

Arguments

x_data	The observed data
w	Vector of mixture weights
mu	Vector of marginal means
Lambda	Factor loadings
Sigma	Common covariance matrix of the errors per cluster
z	Allocation vector

Value

Log-likelihood value

Author(s)

Panagiotis Papastamoulis

observed.log.likelihood0_q0_sameSigma *Log-likelihood of the mixture model for $q = 0$ and same variance of errors*

Description

Log-likelihood of the mixture model evaluated only at the alive components.

Usage

```
observed.log.likelihood0_q0_sameSigma(x_data, w, mu, Sigma, z)
```

Arguments

x_data	The observed data
w	Vector of mixture weights
mu	Vector of marginal means
Sigma	Covariance matrix of the errors per cluster
z	Allocation vector

Value

Log-likelihood value

Author(s)

Panagiotis Papastamoulis

---

observed.log.likelihood0_Sj
Log-likelihood of the mixture model

Description

Log-likelihood of the mixture model evaluated only at the alive components.

Usage

```
observed.log.likelihood0_Sj(x_data, w, mu, Lambda, Sigma, z)
```

Arguments

x_data	The observed data
w	Vector of mixture weights
mu	Vector of marginal means
Lambda	Factor loadings
Sigma	Covariance matrix of the errors per cluster
z	Allocation vector

Value

Log-likelihood value

Author(s)

Panagiotis Papastamoulis

observed.log.likelihood0_Sj_q0 *Log-likelihood of the mixture model for $q = 0$*

Description

Log-likelihood of the mixture model evaluated only at the alive components.

Usage

```
observed.log.likelihood0_Sj_q0(x_data, w, mu, Sigma, z)
```

Arguments

x_data	The observed data
w	Vector of mixture weights
mu	Vector of marginal means
Sigma	Covariance matrix of the errors per cluster
z	Allocation vector

Value

Log-likelihood value

Author(s)

Panagiotis Papastamoulis

overfittingMFA *Basic MCMC sampler*

Description

Gibbs sampling for fitting a mixture model of factor analyzers.

Usage

```
overfittingMFA(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_CCC *Basic MCMC sampler for CCC*

Description

Gibbs sampling for fitting a CCC mixture model of factor analyzers.

Usage

```
overfittingMFA_CCC(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_CCU *Basic MCMC sampler for CCU*

Description

Gibbs sampling for fitting a CCU mixture model of factor analyzers.

Usage

```
overfittingMFA_CCU(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_CUC *Basic MCMC sampler for CUC*

Description

Gibbs sampling for fitting a CUC mixture model of factor analyzers.

Usage

```
overfittingMFA_CUC(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_CUU *Basic MCMC sampler for CUU*

Description

Gibbs sampling for fitting a CUU mixture model of factor analyzers.

Usage

```
overfittingMFA_CUU(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_missing_values

Basic MCMC sampler for the case of missing data

Description

Gibbs sampling for fitting a mixture model of factor analyzers.

Usage

```
overfittingMFA_missing_values(missing_entries, x_data, originalX, outputDirectory, Kmax,
m, thinning, burn, g, h, alpha_prior, alpha_sigma,
beta_sigma, start_values, q, zStart, gibbs_z)
```

Arguments

missing_entries	list which contains the row number (1st entry) and column indexes (subsequent entries) for every row containing missing values.
x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_Sj *Basic MCMC sampler using different error variance per component*

Description

Gibbs sampling for fitting a mixture model of factor analyzers.

Usage

```
overfittingMFA_Sj(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_Sj_missing_values

Basic MCMC sampler for the case of missing data and different error variance

---

Description

Gibbs sampling for fitting a mixture model of factor analyzers.

Usage

```
overfittingMFA_Sj_missing_values(missing_entries, x_data, originalX,
outputDirectory, Kmax,
m, thinning, burn, g, h, alpha_prior, alpha_sigma,
beta_sigma, start_values, q, zStart, gibbs_z)
```

Arguments

missing_entries	list which contains the row number (1st entry) and column indexes (subsequent entries) for every row containing missing values.
x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_UCC *Basic MCMC sampler for CCC*

Description

Gibbs sampling for fitting a UCC mixture model of factor analyzers.

Usage

```
overfittingMFA_UCC(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfittingMFA_UUC *Basic MCMC sampler for UUC*

Description

Gibbs sampling for fitting a UUC mixture model of factor analyzers.

Usage

```
overfittingMFA_UUC(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma,
start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfitting_q0 *MCMC sampler for* $q = 0$

Description

Gibbs sampling for fitting a mixture model with diagonal covariance structure.

Usage

```
overfitting_q0(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma, start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

overfitting_q0_sameSigma

MCMC sampler for $q = 0$ and same error variance parameterization

---

Description

Gibbs sampling for fitting a mixture model with diagonal covariance structure.

Usage

```
overfitting_q0_sameSigma(x_data, originalX, outputDirectory, Kmax, m, thinning, burn,
g, h, alpha_prior, alpha_sigma, beta_sigma, start_values, q, zStart, gibbs_z)
```

Arguments

x_data	normalized data
originalX	observed raw data (only for plotting purpose)
outputDirectory	Name of the output folder
Kmax	Number of mixture components
m	Number of iterations
thinning	Thinning of chain
burn	Burn-in period
g	Prior parameter g. Default value: $g = 2$.
h	Prior parameter h. Default value: $h = 1$.
alpha_prior	Parameters of the Dirichlet prior distribution of mixture weights.
alpha_sigma	Prior parameter α. Default value: $\alpha = 2$.
beta_sigma	Prior parameter β. Default value: $\beta = 1$.
start_values	Optional (not used)
q	Number of factors.
zStart	Optional (not used)
gibbs_z	Optional

Value

List of files

Author(s)

Panagiotis Papastamoulis

---

plot.fabMix.object *Plot function*

Description

This function plots fabMix function.

Usage

```
## S3 method for class 'fabMix.object'
plot(x, what, variableSubset, ...)
```

Arguments

x	An object of class fabMix.object, which is returned by the fabMix function.
what	One of the "BIC", "classification_matplot", "classification_pairs", "correlation", "regularized_expression". The plot will display the BIC values per model and number of factors (along with the most probable number of clusters as text), a matplot per cluster for the selected model, scatterplots pairs, the estimated correlation matrix per cluster, and the estimated regularized expression of each variable to the factor space for the selected model, respectively.
variableSubset	An optional subset of the variables. By default, all variables are selected.
...	ignored.

Details

The function plots the BIC values obtained for each of the fitted models, visualizes the clusters.

Author(s)

Panagiotis Papastamoulis

---

print.fabMix.object *Print function*

Description

This function prints a summary of objects returned by the fabMix function.

Usage

```
## S3 method for class 'fabMix.object'
print(x, printSubset, ...)
```

Arguments

x	An object of class fabMix.object, which is returned by the fabMix function.
printSubset	Logical indicating whether to print the header or the whole matrix of estimates. Default value: TRUE.
...	ignored.

Details

The function prints the estimated distribution of the number of clusters, the estimated number of observations assigned to each cluster after post-processing the output with three label switching algorithms, as well as the header of the posterior mean estimates of θ_{kj} (probability of success for cluster k and feature j) (conditionally on the most probable number of clusters).

Author(s)

Panagiotis Papastamoulis

readLambdaValues *Read Lambda values.*

Description

Function to read Lambda values from file.

Usage

```
readLambdaValues(myFile,K,p,q)
```

Arguments

myFile	File containing Lambda values
K	Number of components
p	Number of variables
q	Number of factors

Value

$K \times p \times q$ array of factor loadings.

Author(s)

Panagiotis Papastamoulis

simData *Synthetic data generator*

Description

Simulate data from a multivariate normal mixture using a mixture of factor analyzers mechanism.

Usage

```
simData(sameSigma, sameLambda, p, q, K.true, n, loading_means, loading_sd, sINV_values)
```

Arguments

sameSigma	Logical.
sameLambda	Logical.
p	The dimension of the multivariate normal distribution ($p > 1$).
q	Number of factors. It should be strictly smaller than p.
K.true	The number of mixture components (clusters).
n	Sample size.
loading_means	A vector which contains the means of blocks of factor loadings. Default: `loading_means = c(-30,-20,-10,10, 20, 30)`.
loading_sd	A vector which contains the standard deviations of blocks of factor loadings. Default: `loading_sd <- rep(2, length(loading_means))`.
sINV_values	A vector which contains the values of the diagonal of the (common) inverse covariance matrix, if `sigmaTrue = TRUE`. An $K \times p$ matrix which contains the values of the diagonal of the inverse covariance matrix per component, if `sigmaTrue = FALSE`. Default: `sINV_values = rgamma(p, shape = 1, rate = 1)`.

Value

A list with the following entries:

data	$n \times p$ array containing the simulated data.
class	n-dimensional vector containing the class of each observation.
factorLoadings	$K.true \times p \times q$-array containing the factor loadings Λ_{krj} per cluster k, feature r and factor j, where $k = 1, \ldots, K$; $r = 1, \ldots, p$; $j = 1, \ldots, q$.
means	$K.true \times p$ matrix containing the marginal means μ_{kr}, $k = 1, \ldots, K$; $r = 1, \ldots, p$.
variance	$p \times p$ diagonal matrix containing the variance of errors σ_{rr}, $r = 1, \ldots, p$. Note that the same variance of errors is assumed for each cluster.
factors	$n \times q$ matrix containing the simulated factor values.
weights	$K.true$-dimensional vector containing the weight of each cluster.

Note

The marginal variance for cluster k is equal to $\Lambda_k \Lambda_k^T + \Sigma$.

Author(s)

Panagiotis Papastamoulis

update_all_y *Gibbs sampling for y*

Description

Gibbs sampling for y

Usage

```
update_all_y(x_data, mu, SigmaINV, Lambda, z)
```

Arguments

x_data	Data
mu	Marginal means
SigmaINV	Precision matrix
Lambda	Factor loadings
z	Allocation vector

Value

A matrix with generated factors

Author(s)

Panagiotis Papastamoulis

update_all_y_Sj *Gibbs sampling for y*

Description

Gibbs sampling for y

Usage

```
update_all_y_Sj(x_data, mu, SigmaINV, Lambda, z)
```

Arguments

x_data	Data
mu	Marginal means
SigmaINV	Precision matrix per component
Lambda	Factor loadings
z	Allocation vector

Value

A matrix with generated factors

Author(s)

Panagiotis Papastamoulis

`update_OmegaINV` *Gibbs sampling for* $\Omega\hat{}-1$

Description

Gibbs sampling for Ω^{-1}

Usage

```
update_OmegaINV(Lambda, K, g, h)
```

Arguments

`Lambda`	Factor loadings
`K`	Number of components
`g`	Prior parameter
`h`	Prior parameter

Value

Ω^{-1}

Author(s)

Panagiotis Papastamoulis

update_OmegaINV_Cxx *Gibbs sampling for* $\Omega\hat{}-1$ *for Cxx model*

Description

Gibbs sampling for Ω^{-1} for Cxx model

Usage

```
update_OmegaINV_Cxx(Lambda, K, g, h)
```

Arguments

Lambda	Factor loadings
K	Number of components
g	Prior parameter
h	Prior parameter

Value

Ω^{-1}

Author(s)

Panagiotis Papastamoulis

update_SigmaINV_faster
Gibbs sampling for $\Sigma\hat{}-1$

Description

Gibbs sampling for Σ^{-1}

Usage

```
update_SigmaINV_faster(x_data, z, y, Lambda, mu, K, alpha_sigma, beta_sigma)
```

Arguments

x_data	Data
z	Allocation vector
y	Factors
Lambda	Factor loadings
mu	Marginal means
K	Number of components
alpha_sigma	Prior parameter
beta_sigma	Prior parameter

Value

Σ^{-1}

Author(s)

Panagiotis Papastamoulis

---

update_SigmaINV_faster_q0

Gibbs sampling for $\Sigma\hat{}-1$ per component for $q = 0$

Description

Gibbs sampling for Σ^{-1} per component

Usage

```
update_SigmaINV_faster_q0( z, mu, K, alpha_sigma, beta_sigma, x_data)
```

Arguments

z	Allocation vector
mu	Marginal means
K	Number of components
alpha_sigma	Prior parameter
beta_sigma	Prior parameter
x_data	Data

Value

Σ^{-1}

Author(s)

Panagiotis Papastamoulis

update_SigmaINV_faster_q0_sameSigma
Gibbs sampling for $\Sigma\hat{}-1$ *per component for* $q = 0$

Description

Gibbs sampling for Σ^{-1} per component

Usage

```
update_SigmaINV_faster_q0_sameSigma( z, mu, K, alpha_sigma, beta_sigma, x_data)
```

Arguments

z	Allocation vector
mu	Marginal means
K	Number of components
alpha_sigma	Prior parameter
beta_sigma	Prior parameter
x_data	Data

Value

Σ^{-1}

Author(s)

Panagiotis Papastamoulis

update_SigmaINV_faster_Sj
Gibbs sampling for $\Sigma\hat{}-1$ *per component*

Description

Gibbs sampling for Σ^{-1} per component

Usage

```
update_SigmaINV_faster_Sj(x_data, z, y, Lambda, mu, K, alpha_sigma, beta_sigma)
```

Arguments

x_data	Data
z	Allocation vector
y	Factors
Lambda	Factor loadings
mu	Marginal means
K	Number of components
alpha_sigma	Prior parameter
beta_sigma	Prior parameter

Value

Σ^{-1}

Author(s)

Panagiotis Papastamoulis

---

update_SigmaINV_xCC *Gibbs sampling for $\Sigma\hat{}-1$ for xCC models*

Description

Gibbs sampling for Σ^{-1} for xCC models

Usage

```
update_SigmaINV_xCC(x_data, z, y, Lambda, mu, K, alpha_sigma, beta_sigma)
```

Arguments

x_data	Data
z	Allocation vector
y	Factors
Lambda	Factor loadings
mu	Marginal means
K	Number of components
alpha_sigma	Prior parameter
beta_sigma	Prior parameter

Value

Σ^{-1}

Author(s)

Panagiotis Papastamoulis

`update_SigmaINV_xUC` *Gibbs sampling for* $\Sigma\hat{}-1$ *per component for xUC models*

Description

Gibbs sampling for Σ^{-1} per component for xUC models

Usage

```
update_SigmaINV_xUC(x_data, z, y, Lambda, mu, K, alpha_sigma, beta_sigma)
```

Arguments

`x_data`	Data
`z`	Allocation vector
`y`	Factors
`Lambda`	Factor loadings
`mu`	Marginal means
`K`	Number of components
`alpha_sigma`	Prior parameter
`beta_sigma`	Prior parameter

Value

Σ^{-1}

Author(s)

Panagiotis Papastamoulis

update_z2 *Collapsed Gibbs for z using matrix inversion lemma*

Description

Collapsed Gibbs for z using matrix inversion lemma

Usage

```
update_z2(w, mu, Lambda, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
SigmaINV	Precision matrix
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

update_z2_Sj *Collapsed Gibbs for z using matrix inversion lemma*

Description

Collapsed Gibbs for z using matrix inversion lemma

Usage

```
update_z2_Sj(w, mu, Lambda, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
SigmaINV	Precision matrix per component
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

---

update_z4 *Collapsed Gibbs for z*

Description

Collapsed Gibbs for z

Usage

```
update_z4(w, mu, Lambda, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
SigmaINV	Precision matrix
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

update_z4_Sj *Collapsed Gibbs for z*

Description

Collapsed Gibbs for z

Usage

```
update_z4_Sj(w, mu, Lambda, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
SigmaINV	Precision matrix per component
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

update_z_b *Gibbs sampling for z*

Description

Gibbs sampling for z

Usage

```
update_z_b(w, mu, Lambda, y, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
y	Matrix of factors
SigmaINV	Precision matrix
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

---

update_z_b_Sj *Gibbs sampling for z*

Description

Gibbs sampling for z

Usage

```
update_z_b_Sj(w, mu, Lambda, y, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
Lambda	Factor loadings
y	Matrix of factors
SigmaINV	Precision matrix per component
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

update_z_q0 *Gibbs sampling for z for $q = 0$*

Description

Gibbs sampling for z

Usage

```
update_z_q0(w, mu, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
SigmaINV	Precision matrix per component
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

update_z_q0_sameSigma *Gibbs sampling for z for $q = 0$*

Description

Gibbs sampling for z

Usage

```
update_z_q0_sameSigma(w, mu, SigmaINV, K, x_data)
```

Arguments

w	Mixture weights
mu	Marginal means
SigmaINV	Precision matrix per component
K	Number of components
x_data	Data

Value

Allocation vector

Author(s)

Panagiotis Papastamoulis

`waveDataset1500` *Wave dataset*

Description

A subset of 1500 randomly sampled observations from the wave dataset (version 1), available from the UCI machine learning repository. It contains 3 classes of waves (variable `class` with values "1", "2" and "3") and 21 attributes. Each class is generated from a combination of 2 of 3 base waves with noise.

Usage

```
waveDataset1500
```

Format

A data frame with 1500 rows and 22 columns. The first column denotes the class of each observation.

Source

https://archive.ics.uci.edu/ml/datasets/Waveform+Database+Generator+(Version+1)

References

Lichman, M. (2013). UCI Machine Learning Repository http://archive.ics.uci.edu/ml. Irvine, CA: University of California, School of Information and Computer Science.

Breiman,L., Friedman,J.H., Olshen,R.A. and Stone,C.J. (1984). Classification and Regression Trees. Wadsworth International Group: Belmont, California.

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