

Package ‘DEBBI’

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Title Differential Evolution-Based Bayesian Inference

Version 0.1.0

Description Bayesian inference algorithms based on the population-based “differential evolution” (DE) algorithm. Users can obtain posterior mode (MAP) estimates via DEMAP, posterior samples via DEMCMC, and variational approximations via DEVI.

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URL <https://github.com/bmgaldo/DEBBI>

BugReports <https://github.com/bmgaldo/DEBBI>

Encoding UTF-8

Imports stats, doParallel, randtoolbox, numDeriv

Depends R (>= 3.5.0)

RoxygenNote 7.1.2

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AlgoParamsDEMAP

AlgoParamsDEMAP

Description

get control parameters for DEMAP function

Usage

```
AlgoParamsDEMAP(
  n_params,
  n_chains = NULL,
  n_iter = 1000,
  init_sd = 0.01,
  init_center = 0,
  n_cores_use = 1,
  step_size = NULL,
  jitter_size = 1e-06,
  crossover_rate = 1,
  parallel_type = "none",
  return_trace = FALSE,
  thin = 1
)
```

Arguments

<code>n_params</code>	number of free parameters estimated
<code>n_chains</code>	number of particle chains, $3 \times n_params$ is the default value
<code>n_iter</code>	number of iterations to run the sampling algorithm, 1000 is default
<code>init_sd</code>	positive scalar or n_params -dimensional numeric vector, determines the standard deviation of the Gaussian initialization distribution
<code>init_center</code>	scalar or n_params -dimensional numeric vector that determines the mean of the Gaussian initialization distribution
<code>n_cores_use</code>	number of cores used when using parallelization.
<code>step_size</code>	positive scalar, jump size in DE crossover step, default is $2.38/\sqrt{2 \times n_params}$.
<code>jitter_size</code>	positive scalar, noise is added during crossover step from $\text{Uniform}(-\text{jitter_size}, \text{jitter_size})$ distribution. $1e-6$ is the default value.
<code>crossover_rate</code>	number on the interval $(0,1]$. Determines the probability a parameter on a chain is updated on a given crossover step, sampled from a Bernoulli distribution.
<code>parallel_type</code>	string specifying parallelization type. 'none', 'FORK', or 'PSOCK' are valid values. 'none' is default value.
<code>return_trace</code>	logical, if true, function returns particle trajectories. This is helpful for diagnosing convergence or debugging model code. Function will return an iteration/thin $\times n_chains \times n_params$ array and the estimated ELBO of each particle in a iteration/thin $\times n_chains$ array.

thin positive integer, only every 'thin'-th iteration will be stored. Default value is 1. Increasing thin will reduce the memory required, while running chains for longer.

Value

list of control parameters for the DEMAP function

AlgoParamsDEMCMC	<i>AlgoParamsDEMCMC</i>
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Description

AlgoParamsDEMCMC

Usage

```
AlgoParamsDEMCMC(
  n_params,
  n_chains = NULL,
  param_names = NULL,
  n_iter = 1000,
  init_sd = 0.01,
  init_center = 0,
  n_cores_use = 1,
  step_size = NULL,
  jitter_size = 1e-06,
  parallel_type = "none",
  burnin = 0,
  thin = 1
)
```

Arguments

n_params	number of free parameters estimated
n_chains	number of MCMC chains, 3*n_params is the default value
param_names	optional vector of parameter names
n_iter	number of iterations to run the sampling algorithm, 1000 is default
init_sd	positive scalar or n_params-dimensional numeric vector, determines the standard deviation of the Gaussian initialization distribution
init_center	scalar or n_params-dimensional numeric vector, determines the mean of the Gaussian initialization distribution
n_cores_use	number of cores used when using parallelization.
step_size	positive scalar, jump size in DE crossover step, default is $2.38/\sqrt{2*n_params}$ which is optimal for multivariate Gaussian target distribution (ter Braak, 2006)

jitter_size	positive scalar, noise is added during crossover step from Uniform(-jitter_size,jitter_size) distribution. 1e-6 is the default value.
parallel_type	string specifying parallelization type. 'none','FORK', or 'PSOCK' are valid values. 'none' is default value.
burnin	number of initial iterations to discard. Default value is 0.
thin	positive integer, only every 'thin'-th iteration will be stored. Default value is 1. Increasing thin will reduce the memory required, while running chains for longer.

Value

list of control parameters for the DEMCMC function

AlgoParamsDEVI	<i>AlgoParamsDEVI</i>
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Description

get control parameters for DEVI function

Usage

```
AlgoParamsDEVI(
  n_params,
  param_names = NULL,
  n_chains = NULL,
  n_iter = 1000,
  init_sd = 0.01,
  init_center = 0,
  n_cores_use = 1,
  step_size = NULL,
  jitter_size = 1e-06,
  parallel_type = "none",
  use_QMC = TRUE,
  purify = NULL,
  quasi_rand_seq = "halton",
  n_samples_ELBO = 10,
  LRVB_correction = TRUE,
  n_samples_LRVB = 25,
  neg_inf = -750,
  thin = 1,
  burnin = 0,
  return_trace = FALSE,
  crossover_rate = 1
)
```

Arguments

<code>n_params</code>	number of free parameters estimated
<code>param_names</code>	optional vector of parameter names
<code>n_chains</code>	number of particle chains used for optimization, $3 \times n_params$ is the default value
<code>n_iter</code>	number of iterations to run the sampling algorithm, 1000 is default
<code>init_sd</code>	positive scalar or n_params -dimensional numeric vector, determines the standard deviation of the Gaussian initialization distribution
<code>init_center</code>	scalar or n_params -dimensional numeric vector, determines the mean of the Gaussian initialization distribution
<code>n_cores_use</code>	number of cores used when using parallelization.
<code>step_size</code>	positive scalar, jump size in DE crossover step, default is $2.38/\sqrt{2 \times n_params}$.
<code>jitter_size</code>	positive scalar, noise is added during crossover step from <code>Uniform(-jitter_size, jitter_size)</code> distribution. $1e-6$ is the default value.
<code>parallel_type</code>	string specifying parallelization type. 'none', 'FORK', or 'PSOCK' are valid values. 'none' is default value.
<code>use_QMC</code>	logical, if true, a quasi-Monte Carlo estimator is used to estimate ELBO during optimization. default is TRUE.
<code>purify</code>	an integer, every 'purify'-th iteration, the Monte Carlo estimator of the ELBO is recalculated. This can help deal with noisy and outlier estimates of the ELBO. Default value is 25. If <code>use_QMC</code> is TRUE, purification is disabled as it is redundant.
<code>quasi_rand_seq</code>	type of low discrepancy sequence used for quasi Monte Carlo integration, either 'sobol' or 'halton'. LRVB correction always use QMC. Default is 'sobol'.
<code>n_samples_ELBO</code>	number of samples used for the Monte Carlo estimator of the ELBO (the objective function). default is 10.
<code>LRVB_correction</code>	logical, if true, LRVB covariance correction (Giordano, Brodderick, & Jordan 2018; Galdo, Bahg, & Turner 2020) is attempted.
<code>n_samples_LRVB</code>	number of samples used for LRVB correction. default is 25.
<code>neg_inf</code>	if density for a given value of theta is numerically 0 for q, this value is assigned for log density. This helps with numeric stability of algorithm. Default value is -750.
<code>thin</code>	positive integer, only every 'thin'-th iteration will be stored. Default value is 1. Increasing thin will reduce the memory required, while running algorithm for longer.
<code>burnin</code>	number of initial iterations to discard. Default value is 0.
<code>return_trace</code>	logical, if true, function returns particle trajectories. This is helpful for diagnosing convergence or debugging model code. Function will return an iteration/thin $\times n_chains \times n_params$ array and the estimated ELBO of each particle in a iteration/thin $\times n_chains$ array.
<code>crossover_rate</code>	number on the interval (0,1]. Determines the probability a parameter on a chain is updated on a given crossover step, sampled from a Bernoulli distribution.

Value

list of control parameters for the DEVI function

DEMAP	<i>DEMAP</i>
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Description

DE optimization for maximum a posteriori (MAP) estimation; this function tries to find the posterior mode.

Usage

```
DEMAP(LogPostLike, control_params = AlgoParamsDEMAP(), ...)
```

Arguments

LogPostLike	function whose first argument is an <code>n_params</code> -dimensional model parameter vector and returns (scalar) sum of log prior density and log likelihood for the parameter vector.
control_params	control parameters for DE algorithm. see AlgoParamsDEMAP function documentation for more details.
...	additional arguments to pass LogPostLike

Value

list contain posterior samples from DEMCMC in a `n_iters_per_chain` by `n_chains` by `n_params` array and the log likelihood of each sample in a `n_iters_per_chain` by `n_chains` array.

Examples

```
# simulate from model
dataExample <- matrix(stats::rnorm(100, c(-1, 1), c(1, 1)), nrow = 50, ncol = 2, byrow = TRUE)

# list parameter names
param_names_example <- c("mu_1", "mu_2")

# log posterior likelihood function = log likelihood + log prior | returns a scalar
LogPostLikeExample <- function(x, data, param_names) {
  out <- 0

  names(x) <- param_names

  # log prior
  out <- out + sum(dnorm(x["mu_1"], 0, sd = 1, log = TRUE))
  out <- out + sum(dnorm(x["mu_2"], 0, sd = 1, log = TRUE))

  # log likelihoods
```

```

    out <- out + sum(dnorm(data[, 1], x["mu_1"], sd = 1, log = TRUE))
    out <- out + sum(dnorm(data[, 2], x["mu_2"], sd = 1, log = TRUE))

    return(out)
}

# Get map estimates
DEMAP(
  LogPostLike = LogPostLikeExample,
  control_params = AlgoParamsDEMAP(
    n_params = length(param_names_example),
    n_iter = 1000,
    n_chains = 12
  ),
  data = dataExample,
  param_names = param_names_example
)

```

DEMCMC

*DEMCMC***Description**

Sample from posterior using Differential Evolution Markov Chain Monte Carlo

Usage

```
DEMCMC(LogPostLike, control_params = AlgoParamsDEMCMC(), ...)
```

Arguments

LogPostLike	function whose first argument is an <code>n_params</code> -dimensional model parameter vector and returns (scalar) sum of log prior density and log likelihood for the parameter vector.
control_params	control parameters for DEMCMC algorithm. see AlgoParamsDEMCMC function documentation for more details. You must specify ' <code>n_params</code> ' here.
...	additional arguments to pass LogPostLike

Value

list contain posterior samples from DEMCMC in a '`n_samples_per_chain`' by '`n_chains`' by `n_params` array and the log posterior likelihood of each sample in a '`n_samples_per_chain`' by '`n_chains`' array.

Examples

```
# simulate from model
dataExample <- matrix(stats::rnorm(100, c(-1, 1), c(1, 1)), nrow = 50, ncol = 2, byrow = TRUE)
#
# list parameter names
param_names_example <- c("mu_1", "mu_2")

# log posterior likelihood function = log likelihood + log prior | returns a scalar
LogPostLikeExample <- function(x, data, param_names) {
  out <- 0

  names(x) <- param_names

  # log prior
  out <- out + sum(dnorm(x["mu_1"], 0, sd = 1, log = TRUE))
  out <- out + sum(dnorm(x["mu_2"], 0, sd = 1, log = TRUE))

  # log likelihoods
  out <- out + sum(dnorm(data[, 1], x["mu_1"], sd = 1, log = TRUE))
  out <- out + sum(dnorm(data[, 2], x["mu_2"], sd = 1, log = TRUE))

  return(out)
}

# Sample from posterior
DEMCMC(
  LogPostLike = LogPostLikeExample,
  control_params = AlgoParamsDEMCMC(
    n_params = length(param_names_example),
    n_iter = 1000,
    n_chains = 12
  ),
  data = dataExample,
  param_names = param_names_example
)
```

DEVI

DEVI

Description

DE optimization for mean-field variational inference. Minimizes the KL divergence (maximizes the ELBO) between $q(\theta|\lambda)$ and the target posterior $p(\theta|data)$. For a tutorial on variational inference check out Galdo, Bahg, & Turner 2020.

Usage

```
DEVI(LogPostLike, control_params = AlgoParamsDEVI(), ...)
```

Arguments

`LogPostLike` function whose first argument is an `n_params`-dimensional model parameter vector and returns (scalar) sum of log prior density and log likelihood for the parameter vector.

`control_params` control parameters for DE algorithm. see [AlgoParamsDEVI](#) function documentation for more details.

`...` additional arguments to pass `LogPostLike`

Value

list contain mean in a `n_iters_per_chain` by `n_chains` by `2*n_params_model` array and the ELBO of each sample in a `n_iters_per_chain` by `n_chains` array.

Examples

```
# simulate from model
dataExample <- matrix(stats::rnorm(100, c(-1, 1), c(1, 1)), nrow = 50, ncol = 2, byrow = TRUE)
## list parameter names
param_names_example <- c("mu_1", "mu_2")

# log posterior likelihood function = log likelihood + log prior | returns a scalar
LogPostLikeExample <- function(x, data, param_names) {
  out <- 0

  names(x) <- param_names

  # log prior
  out <- out + sum(dnorm(x["mu_1"], 0, sd = 1, log = TRUE))
  out <- out + sum(dnorm(x["mu_2"], 0, sd = 1, log = TRUE))

  # log likelihoods
  out <- out + sum(dnorm(data[, 1], x["mu_1"], sd = 1, log = TRUE))
  out <- out + sum(dnorm(data[, 2], x["mu_2"], sd = 1, log = TRUE))

  return(out)
}

# Get variational approximation
DEVI(
  LogPostLike = LogPostLikeExample,
  control_params = AlgoParamsDEVI(
    n_params = length(param_names_example),
    n_iter = 200,
    n_chains = 12
  ),
  data = dataExample,
  param_names = param_names_example
)
```

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