

Package ‘FMMcsVS’

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Type Package

Title Bayesian Finite Mixture Regression Models with Cluster-Specific Variable Selections

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Description Different MCMC algorithms for different Bayesian mixture models in a regression setup, including clustering, variable selection and regression coefficient estimations.

Depends Matrix,
caret,
MCMCprecision,
truncdist,
faux,
mcclust.ext,
philentropy,
aricode,
gsl,
label.switching,
mclust,
clustvarsel,
coda,
flexmix,
BNPmix,
factoextra,
MASS,
R (\geq 2.10)

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LazyData true

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coffee	<i>Coffee Data</i>
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Description

Data on the chemical composition of coffee samples collected from around the world, comprising 43 samples from 29 countries. Each sample is either of the Arabica or Robusta variety. Twelve of the thirteen chemical constituents reported in the study are given. The omitted variable is total chlorogenic acid; it is generally the sum of the chlorogenic, neochlorogenic and isochlorogenic acid values.

Usage

coffee

Format

A data frame with 43 observations and 14 variables. The first two columns contain Variety and Country, respectively, while the remaining 12 columns contain the chemical properties. The Variety is either (1) Arabica or (2) Robusta

Note

The German to English translations of the variable names were carried out by Dr. Sharon M. McNicholas.

Source

Streuli, H. (1973). Der heutige stand der kaffeechemie. In Association Scientifique Internationale du Cafe, 6th International Colloquium on Coffee Chemisrty, Bogata, Columbia, pp. 61–72.

Examples

```
coffee
variety_coffee <- coffee$Variety
```

country	<i>Country Data</i>
---------	---------------------

Description

Country profile data with 4 socio-economic and 3 public health indicators that measure the overall development of a country in 2019. The factors are chosen from the World Development Indicators (WDI) database of the World Bank.

Usage

```
country
```

Format

A data frame with 157 observations and 8 variables:

country_name country name

gdppc The GDP per capita. Calculated as the Total GDP divided by the total population.

health Current health expenditure (percentage of GDP).

import Imports of goods and services (percentage of the GDP).

income Adjusted net national income per capita (current USD).

inflation Inflation, GDP deflator (annual percentage), the measurement of the annual growth rate of the total GDP.

life_exp Life expectancy at birth, total (years). The average number of years a new born child would live if the current mortality patterns are to remain the same.

total_fert Fertility rate, total (births per woman). The number of children that would be born to each woman if the current age-fertility rates remain the same.

Note

Due to the COVID-19 pandemic, the current database does not update health expenditure data after 2019, so we collect the data for all indicators of all countries in year 2019. This will exclude some important countries, eg, the United Kingdom, since the income data for UK is only updated through 2018. After removing countries with missing values in any of the indicators, 157 countries remain in the data.

Source

World Bank Open Data, <https://data.worldbank.org/>

Examples

```
country
name_country <- country$country_name
```

data_gen_func*Generate Simulation Data*

Description

Generate the simulation data for various models. The data consists of single-dim response y and multi-dim covariates X . For subjects in different clusters, $(\alpha, \beta, \zeta, \lambda)$ are different. The function can generate correlated data, by specifying the correlation matrix among X , and balanced/inbalanced data by specifying the cluster probabilities.

For the split model, the data consists of y , fixed covariates W and random covariates Z . For subjects in different clusters, $(\alpha, \beta, \psi, \zeta, \lambda)$ are different.

For covariates following t and Gamma distributions, only independent covariates are generated.

Usage

```
data_gen_func(n = 500, alpha_true = c(0.1, -0.6, 0.5),
             beta_true = rbind(c(0, 0, -0.5, 0, 0.5, 0),
                               c(-0.7, 0, 0.4, 0, 0, 0),
                               c(0.6, 0, 0, 0, -0.4, 0)),
             lambda = c(2, 2, 2), zeta_sep = 1,
             eta = 1, sample_prob = c(1, 1, 1),
             cor_mtx = NULL, rho = rep(0, 3))

data_gen_split(n = 500, alpha_true = c(0.1, -0.6, 0.5),
              beta_true = rbind(c(0, 0, -0.5, 0, 0.5, 0),
                                c(-0.7, 0, 0.4, 0, 0, 0),
                                c(0.6, 0, 0, 0, -0.4, 0)),
              psi_true = rep(c(-1, 0, 1), 3),
              W_mean = 0, lambda = c(2, 2, 2), zeta_sep = 1,
              eta = 1, sample_prob = c(1, 1, 1), rho = 0.5)

data_gen_gamma <- function(n = 500, alpha_true = c(0.1, -0.6, 0.5),
                          beta_true = rbind(c(0, 0, -0.5, 0, 0.5, 0),
                                              c(-0.7, 0, 0.4, 0, 0, 0),
                                              c(0.6, 0, 0, 0, -0.4, 0)),
                          lambda = c(2, 2, 2), zeta_sep = 1,
                          sample_prob = c(1, 1, 1))
data_gen_t <- function(n = 500, alpha_true = c(0.1, -0.6, 0.5),
                     beta_true = rbind(c(0, 0, -0.5, 0, 0.5, 0),
                                         c(-0.7, 0, 0.4, 0, 0, 0),
                                         c(0.6, 0, 0, 0, -0.4, 0)),
                     lambda = c(2, 2, 2), zeta_sep = 1,
                     sample_prob = c(1, 1, 1))
```

Arguments

n sample size

<code>alpha_true</code>	numeric vector, each element represents intercept for each cluster, length should equal the number of clusters
<code>beta_true</code>	numeric matrix, each row represents regression coefficients for each cluster, number of rows should equal the number of clusters
<code>lambda</code>	numeric vector, each element represent precision of response y in each cluster
<code>zeta_sep</code>	numeric, the difference of zeta values in different clusters, the true zeta values are set to be $\text{seq}(0, \text{true_M}-1) * \text{zeta_sep}$
<code>eta</code>	numeric, precision of covariates X, for split model, the precision of Z
<code>sample_prob</code>	numeric vector, ratios of sizes of clusters
<code>cor_mtx</code>	numeric matrix, correlation matrix of X, default value is for D=6, M=3, X1-X2, X3-X4, X5-X6 have blocked-wise correlation with coefficients given by rho (no specified in definition)
<code>rho</code>	numeric vector, correlation coefficients among X in the case described above. For split model, rho is a numeric value, which is the pair-wise correlation among columns of W

Details

For `data_gen_func`:

The regression coefficients alpha, beta, the mean of X, zeta, and the precision of y, lambda are cluster-specific.

If no `cor_mtx` value is specified: 1) if rho is a vector of zero's, generated samples have independent X's; 2) if non-zero values are specified in rho, it's assumed to be the D=6, M=3 case, with a blocked-wise correlations of X1-X2, X3-X4 and X5-X6, and correlation coefficients given by rho, if a different setup (eg, different D or M) is desired, `cor_mtx` must be specified.

The true value of $\text{\$}\backslash\text{zeta}\text{\$}$ for the three groups are 0, `zeta_sep` and $2 * \text{zeta_sep}$ respectively.

For `data_gen_split`:

The regression coefficients alpha, beta for W, psi for Z, the mean of Z, zeta, and the precision of y, lambda, are cluster-specific.

W has a pair-wise correlation of rho.

For `data_gen_gamma`:

Covariates are independently following a $\text{Gamma}(z^2, z)$ distributions, where for the M groups, $z = 2 + m * \text{zeta_sep}$, $m = 0, 1, \dots, M-1$.

For `data_gen_t`:

Covariates are independently following the $t_M / \sqrt{M/(M-2)} + z$, where t_M is the Student's t-distribution with a degree of freedom of M, $z = m * \text{zeta_sep}$, $m = 0, 1, \dots, M-1$.

Value

A list is returned:

<code>y</code>	numeric vector, response with length n
<code>X</code>	numeric matrix, covariates matrix with dimension $M * D$
<code>W</code>	numeric matrix, fixed covariates matrix in split model
<code>Z</code>	numeric matrix, random covariates matrix in split model

index	numeric vector with length n, group membership indicator ranging from 1 to M for each subject
alpha_true	specified true alpha values
beta_true	specified true beta values
psi_true	specified true psi values for split model

Author(s)

Zhen Wang <zwangiowa@gmail.com>

Examples

```
##generate data with independent X
sample_data_1 <- data_gen_func()

##generate inbalanced data with ratio 1:2:10
sample_data_2 <- data_gen_func(sample_prob = c(1,2,10))

##generate data where D=6, M=3, and X has blocked-wise correlations of X1-X2, X3-X4 and X5-X6, and correlation c
sample_data_3 <- data_gen_func(rho = c(0.2,0.5,0.8))

##generate data where correlation matrix among X is given by:
require(Matrix)
cor_mtx = bdiag(matrix(c(1/1, 0.8,0.8, 1/1), 2, 2),
                 matrix(c(1/1, 0.2,0.2, 1/1), 2, 2),
                 matrix(c(1/1, 0.9,0.9, 1/1), 2, 2))
sample_data_4 <- data_gen_func(cor_mtx=cor_mtx)

##generate data for the split model, where W has pair-wise correlation of 0.3
sample_split_data_1 <- data_gen_split(rho=0.3)

##generate data with gamma covariates
sample_data_gamma <- data_gen_gamma()

##generate data with t covariates
sample_data_t <- data_gen_t()
```

flea

Flea Beatles Measurements

Description

This data is from a paper by A. A. Lubischew, "On the Use of Discriminant Functions in Taxonomy", Biometrics, Dec 1962, pp.455-477.

Usage

```
flea
```

Format

A data frame with 74 observations and 7 variables:

tars1 width of the first joint of the first tarsus in microns (the sum of measurements for both tarsi)

tars2 the same for the second joint

head the maximal width of the head between the external edges of the eyes in 0.01 mm

aede1 the maximal width of the aedeagus in the fore-part in microns

aede2 the front angle of the aedeagus (1 unit = 7.5 degrees)

aede3 the aedeagus width from the side in microns

species which species is being examined - concinna, heptapotamica, heikertingeri

Examples

```
head(flea)
```

happy

Happy Score Data

Description

Happy scores and six life quality factors for 149 countries in 2021.

Usage

```
happy
```

Format

A data frame with 149 observations and 8 variables:

country country name

Happy.Score happy score for countries in 2021, it is the national average response to the question of life evaluations.

GDP.per.cap The GDP per capita. Calculated as the Total GDP divided by the total population.

Socl.spprt Social support (or having someone to count on in times of trouble) is the national average of the binary responses (either 0 or 1) to the GWP question “If you were in trouble, do you have relatives or friends you can count on to help you whenever you need them, or not?”.

Life.expt Healthy Life Expectancy (HLE), based on the data extracted from the World Health Organization’s (WHO) Global Health Observatory data repository.

Freedom Freedom to make life choices is the national average of responses to the GWP question “Are you satisfied or dissatisfied with your freedom to choose what you do with your life?”.

Generosity Generosity is the residual of regressing national average of response to the GWP question “Have you donated money to a charity in the past month?” on GDP per capita.

Corruption Corruption Perception. The measure is the national average of the survey responses to two questions in the GWP: “Is corruption widespread throughout the government or not” and “Is corruption widespread within businesses or not?”.

Source

World Happiness Report 2021, <https://worldhappiness.report/ed/2021/>.

Examples

```
happy
name_country <- happy$Country
```

posterior_inf

Posterior Inference for Various Models

Description

Calculate different metrics defined in the paper to evaluate performance of different models in : clustering accuracy, parameter estimation accuracy and variable selection accuracy. Besides the N-IFPP models, functions for inference of RPMS, BNPmix (P-Y mixture models, M9 and M10) and mclust (model-based clustering, M7) are also included.

Usage

```
post_inf(sim_res, scl = 10001:1e5, data)

post_inf_rpms(sim_res, scl = 10001:1e5, data)

post_inf_bnpmix(sim_res, data)

mclust_vs(data)
```

Arguments

<code>sim_res</code>	list, MCMC simulation results, return from the simulation functions defined in the package
<code>scl</code>	numeric vector, index of remaining samples after burn-in
<code>data</code>	list, the data input used to run the simulations

Details

post_inf runs inference for N-IFPP models, post_inf_rpms runs inference results for the RPMS, post_inf_bnpmix runs inference for the two P-Y mixture models (M9, M10). The current versions of the functions can only handle the case of $M_{\text{true}} = 3$ as defined in the default simulation data setups.

mclust_vs does inference for the model-based clustering model (M7) introduced by Fraley and Raftrey (2002, JASA), two different types of VS procedures were implemented to adapt to the regression setup, which is not considered in the original model. The two methods are: 1) implemented the variable selection methods by the clustvarsel package, obtain the VS and clustering results, then a common set of variables are selected for all clusters, fit a linear regression model within each cluster to obtain estimates of alpha and beta; 2) run the mclust package for clustering with both the response and covariates, based on the clustering results with fixed $K=3$, run the variable selection procedure by the clustvarsel package for each cluster so a cluster-specific VS can be achieved, then within each cluster, fit a linear regression model to obtain estimates of alpha and beta;

Value

post_inf, post_inf_rpms and post_inf_bnpmix return:

true_size	numeric vector, the true cluster sizes in the data
measure_km_ari	numeric, a measurement defined to measure how far away the clusters are from each other, calculated by the ARI values of the true clustering membership and the clustering result of the K-means method, a larger value generally means more distant clusters, thus easier clustering problem
auto_corr_k	numeric vector, auto-correlations for posterior of K
geweke_k	numeric, the Z-statistic for the geweke diagnostics for posterior of K
post_k	numeric vector, posterior distribution of K
mse_a1, mse_a2, mse_a3	numeric vectors, the quartiles of the mean squared errors of alpha for the 3 clusters
mse_b1, mse_b2, mse_b3	numeric vectors, the quartiles of the mean squared errors of beta for the 3 clusters
ARI	numeric vector, the (0, 0.05, 0.1, 0.15, 0.2, 0.25, 0.5, 0.75, 1) quantiles of ARI values calculated by all posterior samples of c
c_bin	numeric vector, sizes of clusters obtained by the Binder loss
c_vi	numeric vector, sizes of clusters obtained by the VI loss
ari_point_bin	numeric, ARI values calculated by c_bin and the truth
ari_point_vi	numeric, ARI values calculated by c_vi and the truth
FS	numeric vector, the mean False Sparsity for each cluster
MS	numeric vector, the mean Missed Sparsity for each cluster
ARI_bnpmix	numeric vector, the ARI values calculated from all posterior samples of c by BNPmix

mclust_vs returns:

clust_no_g	numeric vector, the sizes of clusters by mclust when G is not pre-specified
------------	---

<code>ari_no_g</code>	numeric, ARI values for clustering results by mclust when G is not pre-specified
<code>cluster_no_g_vs</code>	numeric, the sizes of clusters by clustvarsel when G is not pre-specified
<code>ari_no_g_vs</code>	numeric, ARI values for clustering results by clustvarsel when G is not pre-specified
<code>clust_g3</code>	numeric vector, the sizes of clusters by mclust when G is fixed at 3
<code>ari_g3</code>	numeric, ARI values for clustering results by mclust when G is fixed at 3
<code>clust_g3_vs</code>	numeric, ARI values for clustering results by clustvarsel when G is fixed at 3
<code>ari_g3_vs</code>	numeric, ARI values for clustering results by clustvarsel when G is fixed at 3
<code>se_a1.1, se_a2.1, se_a3.1</code>	squared-errors of alpha values for M7, method 1)
<code>se_b1.1, se_b2.1, se_b3.1</code>	squared-errors of alpha values for M7, method 1)
<code>FS1.1, FS2.1, FS3.1</code>	FS values for each cluster for M7, method 1)
<code>MS1.1, MS2.1, MS3.1</code>	MS values for each cluster for M7, method 1)
<code>se_a1.2, se_a2.2, se_a3.2</code>	squared-errors of alpha values for M7, method 2)
<code>se_b1.2, se_b2.2, se_b3.2</code>	squared-errors of alpha values for M7, method 2)
<code>FS1.2, FS2.2, FS3.2</code>	FS values for each cluster for M7, method 2)
<code>MS1.2, MS2.2, MS3.2</code>	MS values for each cluster for M7, method 2)

Note

The MSE values, FS, MS values for VS accuracy are calculated with samples of $K = M = 3$. ARI values are calculated with all posterior samples.

Author(s)

Zhen Wang <zwangiowa@gmail.com>

References

- Chris Fraley and Adrian E Raftery. Model-based clustering, discriminant analysis, and density estimation. *Journal of the American Statistical Association*, 97(458):611–631, 2002.
- Riccardo Corradin, Antonio Canale, and Bernardo Nipoti. Bnpx: An r package for bayesian non-parametric modeling via pitman-yor mixtures. *Journal of Statistical Software*, 100(15):1–33, 2021.

Examples

```
##generate simulation data
simulation_data <- data_gen_func()

##FBMM with VS, hyper-prior for beta_bel
simulation_1 <- simulation_func(simulation_data$X, simulation_data$y, prior="Bessel")

##run posterior inference, burn-in the first 20k samples
post_fbmm_vs_1 <- post_inf(simulation_1, 20001:1e5, simulation_data)
```

ppd_sim	<i>Calculate the Posterior Predictive Density (ppd) for New Data Subjects</i>
---------	---

Description

A function to calculate the posterior predictive density for a new subject with response y^* and covariates X^* .

Usage

```
ppd_sim(sim.res, x.new, y.new, scl=10001:1e5, thin=20)
```

Arguments

sim.res	list, MCMC simulation results returned by simulation_func, simulation_split or simulation_func_rpms
x.new	numeric vector of length D, covariates for the new subject
y.new	numeric, the new y value to estimate the ppd at
scl	numeric vector, the index of remaining samples after burn-in
thin	numeric, the thinning factor to speed up calculations, posterior samples for every thin-th iterations are used to calculate ppd

Details

See the Appendix of the paper for formulas of the ppd.

Value

The ppd of x.new at y.new is returned.

Author(s)

Zhen Wang <zwangiowa@gmail.com>

Examples

```
##generate simulation data
simulation_data <- data_gen_func()

##FBMM without VS, hyper-prior for beta_bel
simulation_1 <- simulation_func(simulation_data$X, simulation_data$y, prior="Bessel", SS=F)

x.new <- rnorm(6)
y.new <- 0.5
##calculate ppd
ppd.new.sample <- ppd_sim(simulation_1, x.new, y.new, scl=10001:1e5, thin=20)
```

r_sq_post

Calculate R-square Values for all MCMC Samples

Description

A function to calculate the R-square values for all posterior samples returned by MCMC simulations.

Usage

```
r_sq_post(sim_res, data, scl=10001:1e5)
```

Arguments

sim_res	list, MCMC simulation results returned by simulation_func, simulation_split or simulation_func_rpms
data	dataframe, consists of response y and covariates X
scl	numeric vector, the index of remaining samples after burn-in

Details

Given the posterior samples of K, and c, for each iteration, within each cluster, fit a linear regression, obtain the R^2 value, then for each iteration, we obtain k_post R^2 values for k_post different clusters. Do such calculations for all posterior samples.

Value

The R-square values for all clusters in all posterior samples are returned as a numeric vector.

Author(s)

Zhen Wang <zwangiowa@gmail.com>

Examples

```
##generate simulation data
simulation_data <- data_gen_func()

##FBMM without VS, hyper-prior for beta_bel
simulation_1 <- simulation_func(simulation_data$X, simulation_data$y, prior="Bessel", SS=F)

##calculate the R^2 values
r_sq_samples <- r_sq_post(simulation_1, simulation_data)
```

simulation_func	<i>MCMC Simulations for the Bayesian Finite Mixture Regression Models with Cluster-Specific Variable Selection</i>
-----------------	--

Description

A simulation function running MCMC simulations for various models with different algorithms. Accepted data should be one-dim response y and multi-dim ($D \geq 1$) covariates X expect for the split model, for split model, W and Z should be specified instead of a single covariates matrix X .

Usage

```
simulation_func(X, y, prior = "Dirichlet", SS = TRUE, N = 1e5, gamma_hyperprior = TRUE,
               gamma_fixed = 1, a_gamma = 10, b_gamma = 10, a_unif = 0,
               a_w = 1, b_w = 1, Lambda = 3, a_bessel = 2,
               b_bessel_hyperprior = TRUE, b_bessel_fixed = 1.1, a_b_bessel = 1, b_b_bessel = 10,
               mu = 0, a_tau = 1, b_tau = 1, a_zeta = 0, b_zeta = 1,
               a_lambda = 5, b_lambda = 2, a_alpha = 0, b_alpha = 0.01,
               a_eta = 5, b_eta = 2, L_dynamic = 10, M_init = 6,
               lambda_init = 2, alpha_init = 0)

{
  \item{X}{numeric matrix, covariates matrix}
}
\item{W}{numeric matrix, covariates matrix that is fixed in a split model}
```

```

\item{X}{numeric matrix, covariates matrix that is modeled with Gaussian distributions in a split mo
\item{y}{numeric vector, response}
\item{prior}{char, priors on mixture weights, or different algorithm. "Dirichlet" for FDMM with cond
\item{SS}{logical, if TRUE, spike and slab prior is specified on beta for cluster-specific variable s
\item{N}{numeric, number of iterations}
\item{gamma_hyperprior}{logical, if TRUE, for FDMM with conditional and marginal algorithm, a hyper
\item{gamma_fixed}{numeric, if gamma_hyperprior=FALSE, the fixed value for gamma}
\item{a_gamma, b_gamma}{numeric, if gamma_hyperprior=TRUE, a Gamma(a_gamma, b_gamma) is assigned to
\item{a_unif}{numeric, for FUMM, the unnormalised mixture weights S follows Unif(a_unif, 1),  $0 \leq a_{unif}$ 
\item{a_w, b_w}{numeric, a Beta(a_w, b_w) is assigned to the SS weights w}
\item{Lambda}{numeric,  $M \sim \text{Poisson}_1(\text{Lambda})$ , a shifted Poisson distribution}
\item{a_bessel}{numeric, for FBMM, alpha_bel is fixed at this value}
\item{b_bessel_hyperprior}{logical, if TRUE, a hyperprior is specified to beta_bel}
\item{b_bessel_fixed}{numeric, if b_bessel_hyperprior=FALSE, beta_bel is fixed at this value}
\item{a_b_bessel, b_b_bessel}{numeric, if b_bessel_hyperprior=TRUE, a Gamma(a_b_bessel, b_b_bessel)
\item{mu}{numeric, mean of the slab part of the SS prior on beta}
\item{a_tau, b_tau}{numeric, a Gamma(a_tau, b_tau) hyper-prior for the precision parameter tau of th
\item{a_zeta, b_zeta}{numeric, a Normal(a_tau, b_tau) hyper-prior for the mean parameter of X, zeta}
\item{a_lambda, b_lambda}{numeric, a Gamma(a_lambda, b_lambda) hyper-prior for the precision param
\item{a_alpha, b_alpha}{numeric, a Normal(a_alpha, b_alpha) hyper-prior for the intercept parameter
\item{a_eta, b_eta}{numeric, a Gamma(a_eta, b_eta) hyper-prior for the precision parameter of X, eta
\item{a_psi}{numeric, the prior mean for psi, the regression coefficients corresponding to W}
\item{a_4, b_4}{numeric, a Gamma(a_4, b_4) hyper-prior is assigned to b_psi, the prior precision of p
\item{L_dynamic}{numeric, the number of the auxiliary states, L, in Algorithm 8 of Neal (2000, JCGS)}
\item{M_init}{numeric, initial value of M}
\item{lambda_init}{numeric, initial value of lambda}
\item{alpha_init}{numeric, initial value of alpha}
}
{

```

For simulation_func, a total of four different models are considered: within the N-IFPP category: FDM

For the FDMM, both the conditional and marginal algorithms are implemented, for other models, only th

For the marginal algorithm, Algorithm 8 of Neal (2000, JCGS) is applied.

```

}
{

```

```

\item{M_post}{numeric vector of length N, posterior samples of M}
\item{c_post}{numeric matrix of dim N*n, each row represents posterior samples of c for one iteration
\item{k_post}{numeric vector of length N, posterior samples of K}
\item{U}{numeric vector of length N, posterior samples of U, the auxiliary variable for the condition
\item{gamma_post}{numeric vector of length N, posterior samples of the concentration parameter gamma
\item{Beta_post}{list, each element of the list represents posterior samples of all beta components
\item{alpha_post}{list, each element of the list represents posterior samples of all alpha component
\item{zeta_post}{list, each element of the list represents posterior samples of all zeta components
\item{lambda_post}{numeric vector, posterior samples of lambda, precision of y}
\item{eta_post}{numeric vector, posterior samples of eta, precision of X}

```

```

\item{tau_post}{numeric vector, posterior samples of tau, precision of the slab part in SS prior}
\item{w_post}{numeric matrix of dim N*D, each row represents posterior samples of the SS weights, w,
\item{weight_post}{list, each element represents posterior samples of the mixture weights, omega, of
\item{a_lambda}{pre-specified values}
\item{b_lambda}{pre-specified values}
\item{a_eta}{pre-specified values}
\item{b_eta}{pre-specified values}
\item{X}{covariates input}
\item{y}{response input}
}

```

```
{
```

Generalized Mixtures of Finite Mixtures and Telescoping Sampling. Sylvia Frühwirth-Schnatter, G

Radford M. Neal. Markov chain sampling methods for Dirichlet process mixture models. Journal of Comput

Is infinity that far? A Bayesian nonparametric perspective of finite mixture models. Raffaele Argiento

```
}
```

```
{
```

Zhen Wang zwangiowa@gmail.com

```
}
```

```
{
```

```
##generate simulation data
```

```
simulation_data <- data_gen_func()
```

```
##FBMM without VS, hyper-prior for beta_bel
```

```
simulation_1 <- simulation_func(simulation_data$X, simulation_data$y, prior="Bessel", SS=F)
```

```
##Dynamic FDMM with VS, with fixed gamma value of 2
```

```
simulation_2 <- simulation_func(simulation_data$X, simulation_data$y, prior="Dynamic_FDMM",
                               gamma_hyperprior = FALSE, gamma_fixed = 2)
```

```
##FUMM where  $S \sim \text{Unif}(0.2, 1)$ 
```

```
simulation_3 <- simulation_func(simulation_data$X, simulation_data$y, prior="Uniform", a_unif=0.2)
```

simulation_func_rpms *MCMC Simulations for the RPMS Model*

Description

A simulation function running MCMC simulations for the RPMS model introduced in Barcella et al (SIM, 2016). Accepted data should be one-dim response y and multi-dim ($D_i = 1$) covariates X .

Usage

```
simulation_func_rpms(X, y, SS = TRUE, N = 1e5, a_w = 1, b_w = 1, mu = 0, a_tau = 1, b_tau = 1,
  a_zeta = 0, b_zeta = 1, a_lambda = 0.01, b_lambda = 0.005, a_alpha = 0,
  b_alpha = 0.01, a_eta = 5, b_eta = 2, a_adp = 1, b_adp = 1,
  k_init = 6, lambda_init = 2, alpha_dp_init = 1, alpha_init = 0, m_aux = 10)
```

Arguments

<code>X</code>	numeric matrix, covariates matrix
<code>y</code>	numeric vector, response
<code>SS</code>	logical, if TRUE, spike and slab prior is specified on beta for cluster-specific variable selection; if FALSE, a continuous Gaussian prior is specified to beta, no variable selection is implemented
<code>N</code>	numeric, number of iterations
<code>a_w, b_w</code>	numeric, a Beta(<code>a_w</code> , <code>b_w</code>) is assigned to the SS weights w
<code>mu</code>	numeric, mean of the slab part of the SS prior on beta
<code>a_tau, b_tau</code>	numeric, a Gamma(<code>a_tau</code> , <code>b_tau</code>) hyper-prior for the precision parameter τ of the slab part of the SS prior on beta
<code>a_zeta, b_zeta</code>	numeric, a Normal(<code>a_tau</code> , <code>b_tau</code>) hyper-prior for the mean parameter of X , ζ
<code>a_lambda, b_lambda</code>	numeric, a Gamma(<code>a_lambda</code> , <code>b_lambda</code>) hyper-prior for the precision parameter of y , λ
<code>a_alpha, b_alpha</code>	numeric, a Normal(<code>a_alpha</code> , <code>b_alpha</code>) hyper-prior for the intercept parameter α
<code>a_eta, b_eta</code>	numeric, a Gamma(<code>a_eta</code> , <code>b_eta</code>) hyper-prior for the precision parameter of X , η
<code>a_adp, b_adp</code>	numeric, a Gamma(<code>a_adp</code> , <code>b_adp</code>) hyper-prior for the concentration parameter, α , of DP
<code>k_init</code>	numeric, initial value of K
<code>lambda_init</code>	numeric, initial value of λ
<code>alpha_dp_init</code>	numeric, initial value of α of DP
<code>alpha_init</code>	numeric, initial value of α
<code>m_aux</code>	numeric, the number of the auxiliary states, m , in Algorithm 8 of Neal (2000, JCGS)

Details

A marginal algorithm is implemented, Algorithm 8 of Neal (2000, JCGS) is applied in this algorithm.

Value

c_post	numeric matrix of dim $N \times n$, each row represents posterior samples of c for one iteration
k_post	numeric vector of length N , posterior samples of K
alpha_dp_post	numeric vector of length N , posterior samples of concentration parameter α of DP
Beta_post	list, each element of the list represents posterior samples of all beta components in each iteration, of length $M_post \times D$
alpha_post	list, each element of the list represents posterior samples of all alpha components in each iteration, of length M_post
zeta_post	list, each element of the list represents posterior samples of all zeta components in each iteration, of length $M_post \times D$
lambda_post	numeric vector, posterior samples of λ , precision of y
eta_post	numeric vector, posterior samples of η , precision of X
tau_post	numeric vector, posterior samples of τ , precision of the slab part in SS prior
w_post	numeric matrix of dim $N \times D$, each row represents posterior samples of the SS weights, w , in one iteration
a_lambda	pre-specified values
b_lambda	pre-specified values
a_eta	pre-specified values
b_eta	pre-specified values
X	covariates input
y	response input

Author(s)

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References

- William Barcella, Maria De Iorio, Gianluca Baio, and James Malone-Lee. Variable selection in co- variate dependent random partition models: an application to urinary tract infection. *Statistics in Medicine*, 35(8):1373–1389, 2016.
- Radford M. Neal. Markov chain sampling methods for Dirichlet process mixture models. *Journal of Computational and Graphical Statistics*, 9(2):249–265, 2000.

Examples

```
##generate simulation data
simulation_data <- data_gen_func()

##RPMS with VS
simulation_rpms_1 <- simulation_func_rpms(simulation_data$X, simulation_data$y)
```

```
##RPMS without VS
simulation_rpms_2 <- simulation_func_rpms(simulation_data$X, simulation_data$y, SS=F)
```

simulation_split	<i>MCMC Simulations for the Bayesian Finite Mixture Regression Models with Cluster-Specific Variable Selection</i>
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Description

A simulation function running MCMC simulations for various split models. Accepted data should be one-dim response y and multi-dim ($D \geq 1$) covariates W and Z . Here Z are the covariates being modeled in the model, while W are the fixed covariates.

Usage

```
simulation_split(W, Z, y, prior = "Dirichlet", SS = TRUE, N = 1e+05,
  gamma_hyperprior = TRUE, gamma_fixed = 1, a_gamma = 1,
  b_gamma = 1, a_w = 1, b_w = 1, Lambda = 3, a_bessel = 2,
  b_bessel_hyperprior = TRUE, b_bessel_fixed = 1.1,
  a_b_bessel = 1, b_b_bessel = 10, mu = 0, a_tau = 1,
  b_tau = 1, a_zeta = 0, b_zeta = 1, a_psi = 0, a_4 = 1, b_4 = 1,
  a_lambda = 0.01, b_lambda = 0.005, a_alpha = 0, b_alpha = 0.01,
  a_eta = 5, b_eta = 2, M_init = 6, lambda_init = 2,
  alpha_init = 0)
```

Arguments

W	numeric matrix, covariates matrix that is fixed in a split model
X	numeric matrix, covariates matrix that is modeled with Gaussian distributions in a split model
y	numeric vector, response
prior	char, priors on mixture weights, or different algorithm. "Dirichlet" for FDMM with conditional algorithm, "Bessel" for FBMM with conditional algorithm.
SS	logical, if TRUE, spike and slab prior is specified on beta for cluster-specific variable selection; if FALSE, a continuous Gaussian prior is specified to beta, no variable selection is implemented
N	numeric, number of iterations

<code>gamma_hyperprior</code>	logical, if TRUE, for FDMMM with conditional and marginal algorithm, a hyper-prior is specified to the concentration parameter gamma, otherwise, gamma is fixed
<code>gamma_fixed</code>	numeric, if <code>gamma_hyperprior</code> =FALSE, the fixed value for gamma
<code>a_gamma, b_gamma</code>	numeric, if <code>gamma_hyperprior</code> =TRUE, a $\text{Gamma}(a_gamma, b_gamma)$ is assigned to gamma
<code>a_w, b_w</code>	numeric, a $\text{Beta}(a_w, b_w)$ is assigned to the SS weights w
<code>Lambda</code>	numeric, $M \sim \text{Poisson}_1(\text{Lambda})$, a shifted Poisson distribution
<code>a_bessel</code>	numeric, for FBMM, <code>alpha_bel</code> is fixed at this value
<code>b_bessel_hyperprior</code>	logical, if TRUE, a hyperprior is specified to <code>beta_bel</code>
<code>b_bessel_fixed</code>	numeric, if <code>b_bessel_hyperprior</code> =FALSE, <code>beta_bel</code> is fixed at this value
<code>a_b_bessel, b_b_bessel</code>	numeric, if <code>b_bessel_hyperprior</code> =TRUE, a $\text{Gamma}(a_b_bessel, b_b_bessel)$ hyper-prior is assigned to <code>beta_bel-1</code>
<code>mu</code>	numeric, mean of the slab part of the SS prior on beta
<code>a_tau, b_tau</code>	numeric, a $\text{Gamma}(a_tau, b_tau)$ hyper-prior for the precision parameter tau of the slab part of the SS prior on beta
<code>a_zeta, b_zeta</code>	numeric, a $\text{Normal}(a_tau, b_tau)$ hyper-prior for the mean parameter of X, zeta
<code>a_lambda, b_lambda</code>	numeric, a $\text{Gamma}(a_lambda, b_lambda)$ hyper-prior for the precision parameter of y, lambda
<code>a_alpha, b_alpha</code>	numeric, a $\text{Normal}(a_alpha, b_alpha)$ hyper-prior for the intercept parameter alpha
<code>a_eta, b_eta</code>	numeric, a $\text{Gamma}(a_eta, b_eta)$ hyper-prior for the precision parameter of X, eta
<code>a_psi</code>	numeric, the prior mean for psi, the regression coefficients corresponding to W
<code>a_4, b_4</code>	numeric, a $\text{Gamma}(a_4, b_4)$ hyper-prior is assigned to <code>b_psi</code> , the prior precision of psi
<code>M_init</code>	numeric, initial value of M
<code>lambda_init</code>	numeric, initial value of lambda
<code>alpha_init</code>	numeric, initial value of alpha

Details

Conditional algorithms are implemented for both FDMM and FBMM with variable selection.

Value

<code>M_post</code>	numeric vector of length N, posterior samples of M
<code>c_post</code>	numeric matrix of dim $N \times n$, each row represents posterior samples of c for one iteration

k_post	numeric vector of length N, posterior samples of K
U	numeric vector of length N, posterior samples of U, the auxiliary variable for the conditional algorithm of N-IFPP models
gamma_post	numeric vector of length N, posterior samples of the concentration parameter gamma in both FDMM models
Beta_post	list, each element of the list represents posterior samples of all beta components in each iteration, of length M_post*D
alpha_post	list, each element of the list represents posterior samples of all alpha components in each iteration, of length M_post
zeta_post	list, each element of the list represents posterior samples of all zeta components in each iteration, of length M_post*D
lambda_post	numeric vector, posterior samples of lambda, precision of y
eta_post	numeric vector, posterior samples of eta, precision of X
tau_post	numeric vector, posterior samples of tau, precision of the slab part in SS prior
w_post	numeric matrix of dim N*D, each row represents posterior samples of the SS weights, w, in one iteration
weight_post	list, each element represents posterior samples of the mixture weights, omega, of length M_post
psi_post	numeric matrix, each row represents posterior samples of psi for each iteration
b_psi_post	numeric vector, posterior samples of b_psi
a_lambda	pre-specified values
b_lambda	pre-specified values
a_eta	pre-specified values
b_eta	pre-specified values
y	response input
W	covariates input that is fixed
Z	covariates input that is modeled

Note

For current version of simulation_split, it's required that $\dim(W) \geq 2$ and $\dim(Z) \geq 3$.

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References

Is infinity that far? A Bayesian nonparametric perspective of finite mixture models. Raffaele Argiento and Maria De Iorio, Ann. Statist. 50(5): 2641-2663, 2022.

Examples

```
##generate data for the split model
data_split <- data_gen_split()

###split model with FBMM
simulation_split_fbmm <- simulation_split(data_split$W, data_split$Z, data_split$y, prior = "Bessel")
```