

# Package ‘BaPreStoPro’

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**Type** Package

**Title** Bayesian Prediction of Stochastic Processes

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**Description** Bayesian inference and prediction for several kinds of stochastic processes, e.g. jump diffusion, (mixed) diffusion models.

**License** GPL (>= 2)

**Depends** mvtnorm,stats,methods

**LazyData** TRUE

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BaPreStoPro-package	<i>Bayesian Prediction of Stochastic Processes</i>
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## Description

This package contains simulate, estimate and predict methods for jump diffusions, diffusions, mixed diffusions, hidden (mixed) diffusion models and regression models for comparison.

## Details

Package:	BaPreStoPro
Type:	Package
Version:	1.0
Date:	2016-04-25
License:	GLP-2, GLP-3

An overview of how to use the package, including the most important functions

## Author(s)

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## References

Bayesian Prediction of Crack Growth Based on a Hierarchical Diffusion Model. S. Hermann, K. Ickstadt and C. Mueller, *appearing in: Applied Stochastic Models in Business and Industry 2016*.

## Examples

```
model <- set.to.class("Diffusion", parameter = list(phi = 0.5, gamma2 = 0.01))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5, plot.series = TRUE)
est_diff <- estimate(model, t, data, 100) # better: 10000
plot(est_diff)
pred_diff <- predict(est_diff, b.fun.mat = function(phi, t, y) phi[,1]) # much faster
```

---

ad.propSd	<i>Helping function</i>
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---

**Description**

Adaptive MCMC

**Usage**

```
ad.propSd(chain, propSd, iteration, lower = 0.3, upper = 0.6,
  delta.n = function(n) min(0.05, 1/sqrt(n)))
```

**Arguments**

chain	Markov chain
propSd	current proposal standard deviation
iteration	current iteration (batch)
lower	lower bound
upper	upper bound
delta.n	function of batch number

**Value**

adjusted proposal standard deviation

---

BinSearch	<i>Binary Search Algorithm</i>
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---

**Description**

Binary Search Algorithm

**Usage**

```
BinSearch(Fun, len, candArea, grid = 1e-05, method = c("vector", "free"))
```

**Arguments**

Fun	cumulative distribution function
len	number of samples
candArea	candidate area
grid	fineness degree
method	vectorial ("vector") or not ("free")

**Value**

vector of samples

**Examples**

```
test <- BinSearch(function(x) pnorm(x, 5, 1), 1000, candArea = c(0, 10), method = "free")
plot(density(test))
curve(dnorm(x, 5, 1), col = 2, add = TRUE)
```

---

class.to.list	<i>Builds a list from class</i>
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---

**Description**

Class to list

**Usage**

```
class.to.list(cl)
```

**Arguments**

cl	class
----	-------

**Value**

list

---

diagnostic	<i>Calcucation of burn-in phase and thin rate</i>
------------	---------------------------------------------------

---

**Description**

Proposal for burn-in and thin rate

**Usage**

```
diagnostic(chain, dependence = 0.8, m = 10)
```

**Arguments**

chain	vector of Markov chain samples
dependence	allowed dependence for the chain
m	number of blocks

---

dNtoTimes

*Transformation of counting process to vector of event times*


---

**Description**

Transformation of vector of counting process to event times.

**Usage**

```
dNtoTimes(dN, t)
```

**Arguments**

dN                      vector of differences of counting process  
t                        times of counting process

**Value**

vector of event times

**Examples**

```
t <- seq(0, 1, by = 0.01)
process <- simN(t, c(5, 0.5), len = 1)$N
times <- dNtoTimes(diff(process), t)
```

---

drawSDE

*Function for simulating diffusion process*


---

**Description**

Simulation of process defined by  $dY_t = b(\phi, t, Y_t)dt + \gamma \text{sigmaTilde}(t, Y_t)dW_t$ .

**Usage**

```
drawSDE(phi, gamma2, t, b, fODE, sigmaTilde, mw = 10,
        strictly.positive = TRUE)
```

**Arguments**

phi                      parameter  $\phi$   
gamma2                   parameter  $\gamma^2$   
t                        vector of time points  
b                        drift function  
fODE                    function for the starting point dependent on phi, for fixed y0: fODE = function(phi, t) y0  
sigmaTilde              variance function  $s(\gamma, t, y) = \gamma \text{sigmaTilde}(t, y)$   
mw                      mesh width to simulate the time-continuity  
strictly.positive       if TRUE, only positive values for process  $Y_t$

**Value**

data series in t

---

estimate	<i>Estimation</i>
----------	-------------------

---

**Description**

Method for the S4 classes

**Usage**

```
estimate(model.class, ...)
```

**Arguments**

model.class	class
...	parameters dependent on the model class

---

estimate, Diffusion-method
<i>Estimation for diffusion process</i>

---

**Description**

Bayesian estimation of the parameters of the stochastic process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ .

**Usage**

```
## S4 method for signature 'Diffusion'
estimate(model.class, t, data, nMCMC)
```

**Arguments**

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

**Examples**

```
model <- set.to.class("Diffusion", parameter = list(phi = 0.5, gamma2 = 0.01))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5, plot.series = TRUE)
est_diff <- estimate(model, t, data, 1000)
plot(est_diff)
```

---

estimate,hiddenDiffusion-method

*Estimation for noisy / hidden diffusion process*


---

## Description

Bayesian estimation of the model,  $Z_i = Y_{t_i} + \epsilon_i$ ,  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ .

## Usage

```
## S4 method for signature 'hiddenDiffusion'
estimate(model.class, t, data, nMCMC, Npart = 100)
```

## Arguments

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain
Npart	number of particles in the particle Gibbs sampler

## Examples

```
model <- set.to.class("hiddenDiffusion", y0.fun = function(phi, t) 0.5,
  parameter = list(phi = 5, gamma2 = 1, sigma2 = 0.1))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est <- estimate(model, t, data$Z, 100) # nMCMC should be much larger!
plot(est)
## Not run:
# OU
b.fun <- function(phi, t, y) phi[1]-phi[2]*y
model <- set.to.class("hiddenDiffusion", y0.fun = function(phi, t) 0.5,
  parameter = list(phi = c(10, 5), gamma2 = 1, sigma2 = 0.1),
  b.fun = b.fun, sT.fun = function(t, x) 1)
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est <- estimate(model, t, data$Z, 1000)
plot(est)

## End(Not run)
```



---

estimate,hiddenmixedDiffusion-method

*Estimation for noisy/hidden mixed diffusion process*


---

### Description

Bayesian estimation of a stochastic process  $Z_{ij} = Y_{t_{ij}} + \epsilon_{ij}$ ,  $dY_t = b(\phi_j, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ ,  $\phi_j \sim N(\mu, \Omega)$ .

### Usage

```
## S4 method for signature 'hiddenmixedDiffusion'
estimate(model.class, t, data, nMCMC,
  Npart = 100)
```

### Arguments

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain
Npart	number of particles in the particle Gibbs sampler

### Examples

```
mu <- c(5, 1); Omega <- c(0.9, 0.04)
phi <- cbind(rnorm(21, mu[1], sqrt(Omega[1])), rnorm(21, mu[2], sqrt(Omega[2])))
y0.fun <- function(phi, t) phi[2]
model <- set.to.class("hiddenmixedDiffusion", y0.fun = y0.fun,
  b.fun = function(phi, t, y) phi[1],
  parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 1, sigma2 = 0.01))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
## Not run:
est <- estimate(model, t, data$Z[1:20,], 2000)
plot(est)

## End(Not run)
```

---

estimate,jumpDiffusion-method

*Estimation for jump diffusion process*


---

### Description

Bayesian estimation of a stochastic process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t + h(\eta, t, Y_t)dN_t$ .

**Usage**

```
## S4 method for signature 'jumpDiffusion'
estimate(model.class, t, data, nMCMC)
```

**Arguments**

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

**Examples**

```
model <- set.to.class("jumpDiffusion", Lambda = function(t, xi) (t/xi[2])^xi[1],
  parameter = list(theta = 0.1, phi = 0.05, gamma2 = 0.1, xi = c(3, 1/4)))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5, plot.series = TRUE)
est <- estimate(model, t, data, 1000)
plot(est)
```

---

estimate, Merton-method

*Estimation for jump diffusion process*

---

**Description**

Bayesian estimation of a stochastic process  $Y_t = y_0 \exp(\phi t - \gamma^2/2t + \gamma W_t + \log(1 + \theta)N_t)$ .

**Usage**

```
## S4 method for signature 'Merton'
estimate(model.class, t, data, nMCMC)
```

**Arguments**

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

**Examples**

```

model <- set.to.class("Merton", parameter = list(thetaT = 0.1, phi = 0.05, gamma2 = 0.1, xi = 10))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5, plot.series = TRUE)
est <- estimate(model, t, data, 1000)
plot(est)
## Not run:
est_hidden <- estimate(model, t, data$Y, 1000)
plot(est_hidden)

## End(Not run)

```

---

estimate,mixedDiffusion-method

*Estimation for mixed diffusion process*


---

**Description**

Bayesian estimation of a stochastic process  $dY_t = b(\phi_j, t, Y_t)dt + s(\gamma, t, Y_t)dW_t, \phi_j \sim N(\mu, \Omega)$ .

**Usage**

```

## S4 method for signature 'mixedDiffusion'
estimate(model.class, t, data, nMCMC)

```

**Arguments**

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

**Examples**

```

mu <- 2; Omega <- 0.4; phi <- matrix(rnorm(21, mu, sqrt(Omega)))
model <- set.to.class("mixedDiffusion",
  parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1),
  b.fun = function(phi, t, x) phi*x, sT.fun = function(t, x) x)
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est <- estimate(model, t, data[1:20,], 100) # nMCMC should be much larger
plot(est)
# OU
b.fun <- function(phi, t, y) phi[1]-phi[2]*y; y0.fun <- function(phi, t) phi[3]
mu <- c(10, 5, 0.5); Omega <- c(0.9, 0.01, 0.01)
phi <- sapply(1:3, function(i) rnorm(21, mu[i], sqrt(Omega[i])))
model <- set.to.class("mixedDiffusion",
  parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1),
  y0.fun = y0.fun, b.fun = b.fun, sT.fun = function(t, x) 1)
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est <- estimate(model, t, data[1:20,], 100) # nMCMC should be much larger
plot(est)

```

---

 estimate,mixedRegression-method

*Estimation for mixed regression model*


---

### Description

Bayesian estimation of the parameter of the regression model  $y_i = f(\phi_j, t_i) + \epsilon_i, \phi_j \sim N(\mu, \Omega)$ .

### Usage

```
## S4 method for signature 'mixedRegression'
estimate(model.class, t, data, nMCMC)
```

### Arguments

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

### Examples

```
mu <- c(10, 5); Omega <- c(0.9, 0.01)
phi <- cbind(rnorm(21, mu[1], sqrt(Omega[1])), rnorm(21, mu[2], sqrt(Omega[2])))
model <- set.to.class("mixedRegression",
  parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1),
  fun = function(phi, t) phi[1]*t + phi[2], sT.fun = function(t) 1)
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est <- estimate(model, t, data[1:20,], 2000)
plot(est)
```

---

 estimate,NHPP-method    *Estimation for Poisson process*


---

### Description

Bayesian estimation of a nonhomogeneous Poisson process.

### Usage

```
## S4 method for signature 'NHPP'
estimate(model.class, t, data, nMCMC)
```

**Arguments**

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

**Examples**

```
model <- set.to.class("NHPP", parameter = list(xi = c(5, 1/2)),
  Lambda = function(t, xi) (t/xi[2])^xi[1])
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est_NHPP <- estimate(model, t, data$Times, 10000)
plot(est_NHPP)
```

---

estimate,Regression-method

*Estimation for regression model*


---

**Description**

Bayesian estimation of the parameter of the regression model  $y_i = f(\phi, t_i) + \epsilon_i$ .

**Usage**

```
## S4 method for signature 'Regression'
estimate(model.class, t, data, nMCMC)
```

**Arguments**

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

**Examples**

```
t <- seq(0,1, by = 0.01)
model <- set.to.class("Regression", fun = function(phi, t) phi[1]*t + phi[2],
  parameter = list(phi = c(1,2), gamma2 = 0.1))
data <- simulate(model, t = t, plot.series = TRUE)
est <- estimate(model, t, data, 1000)
plot(est)
```

---

```
estimate, reg_hiddenNHPP-method
```

*Estimation for regression model dependent on Poisson process*

---

### Description

Bayesian estimation of the parameter of the regression model  $y_i = f(t_i, N_i, \theta) + \epsilon_i$ .

### Usage

```
## S4 method for signature 'reg_hiddenNHPP'
estimate(model.class, t, data, nMCMC)
```

### Arguments

model.class	class of the respective model including all required information, see function set.to.class
t	vector of time points
data	vector or list or matrix of observation variables
nMCMC	length of Markov chain

### Examples

```
t <- seq(0,1, by = 0.01)
model <- set.to.class("reg_hiddenNHPP", fun = function(t, N, theta) exp(theta[1]*t) + theta[2]*N,
  parameter = list(theta = c(2, 2), gamma2 = 0.25, xi = c(3, 0.5)),
  Lambda = function(t, xi) (t/xi[2])^xi[1])
data <- simulate(model, t = t)
est <- estimate(model, t, data, 1000)
plot(est)
## Not run:
est_hid <- estimate(model, t, data$Y, 1000)
plot(est_hid)

## End(Not run)
```

---

```
estReg
```

*Bayesian estimation in mixed nonlinear regression models*

---

### Description

Bayesian estimation of the random effects  $\phi_j$  in the mixed nonlinear regression model  $y_{ij} = f(\phi_j, t_{ij}) + \epsilon_{ij}$ ,  $\epsilon_{ij} \sim N(0, \gamma^2 * s^2(t_{ij}))$ ,  $\phi_j \sim N(\mu, \Omega)$  and the parameters  $\mu, \Omega, \gamma^2$ .

### Usage

```
estReg(t, y, prior, start, fODE, sVar, ipred = 1, cut, len = 1000,
  mod = c("Gompertz", "logistic", "Weibull", "Richards", "Paris", "Paris2"),
  propPar = 0.02)
```

**Arguments**

t	vector of observation times
y	matrix of the M trajectories
prior	list of prior parameters - list(m, v, priorOmega, alpha, beta)
start	list of starting values
fODE	regression function
sVar	variance function
ipred	which of the M trajectories is the one to be predicted
cut	the index how many of the ipred-th series are used for estimation
len	number of iterations of the MCMC algorithm
mod	model out of Gompertz, Richards, logistic, Weibull, only used instead of fODE
propPar	proposal standard deviation of phi is lstart\$mul*propPar

**Value**

phi	samples from posterior of $\phi$
mu	samples from posterior of $\mu$
Omega	samples from posterior of $\Omega$
gamma2	samples from posterior of $\gamma^2$

estReg\_single

*Bayesian stimation in nonlinear regression models***Description**

Bayesian estimation of the parameters of the mixed nonlinear regression model  $y_j = f(\phi, t_j) + \epsilon_j$ ,  $\epsilon_j \sim N(0, \gamma^2 * s^2(t_j))$ .

**Usage**

```
estReg_single(t, y, prior, start, fODE, sVar, len = 1000, mod = "Gompertz")
```

**Arguments**

t	vector of observation times
y	vector of the M trajectories
prior	list of prior parameters - list(mu, Omega, alpha, beta)
start	list of starting values
fODE	regression function
sVar	variance function
len	number of iterations of the MCMC algorithm
mod	model out of Gompertz, Richards, logistic, Weibull, only used instead of fODE

**Value**

phi	estimator of $\phi$
gamma2	estimator of $\gamma^2$

estSDE

*Bayesian estimation in mixed stochastic differential equations***Description**

Bayesian estimation of the random effects  $\phi_i$  in the mixed SDE  $dY_i(t) = b(\phi_i, t, Y_i(t))dt + \gamma s(t, Y_i(t))dW_i(t)$ ,  $\phi_i \sim N(\mu, \Omega)$ ,  $i = 1, \dots, n$  and the parameters  $\mu, \Omega, \gamma^2$ .

**Usage**

```
estSDE(t, y, prior, start, y0.fun, bSDE, sVar, ipred = 1, cut, len = 1000,
      mod = c("Gompertz", "logistic", "Weibull", "Richards", "Paris", "Paris2"),
      propPar = 0.2)
```

**Arguments**

t	vector of observation times
y	matrix or list of the n trajectories
prior	list of prior parameters - list(m, v, alpha.omega, beta.omega, alpha.gamma, beta.gamma)
start	list of starting values
y0.fun	$y_0(\phi, t_1)$ function
bSDE	$b(\phi, t, x)$ drift function
sVar	variance function $s^2$
ipred	which of the n trajectories is the one to be predicted
cut	the index how many of the ipred-th series are used for estimation
len	number of iterations of the MCMC algorithm - chain length
mod	model out of Gompertz, Richards, logistic, Weibull, Paris, Paris2, only used instead of bSDE
propPar	proposal standard deviation of phi is $lstart\$mul*propPar$

**Details**

Simulation from the posterior distribution of the random effect from n independent trajectories of the SDE (the Brownian motions  $W_1, \dots, W_n$  are independent).

**Value**

phi	samples from posterior of $\phi$
mu	samples from posterior of $\mu$
Omega	samples from posterior of $\Omega$
gamma2	samples from posterior of $\gamma^2$

**References**

Hermann et al. (2015)



---

estSDE_single	<i>Bayesian estimation in stochastic differential equations</i>
---------------	-----------------------------------------------------------------

---

**Description**

Bayesian estimation of the parameters in the SDE  $dY(t) = b(\phi, t, Y(t))dt + \gamma s(t, Y(t))dW(t)$ .

**Usage**

```
estSDE_single(t, X, prior, start, bSDE, sVar, len = 1000)
```

**Arguments**

t	vector of observation times
X	vector of the M trajectories
prior	list of prior parameters - list(mu, Omega, alpha, beta)
start	list of starting values
bSDE	drift function
sVar	variance function
len	number of iterations of the MCMC algorithm

**Details**

Simulation from the posterior distribution of the random effect from n independent trajectories of the SDE (the Brownian motions  $W_1, \dots, W_n$  are independent).

**Value**

phi	estimator of $\phi$
gamma2	estimator of $\gamma^2$

---

est_JD_Euler	<i>Metropolis within Gibbs sampler</i>
--------------	----------------------------------------

---

**Description**

Bayesian estimation of the parameter of the jump diffusion process define by SDE  $dXt = b(\phi, t, Xt)dt + s(\gamma, t, Xt)dWt + h(\theta, t, Xt)dNt$ .

**Usage**

```
est_JD_Euler(X, N, t, n = 1000, start, b, s, h, priorRatio, Lambda,
  int = c("Weibull", "Exp"), rangeN = 2, propSd = 0.5, it.xi = 5)
```

**Arguments**

X	vector of observed variables
N	vector of Poisson process variables
t	vector of time points
n	length of Markov chain
start	list of starting values, list(phi, theta, gamma2)
b	drift function
s	variance function
h	jump high function
priorRatio	list of functions for the prior ratio of MH step, if missing: non-informative
Lambda	intensity rate function
int	if Lambda is missing, one of "Weibull" or "Exp"
rangeN	range for candidates for filtering N
propSd	starting value for proposal standard deviation
it.xi	number of iterations of MH inside the Gibbs sampler

**Value**

phi	estimator of $\phi$
gamma2	estimator of $\gamma^2$
theta	estimator of $\theta$
xi	estimator of $\xi$
N	estimator of latent variable $N$ , if not observed
prop	storage of adapted proposal variances

---

est\_Merton

*Gibbs sampler*


---

**Description**

Bayesian estimation of the parameter of the jump diffusion process  $X_t = x_0 \exp(\phi t - \gamma^2/2t + \gamma W_t + \log(1 + \theta)N_t)$ .

**Usage**

```
est_Merton(X, N, t, n = 1000, start, prior, Lambda, rangeN = 2,
  it.xi = 10)
```

**Arguments**

X	vector of observed variables
N	vector of Poisson process variables, optional: if missing, filtering
t	vector of time points
n	length of Markov chain
start	list of starting values, list(thetaT, gamma2)
prior	list of prior values, list(mu_phi, s_phi, mu_th, s_th, alpha, beta)
Lambda	intensity rate function
rangeN	range of candidates for filtering N
it.xi	number of iterations of MH inside the Gibbs sampler

**Value**

phi	estimator of $\phi$
gamma2	estimator of $\gamma^2$
thetaT	estimator of $\log(1 + \theta)$
xi	estimator of $\xi$
N	estimator of latent variable $N$ , if not observed

---

est_NHPP	<i>Metropolis-Hastings sampler</i>
----------	------------------------------------

---

**Description**

Bayesian estimation of the parameter of the intensity rate.

**Usage**

```
est_NHPP(jumpTimes, Tend, start, n = 5000, int = c("Weibull", "Exp"),
  priorRatio, proposal = c("lognormal", "normal"), Lambda)
```

**Arguments**

jumpTimes	vector of jump (or event) times
Tend	last observation vector of the NHPP
start	starting value
n	length of Markov chain
int	one out of "Weibull" or "Exp"
priorRatio	function for prior ratio, if missing: non-informative
proposal	"lognormal" (for positive parameters, default) or "normal"
Lambda	intensity rate function

**Value**

p x n dimensional matrix of posterior samples

## Examples

```
# exponential
Lambda <- function(t, xi){
  xi[2]*exp(xi[1]*t)-xi[2]
}
jumpTimes <- simN(seq(0, 1, by = 0.001), c(4, 0.5), len = 1, Lambda = Lambda)$Times
chain <- est_NHPP(jumpTimes, 1, c(4, 0.5), Lambda = Lambda)
# chain <- est_NHPP(jumpTimes, 1, c(4, 0.5), int = "Exp")
plot(chain[1,], type="l"); abline(h = 4)
plot(chain[2,], type="l"); abline(h = 0.5)

# weibull
Lambda <- function(t, xi){
  (t/xi[2])^xi[1]
}
jumpTimes <- simN(seq(0, 1, by = 0.001), c(3, 0.5), len = 1, Lambda = Lambda)$Times
chain <- est_NHPP(jumpTimes, 1, c(3, 0.5), Lambda = Lambda)
# chain <- est_NHPP(jumpTimes, 1, c(3, 0.5), int = "Weibull")
plot(chain[1,], type="l"); abline(h = 3)
plot(chain[2,], type="l"); abline(h = 0.5)
```

---

est\_reg\_hiddenNHPP      *Gibbs sampler*

---

## Description

Bayesian estimation of the parameter of the regression model  $y_i = f(t_i, N_i, \theta) + \epsilon_i$ .

## Usage

```
est_reg_hiddenNHPP(Y, N, t, fun, n = 1000, start, prior, Lambda,
  int = c("Weibull", "Exp"), rangeN = 2)
```

## Arguments

Y	vector of observation variables
N	vector of Poisson process variables
t	vector of time points
fun	regression function
n	length of Markov chain
start	list of starting values
prior	list of prior values
Lambda	intensity rate function
int	one out of "Weibull" or "Exp", if Lambda is missing
rangeN	range for candidates of filtering N, if unobserved

**Value**

theta	Markov chains of $\theta$
gamma2	Markov chains of $\gamma^2$
N	if hidden: Markov chains of $N$
xi	Markov chains of $\xi$

---

est_SEP	<i>Metropolis-Hastings sampler</i>
---------	------------------------------------

---

**Description**

Bayesian estimation of the parameter of the self-exciting process.

**Usage**

```
est_SEP(jumpTimes, Tend, start, n = 5000, s = 200, f_xi, priorRatio,
        proposals, Lmax = 35)
```

**Arguments**

jumpTimes	vector of event times
Tend	last observation time, if missing => last event time
start	starting value
n	length of Markov chain
s	stress range
f_xi	exp(-f_xi) intensity function
priorRatio	function(old, new) for prior ratio, if missing => noninformative
proposals	list of functions: draw(old) and ratio(new, old)
Lmax	maximal number of tension wires that can break

**Value**

p x n dimensional matrix of posterior samples, p length of vector xi

**Examples**

```
wt <- rexp(20, exp(- 1 + 0.25*200/(35-(0:19))))
jumpTimes <- cumsum(wt)
chain <- est_SEP(jumpTimes, start = c(1, 0.25), s = 200)
par(mfrow = c(2,1))
plot(chain[1,], type = "l")
abline(h = 1, col = 2)
plot(chain[2,], type = "l")
abline(h = 0.25, col = 2)

print("acceptance rate:")
length(unique(chain[1,]))/length(chain[1,])

# or:
```

```

s <- seq(100, 300, length = 10)
jumpTimes <- lapply(1:10, function(a){
  wt <- rexp(20, exp(- 1 + 0.25*s[a]/(35-(0:19))))
  cumsum(wt)
})
chain <- est_SEP(jumpTimes, start = c(1, 0.25), s = s)

```

---

findCandidateArea	<i>Helping function</i>
-------------------	-------------------------

---

### Description

Finding suitable candidate area

### Usage

```
findCandidateArea(VFun, start = 1, pos.support = TRUE, quasi.null = 1e-05)
```

### Arguments

VFun	cumulative distribution function
start	starting point for search
pos.support	if TRUE: only positive support
quasi.null	size of values to be defined as zero, default: 10 <sup>-5</sup>

### Value

adjusted proposal standard deviation

---

getFun	<i>Yields the regression or drift function for the specific models</i>
--------	------------------------------------------------------------------------

---

### Description

Gives out the regression function (ODE) or the drift function (SDE) for the specific model out of Gompertz, Richards, logistic, Weibull

### Usage

```
getFun(model, mod)
```

### Arguments

model	one out of SDE, ODE
mod	one out of Gompertz, Richards, logistic, Weibull

### Value

function

---

getPrior	<i>Builds list of prior parameters</i>
----------	----------------------------------------

---

**Description**

Creation of list of parameters conditional on true values  $\mu$  and  $\gamma^2$

**Usage**

```
getPrior(parameter, model = c("jumpDiffusion", "Merton", "Diffusion",
  "mixedDiffusion", "hiddenDiffusion", "hiddenmixedDiffusion", "reg_hiddenNHPP",
  "NHPP", "Regression", "mixedRegression"))
```

**Arguments**

parameter	list of parameters
model	name of model

**Value**

list of prior values

---

jumpDiffusion-class	<i>S4 class for the jump diffusion process</i>
---------------------	------------------------------------------------

---

**Description**

S4 class for the jump diffusion process

**Slots**

theta
phi ...

---

partFiltering	<i>Estimation function</i>
---------------	----------------------------

---

**Description**

Bayesian estimation of the parameter of the model  $Y_i = X_{t_i} + \epsilon_i$  with  $dX_t = b(\phi, t, X_t)dt + s(\gamma, t, X_t)dW_t$ .

**Usage**

```
partFiltering(t, y, prior, start, est = c("Bayes", "MLE"), len = 1000,
  sigmaTilde, Npart = 10, y0.fun = 1, b.fun = 1, parPropPhi = 5,
  maxIt = 10)
```

**Arguments**

t	vector of time points
y	vector of observation variables
prior	list of prior values
start	list of starting values
est	one out of "Bayes" or "MLE"
len	length of Markov chain
sigmaTilde	variance function $s(\gamma, t, y) = \gamma \text{sigmaTilde}(t, y)$
Npart	number of particles
y0.fun	function for starting point dependent on $\phi$
b.fun	drift function
parPropPhi	parameter for proposal standard deviation
maxIt	maximal iteration of MH step of $\phi$

**Value**

phi	Markov chain of $\phi$
gamma2	Markov chain of $\gamma^2$
sigma2	Markov chain of $\sigma^2$
X	filtered process

**Author(s)**

Simone Hermann and Adeline Leclercq-Samson

---

partFiltering\_mixed     *Estimation function*

---

**Description**

Bayesian estimation of the parameter of the model  $Y_{ij} = X_{t_{ij}} + \epsilon_{ij}$  with  $dX_t = b(\phi_j, t, X_t)dt + s(\gamma, t, X_t)dW_t$ .

**Usage**

```
partFiltering_mixed(t, y, prior, start, len = 1000, sigmaTilde, y0.fun = 1,
  b.fun = 1, Npart = 10, parPropPhi = 5, maxIt = 10)
```

**Arguments**

t	vector of time points
y	matrix of observation variables
prior	list of prior values
start	list of starting values
len	length of Markov chain



sigmaTilde	variance function $s(\gamma, t, y) = \gamma \text{sigmaTilde}(t, y)$
y0.fun	function for starting point dependent on $\phi$
b.fun	drift function
Npart	number of particles
parPropPhi	parameter for proposal standard deviation
maxIt	maximal iteration of MH step of $\phi$

**Value**

phi	Markov chain of $\phi$
mu	Markov chain of $\mu$
Omega	Markov chain of $\Omega$
gamma2	Markov chain of $\gamma^2$
sigma2	Markov chain of $\sigma^2$
x	filtered process

**Author(s)**

Simone Hermann and Adeline Leclercq-Samson

---

phi_ij	<i>Helping function</i>
--------	-------------------------

---

**Description**

Returns the ij th matrix entry from a list

**Usage**

```
phi_ij(phi, i, j)
```

**Arguments**

phi	list with each entry a matrix
i	row
j	column

**Value**

vector of samples

---

plot,est.Diffusion-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.Diffusion

## Usage

```
## S4 method for signature 'est.Diffusion'
plot(x, par.options, style = c("chains", "acf",
    "density"), par2plot, reduced = TRUE, thinning, burnIn, priorMeans = TRUE,
    col.priorMean = 2, lty.priorMean = 1, newwindow = FALSE, ...)
```

## Arguments

x	est.Diffusion class
par.options	list of options for function par()
style	one out of "chains", "acf", "density"
par2plot	logical vector, which parameters to be plotted, order: $(\phi, \gamma^2)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
model <- set.to.class("Diffusion", b.fun = function(phi, t, y) phi[1]-phi[2]*y,
    parameter = list(phi = c(10, 1), gamma2 = 0.1))
data <- simulate(model, t = seq(0, 1, by = 0.01), y0 = 0.5, plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.01), data, 1000) # nMCMC small for example
plot(est)
plot(est, burnIn = 100, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(3,1)), xlab = "iteration")
plot(est, style = "acf", main = "", par2plot = c(TRUE, TRUE, FALSE))
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), main = "", par2plot = c(FALSE, FALSE, TRUE))
plot(est, priorMeans = FALSE, newwindow = TRUE)
```

---

plot,est.hiddenDiffusion-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.hiddenDiffusion

## Usage

```
## S4 method for signature 'est.hiddenDiffusion'
plot(x, par.options, style = c("chains",
  "acf", "density"), par2plot, reduced = TRUE, thinning, burnIn,
  priorMeans = TRUE, col.priorMean = 2, lty.priorMean = 1,
  newwindow = FALSE, ...)
```

## Arguments

x	est.hiddenDiffusion class
par.options	list of options for function par()
style	one out of "chains", "acf", "density"
par2plot	logical vector, which parameters to be plotted, order: $(\phi, \gamma^2, \sigma^2, Y)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
model <- set.to.class("hiddenDiffusion", b.fun = function(phi, t, y) phi[1]-phi[2]*y,
  parameter = list(phi = c(10, 1), gamma2 = 1, sigma2 = 0.1),
  y0 = function(phi, t) 0.5)
data <- simulate(model, t = seq(0, 1, by = 0.01), plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.01), data$Y, 100) # nMCMC small for example
plot(est)
plot(est, par2plot = c(rep(FALSE, 3), TRUE, FALSE), ylim = c(0.001, 0.1), par.options = list())
plot(est, burnIn = 10, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(3,1)), xlab = "iteration")
plot(est, style = "acf", main = "", par2plot = c(TRUE, TRUE, FALSE, FALSE))
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), main = "", par2plot = c(FALSE, FALSE, TRUE, TRUE))
plot(est, priorMeans = FALSE, newwindow = TRUE)
```

---

plot.est.hiddenmixedDiffusion-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.hiddenmixedDiffusion

## Usage

```
## S4 method for signature 'est.hiddenmixedDiffusion'
plot(x, par.options, style = c("chains",
  "acf", "density", "int.phi"), par2plot, reduced = TRUE, thinning, burnIn,
  priorMeans = TRUE, col.priorMean = 2, lty.priorMean = 1, level = 0.05,
  phi, newwindow = FALSE, ...)
```

## Arguments

x	est.hiddenmixedDiffusion class
par.options	list of options for function par()
style	one out of "chains", "acf", "density", "int.phi"
par2plot	logical vector, which parameters to be plotted, order: $(\mu, \Omega, \gamma^2, \sigma^2, Y)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
level	level for style = "int.phi"
phi	in the case of simulation study: known values for phi
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
## Not run:
mu <- c(10, 3, 1); Omega = c(1, 0.4, 0.01)
phi <- sapply(1:3, function(i) rnorm(20, mu[i], sqrt(Omega[i])))
model <- set.to.class("hiddenmixedDiffusion", b.fun = function(phi, t, y) phi[1]-phi[2]*y,
  parameter = list(mu = mu, Omega = Omega, phi = phi, gamma2 = 1, sigma2 = 0.1),
  y0 = function(phi, t) phi[3])
data <- simulate(model, t = seq(0, 1, by = 0.02), plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.02), data$Z, 1000)
plot(est, newwindow = TRUE)
plot(est, burnIn = 10, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(2,1)), xlab = "iteration")
plot(est, style = "acf", main = "", par2plot = c(TRUE, TRUE, rep(FALSE, 7)))
plot(est, style = "density", lwd = 2, priorMean = FALSE, par2plot = c(rep(FALSE, 6), TRUE, TRUE, FALSE))
```

```

plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), main = "", par2plot = c(rep(FALSE, 6), TRUE, TRUE))
plot(est, priorMeans = FALSE, newwindow = TRUE)
plot(est, style = "int.phi", phi = phi, par2plot = c(TRUE, FALSE, FALSE))

## End(Not run)

```

---

plot.est.jumpDiffusion-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.jumpDiffusion

## Usage

```

## S4 method for signature 'est.jumpDiffusion'
plot(x, par.options, style = c("chains", "acf",
  "density"), par2plot, reduced = TRUE, thinning, burnIn, priorMeans = TRUE,
  col.priorMean = 2, lty.priorMean = 1, newwindow = FALSE, ...)

```

## Arguments

x	est.jumpDiffusion class
par.options	list of options for function par()
style	one out of "chains", "acf", "density"
par2plot	logical vector, which parameters to be plotted, order: $(\phi, \theta, \gamma^2, \xi, N)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```

model <- set.to.class("jumpDiffusion", Lambda = function(t, xi) (t/xi[2])^xi[1],
  parameter = list(theta = 0.1, phi = 0.05, gamma2 = 0.1, xi = c(3, 1/4)))
data <- simulate(model, t = seq(0, 1, by = 0.01), y0 = 0.5, plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.01), data, 1000) # nMCMC small for example
plot(est)
plot(est, burnIn = 100, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(2, 3)), xlab = "iteration")
# plot only for phi and xi ...
plot(est, style = "acf", main = "", par2plot = c(TRUE, FALSE, FALSE, TRUE, TRUE))
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), par2plot = c(TRUE, rep(FALSE, 4)), main = "")
plot(est, priorMeans = FALSE, newwindow = TRUE)

```

---

plot,est.Merton-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.Merton

## Usage

```
## S4 method for signature 'est.Merton'
plot(x, par.options, style = c("chains", "acf",
  "density"), par2plot, reduced = TRUE, thinning, burnIn, priorMeans = TRUE,
  col.priorMean = 2, lty.priorMean = 1, newwindow = FALSE, ...)
```

## Arguments

x	est.Merton class
par.options	list of options for function par()
style	one out of "chains", "acf", "density"
par2plot	logical vector, which parameters to be plotted, order: $(\phi, \tilde{\theta}, \gamma^2, \xi, N)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
model <- set.to.class("Merton", Lambda = function(t, xi) (t/xi[2])^xi[1],
  parameter = list(thetaT = 0.1, phi = 0.05, gamma2 = 0.1, xi = c(3, 1/4)))
data <- simulate(model, t = seq(0, 1, by = 0.01), y0 = 0.5, plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.01), data, 1000) # nMCMC small for example
plot(est)
plot(est, burnIn = 100, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(2, 3)), xlab = "iteration",
  # plot only for phi and xi ...
  plot(est, style = "acf", main = "", par2plot = c(TRUE, FALSE, FALSE, TRUE, TRUE))
  plot(est, style = "density", lwd = 2, priorMean = FALSE)
  plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
  plot(est, style = "acf", par.options = list(), par2plot = c(TRUE, rep(FALSE, 4)), main = "")
  plot(est, priorMeans = FALSE, newwindow = TRUE)
```

---

plot,est.mixedDiffusion-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.mixedDiffusion

## Usage

```
## S4 method for signature 'est.mixedDiffusion'
plot(x, par.options, style = c("chains", "acf",
  "density", "int.phi"), par2plot, reduced = TRUE, thinning, burnIn,
  priorMeans = TRUE, col.priorMean = 2, lty.priorMean = 1, level = 0.05,
  phi, newwindow = FALSE, ...)
```

## Arguments

x	est.mixedDiffusion class
par.options	list of options for function par()
style	one out of "chains", "acf", "density", "int.phi"
par2plot	logical vector, which parameters to be plotted, order: $(\mu, \Omega, \gamma^2)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
level	level for style = "int.phi"
phi	in the case of simulation study: known values for phi
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
mu <- c(10, 3, 1); Omega = c(1, 0.4, 0.01)
phi <- sapply(1:3, function(i) rnorm(20, mu[i], sqrt(Omega[i])))
model <- set.to.class("mixedDiffusion", b.fun = function(phi, t, y) phi[1]-phi[2]*y,
  parameter = list(mu = mu, Omega = Omega, phi = phi, gamma2 = 0.1),
  y0 = function(phi, t) phi[3], sT.fun = function(t, x) sqrt(abs(x)))
data <- simulate(model, t = seq(0, 1, by = 0.02), plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.02), data, 100) # nMCMC small for example
plot(est, newwindow = TRUE)
plot(est, burnIn = 10, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(2,1)), xlab = "iteration")
plot(est, style = "acf", main = "")
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
```

```
plot(est, style = "acf", par.options = list(), main = "", par2plot = c(rep(FALSE, 6), TRUE))
plot(est, priorMeans = FALSE, newwindow = TRUE)
plot(est, style = "int.phi", phi = phi, par2plot = c(TRUE, FALSE, FALSE))
```

---

plot,est.mixedRegression-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.mixedRegression

## Usage

```
## S4 method for signature 'est.mixedRegression'
plot(x, par.options, style = c("chains",
  "acf", "density", "int.phi"), par2plot, reduced = TRUE, thinning, burnIn,
  priorMeans = TRUE, col.priorMean = 2, lty.priorMean = 1, level = 0.05,
  phi, newwindow = FALSE, ...)
```

## Arguments

x	est.mixedRegression class
par.options	list of options for function par()
style	one out of "chains", "acf", "density", "int.phi"
par2plot	logical vector, which parameters to be plotted, order: $(\mu, \Omega, \gamma^2)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
level	level for style = "int.phi"
phi	in the case of simulation study: known values for phi
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
mu <- c(1, 3); Omega = c(0.4, 0.01)
phi <- sapply(1:2, function(i) rnorm(20, mu[i], sqrt(Omega[i])))
model <- set.to.class("mixedRegression", fun = function(phi, t) phi[1]*t + phi[2],
  parameter = list(mu = mu, Omega = Omega, phi = phi, gamma2 = 0.1))
data <- simulate(model, t = seq(0, 1, by = 0.02), plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.02), data, 100) # nMCMC small for example
plot(est, newwindow = TRUE)
plot(est, burnIn = 10, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(2,1)), xlab = "iteration")
```



```

plot(est, style = "acf", main = "")
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), main = "", par2plot = c(rep(FALSE, 4), TRUE))
plot(est, priorMeans = FALSE, newwindow = TRUE)
plot(est, style = "int.phi", phi = phi, par2plot = c(TRUE, FALSE))

```

---

plot,est.NHPP-method    *Plot method for the Bayesian estimation class object*

---

## Description

Plot method for the S4 class est.NHPP

## Usage

```

## S4 method for signature 'est.NHPP'
plot(x, par.options, style = c("chains", "acf",
    "density"), par2plot, reduced = TRUE, thinning, burnIn, priorMeans = TRUE,
    col.priorMean = 2, lty.priorMean = 1, newwindow = FALSE, ...)

```

## Arguments

x	est.NHPP class
par.options	list of options for function par()
style	one out of "chains", "acf", "density"
par2plot	logical vector, which parameters to be plotted, order: $(\phi, \theta, \gamma^2, \xi, N)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```

model <- set.to.class("NHPP", parameter = list(xi = c(5, 1/2)),
    Lambda = function(t, xi) (t/xi[2])^xi[1])
data <- simulate(model, t = seq(0, 1, by = 0.01), plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.01), data$Times, 10000) # nMCMC small for example
plot(est)
plot(est, burnIn = 1000, thinning = 20)
plot(est, reduced = FALSE, xlab = "iteration")
plot(est, style = "acf", main = "", par2plot = c(TRUE, FALSE), par.options = list(mfrow = c(1, 1)))
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), par2plot = c(FALSE, TRUE), main = "")
plot(est, priorMeans = FALSE, newwindow = TRUE)

```

---

plot,est.Regression-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.Regression

## Usage

```
## S4 method for signature 'est.Regression'
plot(x, par.options, style = c("chains", "acf",
  "density"), par2plot, reduced = TRUE, thinning, burnIn, priorMeans = TRUE,
  col.priorMean = 2, lty.priorMean = 1, newwindow = FALSE, ...)
```

## Arguments

x	est.Regression class
par.options	list of options for function par()
style	one out of "chains", "acf", "density"
par2plot	logical vector, which parameters to be plotted, order: $(\phi, \gamma^2)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
model <- set.to.class("Regression", fun = function(phi, t) phi[1]*t + phi[2],
  parameter = list(phi = c(1, 2), gamma2 = 0.1))
data <- simulate(model, t = seq(0, 1, by = 0.01), plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.01), data, 1000) # nMCMC small for example
plot(est)
plot(est, burnIn = 100, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(3,1)), xlab = "iteration")
plot(est, style = "acf", main = "", par2plot = c(TRUE, TRUE, FALSE))
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), main = "", par2plot = c(FALSE, FALSE, TRUE))
plot(est, priorMeans = FALSE, newwindow = TRUE)
```

---

plot,est.reg\_hiddenNHPP-method

*Plot method for the Bayesian estimation class object*


---

## Description

Plot method for the S4 class est.reg\_hiddenNHPP

## Usage

```
## S4 method for signature 'est.reg_hiddenNHPP'
plot(x, par.options, style = c("chains", "acf",
  "density"), par2plot, reduced = TRUE, thinning, burnIn, priorMeans = TRUE,
  col.priorMean = 2, lty.priorMean = 1, newwindow = FALSE, ...)
```

## Arguments

x	est.reg_hiddenNHPP class
par.options	list of options for function par()
style	one out of "chains", "acf", "density"
par2plot	logical vector, which parameters to be plotted, order: $(\phi, \theta, \gamma^2, \xi, N)$
reduced	logical (1), if TRUE, the chains are thinned and burn-in phase is dropped
thinning	thinning rate, if missing, the proposed one by the estimation procedure is taken
burnIn	burn-in phase, if missing, the proposed one by the estimation procedure is taken
priorMeans	logical(1), if TRUE (default), prior means are marked with a line
col.priorMean	color of the prior mean line, default 2
lty.priorMean	linetype of the prior mean line, default 1
newwindow	logical(1), if TRUE, a new window is opened for the plot
...	optional plot parameters

## Examples

```
model <- set.to.class("reg_hiddenNHPP", fun = function(t, N, theta) exp(theta[1]*t) + theta[2]*N,
  parameter = list(theta = c(2, 2), gamma2 = 0.25, xi = c(3, 0.5)),
  Lambda = function(t, xi) (t/xi[2])^xi[1])
data <- simulate(model, t = seq(0, 1, by = 0.01), plot.series = TRUE)
est <- estimate(model, t = seq(0, 1, by = 0.01), data, 1000) # nMCMC small for example
plot(est)
plot(est, burnIn = 100, thinning = 2)
plot(est, reduced = FALSE, par.options = list(mar = c(5, 4.5, 4, 2) + 0.1, mfrow = c(2, 3)), xlab = "iteration")
plot(est, style = "acf", main = "", par2plot = c(TRUE, FALSE, FALSE, TRUE, TRUE))
plot(est, style = "density", lwd = 2, priorMean = FALSE)
plot(est, style = "density", col.priorMean = 1, lty.priorMean = 2, main = "posterior")
plot(est, style = "acf", par.options = list(), par2plot = c(TRUE, rep(FALSE, 4)), main = "")
plot(est, priorMeans = FALSE, newwindow = TRUE)
```

---

plotQuantiles	<i>Plot function for credibility or prediction intervals</i>
---------------	--------------------------------------------------------------

---

**Description**

Plots intervals.

**Usage**

```
plotQuantiles(input, xlab = "", ylab = "", main = "", l = 0.025,
              u = 0.975, color = 1)
```

**Arguments**

input	list or matrix of samples from posterior or predictive distribution
xlab	a title for the x axis
ylab	a title for the y axis
main	an overall title for the plot
l	lower bound
u	upper bound
color	color of the lines to be drawn

---

postmu	<i>Posterior for mu</i>
--------	-------------------------

---

**Description**

Posterior for parameters  $\mu$

**Usage**

```
postmu(phi, m, v, Omega)
```

**Arguments**

phi	matrix of random effects
m	prior mean
v	prior variance
Omega	variance of the random effects

**Value**

one sample of posterior

---

postOmega	<i>Posterior</i>
-----------	------------------

---

**Description**

Posterior for parameters  $\Omega$

**Usage**

```
postOmega(alpha, beta, phi, mu)
```

**Arguments**

alpha	vector of prior variables
beta	vector of prior variables
phi	matrix of random effects
mu	mean of random effects

**Value**

one sample of posterior

---

postOmega_matrix	<i>Posterior</i>
------------------	------------------

---

**Description**

Posterior for parameters  $\Omega$

**Usage**

```
postOmega_matrix(R, phi, mu)
```

**Arguments**

R	prior matrix of wishart distribution
phi	matrix of random effects
mu	mean of random effects

**Value**

one sample of posterior

---

pred.base	<i>Prediction function</i>
-----------	----------------------------

---

**Description**

Bayesian prediction of the parameter of a stochastic process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t + h(\eta, t, Y_t)dN_t$ .

**Usage**

```
pred.base(samples, VFun, dens, len = 100, x0, method = c("vector", "free"),
  pred.alg = c("Distribution", "Trajectory"), sampling.alg = c("RejSamp",
    "BinSearch"), candArea, grid = 0.001)
```

**Arguments**

samples	MCMC samples
VFun	cumulative distribution function
dens	density function
len	number of samples to be drawn
x0	starting point
method	vectorial ("vector") or not ("free")
pred.alg	prediction algorithm, "Distribution" or "Trajectory"
sampling.alg	sampling algorithm, rejection sampling ("RejSamp") or inversion method ("BinSearch")
candArea	candidate area
grid	fineness degree

**Value**

vector of samples from prediction

---

predict,est.Diffusion-method	<i>Prediction for diffusion process</i>
------------------------------	-----------------------------------------

---

**Description**

Bayesian prediction of a stochastic process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ .

**Usage**

```
## S4 method for signature 'est.Diffusion'
predict(object, t, Euler.interval = FALSE,
  level = 0.05, burnIn, thinning, b.fun.mat, which.series = c("new",
    "current"), y.start, M2pred = 10, cand.length = 1000,
  pred.alg = c("Distribution", "Trajectory", "simpleTrajectory",
    "simpleBayesTrajectory"), method = c("vector", "free"),
  sampling.alg = c("BinSearch", "RejSamp"), sample.length, grid,
  plot.prediction = TRUE)
```

**Arguments**

object	class object of MCMC samples: "est.Diffusion"
t	vector of time points to make predictions for
Euler.interval	if TRUE: simple prediction intervals with Euler are made (in one step each)
level	level of the prediction intervals
burnIn	burn-in period
thinning	thinning rate
b.fun.mat	matrix-wise definition of drift function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
y.start	optional, if missing, first (which.series = "new") or last observation variable ("current") is taken
M2pred	optional, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector")
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleBayesTrajectory"
method	vectorial ("vector") or not ("free")
sampling.alg	sampling algorithm, inversion method ("BinSearch") or rejection sampling ("RejSamp")
sample.length	length of samples to be drawn
grid	fineness degree of approximation
plot.prediction	if TRUE, result are plotted

**Examples**

```

model <- set.to.class("Diffusion", parameter = list(phi = 0.5, gamma2 = 0.01))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5, plot.series = TRUE)
est_diff <- estimate(model, t, data, 1000) # better: 10000
plot(est_diff)
## Not run:
pred_diff <- predict(est_diff, t = seq(0, 1, by = 0.1))
pred_diff <- predict(est_diff, b.fun.mat = function(phi, t, y) phi[,1]) # much faster
pred_diff2 <- predict(est_diff, which.series = "current", b.fun.mat = function(phi, t, y) phi[,1])
pred_diff3 <- predict(est_diff, which.series = "current", y.start = data[51],
                     t = t[seq(51, 100, by = 5)], b.fun.mat = function(phi, t, y) phi[,1])

## End(Not run)
pred_diff <- predict(est_diff, Euler.interval = TRUE, b.fun.mat = function(phi, t, y) phi[,1])
# one step Euler approximation
pred_diff <- predict(est_diff, pred.alg = "simpleTrajectory", sample.length = 100)
for(i in 1:100) lines(t[-1], pred_diff[i,], col = "grey")
pred_diff <- predict(est_diff, pred.alg = "simpleBayesTrajectory")

```

---

predict,est.hiddenDiffusion-method

*Prediction for noisy / hidden diffusion process*


---

## Description

Bayesian prediction of the model,  $Z_i = Y_{t_i} + \epsilon_i$ ,  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ .

## Usage

```
## S4 method for signature 'est.hiddenDiffusion'
predict(object, t, burnIn, thinning, b.fun.mat,
  which.series = c("new", "current"), M2pred = 10, cand.length = 1000,
  pred.alg = c("Distribution", "Trajectory", "simpleTrajectory",
    "simpleBayesTrajectory"), sample.length, grid, plot.prediction = TRUE)
```

## Arguments

object	class object of MCMC samples: "est.hiddenDiffusion"
t	vector of time points to make predictions for
burnIn	burn-in period
thinning	thinning rate
b.fun.mat	matrix-wise definition of drift function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
M2pred	optional, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector")
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleBayesTrajectory"
sample.length	length of samples to be drawn
grid	fineness degree of approximation
plot.prediction	if TRUE, result are plotted

## Examples

```
model <- set.to.class("hiddenDiffusion", parameter = list(phi = 5, gamma2 = 1, sigma2 = 0.1))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est_hiddiff <- estimate(model, t, data$Z, 100) # nMCMC should be much larger!
plot(est_hiddiff)
## Not run:
pred_hiddiff <- predict(est_hiddiff, t = seq(0, 1, by = 0.1))
pred_hiddiff2 <- predict(est_hiddiff, which.series = "current")

## End(Not run)
pred_hiddiff <- predict(est_hiddiff, pred.alg = "simpleTrajectory", sample.length = 100)
pred_hiddiff <- predict(est_hiddiff, pred.alg = "simpleBayesTrajectory")
```



---

predict,est.hiddenmixedDiffusion-method

*Prediction for noisy/hidden mixed diffusion process*


---

## Description

Bayesian prediction of a stochastic process  $Z_{ij} = Y_{t_{ij}} + \epsilon_{ij}$ ,  $dY_t = b(\phi_j, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ ,  $\phi_j \sim N(\mu, \Omega)$ .

## Usage

```
## S4 method for signature 'est.hiddenmixedDiffusion'
predict(object, t, burnIn, thinning,
        b.fun.mat, which.series = c("new", "current"), ind.pred, M2pred = 10,
        cand.length = 1000, pred.alg = c("Distribution", "Trajectory",
        "simpleTrajectory", "simpleBayesTrajectory"), sample.length, grid,
        plot.prediction = TRUE)
```

## Arguments

object	class object of MCMC samples: "est.hiddenmixedDiffusion"
t	vector of time points to make predictions for
burnIn	burn-in period
thinning	thinning rate
b.fun.mat	matrix-wise definition of drift function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
ind.pred	index of series to be predicted, optional, if which.series = "current" and ind.pred missing, the last series is taken
M2pred	optional, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector")
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleBayesTrajectory"
sample.length	length of samples to be drawn
grid	fineness degree of approximation
plot.prediction	if TRUE, result are plotted

## Examples

```
mu <- c(5, 1); Omega <- c(0.9, 0.04)
phi <- cbind(rnorm(21, mu[1], sqrt(Omega[1])), rnorm(21, mu[2], sqrt(Omega[2])))
y0.fun <- function(phi, t) phi[2]
model <- set.to.class("hiddenmixedDiffusion", y0.fun = y0.fun,
                     b.fun = function(phi, t, y) phi[1],
                     parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 1, sigma2 = 0.01))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
```

```
## Not run:
est_hidmixdiff <- estimate(model, t, data$Z[1:20,], 2000)
plot(est_hidmixdiff)
pred1 <- predict(est_hidmixdiff, b.fun.mat = function(phi, t, y) phi[,1])
pred2 <- predict(est_hidmixdiff, pred.alg = "Trajectory", b.fun.mat = function(phi, t, y) phi[,1])
pred3 <- predict(est_hidmixdiff, pred.alg = "simpleTrajectory", sample.length = nrow(pred1$Y))
lines(t, apply(pred1$Z, 2, quantile, 0.025), col = 3)
lines(t, apply(pred1$Z, 2, quantile, 0.975), col = 3)
lines(t, apply(pred2$Z, 2, quantile, 0.025), col = 4)
lines(t, apply(pred2$Z, 2, quantile, 0.975), col = 4)
pred4 <- predict(est_hidmixdiff, pred.alg = "simpleBayesTrajectory")

## End(Not run)
```

---

predict,est.jumpDiffusion-method

*Prediction for jump diffusion process*

---

## Description

Bayesian prediction of a stochastic process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t + h(\eta, t, Y_t)dN_t$ .

## Usage

```
## S4 method for signature 'est.jumpDiffusion'
predict(object, t, burnIn, thinning, Lambda.mat,
  which.series = c("new", "current"), M2pred = 10, cand.length = 1000,
  pred.alg = c("Trajectory", "Distribution", "simpleTrajectory",
    "simpleBayesTrajectory"), pred.alg.N = c("Trajectory", "Distribution"),
  candN = 0:5, sample.length, plot.prediction = TRUE)
```

## Arguments

object	class object of MCMC samples: "est.jumpDiffusion"
t	vector of time points to make predictions for
burnIn	burn-in period
thinning	thinning rate
Lambda.mat	matrix-wise definition of intensity rate function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
M2pred	default 10, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector"), for jump diffusion
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleBayesTrajectory"
pred.alg.N	prediction algorithm, "Distribution", "Trajectory"
candN	vector of candidate area for differences of N, only if pred.alg.N = "Distribution"
sample.length	length of samples to be drawn
plot.prediction	if TRUE, result are plotted

## Examples

```

model <- set.to.class("jumpDiffusion",
  parameter = list(theta = 0.1, phi = 0.05, gamma2 = 0.1, xi = c(3, 1/4)),
  Lambda = function(t, xi) (t/xi[2])^xi[1])
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5, plot.series = TRUE)
est_jd <- estimate(model, t, data, 2000)
plot(est_jd)
## Not run:
pred_jd <- predict(est_jd, Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1])
pred_jd2 <- predict(est_jd, pred.alg = "Distribution", pred.alg.N = "Distribution", Lambda.mat = function(t,
est <- estimate(model, t[1:81], data = list(N = data$N[1:81], Y = data$Y[1:81]), 2000)
pred <- predict(est, t = t[81:101], which.series = "current",
  Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1])
lines(t, data$Y, type="l", lwd = 2)

## End(Not run)
pred_jd4 <- predict(est_jd, pred.alg = "simpleTrajectory", sample.length = 100)
for(i in 1:100) lines(t[-1], pred_jd4$Y[i,], col = "grey")

```

---

predict.est.Merton-method

*Prediction for jump diffusion process*

---

## Description

Bayesian prediction of a stochastic process  $Y_t = y_0 \exp(\phi t - \gamma^2/2t + \gamma W_t + \log(1 + \theta)N_t)$ .

## Usage

```

## S4 method for signature 'est.Merton'
predict(object, t, burnIn, thinning, Lambda.mat,
  which.series = c("new", "current"), M2pred = 10, only.interval = TRUE,
  level = 0.05, cand.length = 1000, pred.alg = c("Distribution",
  "Trajectory", "simpleTrajectory", "simpleBayesTrajectory"), sample.length,
  plot.prediction = TRUE)

```

## Arguments

object	class object of MCMC samples: "est.Merton"
t	vector of time points to make predictions for
burnIn	burn-in period
thinning	thinning rate
Lambda.mat	matrix-wise definition of intensity rate function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
M2pred	default 10, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
only.interval	if TRUE: only calculation of prediction intervals (only for pred.alg = "Distribution")

level	level of the prediction intervals
cand.length	length of candidate samples (if method = "vector"), for jump diffusion
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleBayesTrajectory"
sample.length	length of samples to be drawn
plot.prediction	if TRUE, result are plotted

### Examples

```

cl <- set.to.class("Merton",
  parameter = list(thetaT = 0.1, phi = 0.05, gamma2 = 0.1, xi = c(3, 1/4)),
  Lambda = function(t, xi) (t/xi[2])^xi[1])
t <- seq(0, 1, by = 0.01)
data <- simulate(cl, t = t, y0 = 0.5, plot.series = TRUE)
est <- estimate(cl, t, data, 1000)
plot(est)
## Not run:
pred1 <- predict(est, Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1])
pred2 <- predict(est, Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1], pred.alg = "Trajectory")
pred3 <- predict(est, pred.alg = "simpleTrajectory")
pred4 <- predict(est, pred.alg = "simpleBayesTrajectory")

## End(Not run)

```

---

predict,est.mixedDiffusion-method

*Prediction for mixed diffusion process*

---

### Description

Bayesian prediction of a stochastic process  $dY_t = b(\phi_j, t, Y_t)dt + s(\gamma, t, Y_t)dW_t, \phi_j \sim N(\mu, \Omega)$ .

### Usage

```

## S4 method for signature 'est.mixedDiffusion'
predict(object, t, Euler.interval = FALSE,
  level = 0.05, burnIn, thinning, b.fun.mat, which.series = c("new",
  "current"), y.start, ind.pred, M2pred = 10, cand.length = 1000,
  pred.alg = c("Distribution", "Trajectory", "simpleTrajectory",
  "simpleBayesTrajectory"), sample.length, grid, plot.prediction = TRUE)

```

### Arguments

object	class object of MCMC samples: "est.mixedDiffusion"
t	vector of time points to make predictions for
Euler.interval	if TRUE: simple prediction intervals with Euler are made (in one step each)
level	level of the prediction intervals
burnIn	burn-in period
thinning	thinning rate

b.fun.mat	matrix-wise definition of drift function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
y.start	optional, if missing, first (which.series = "new") or last observation variable ("current") is taken
ind.pred	index of series to be predicted, optional, if which.series = "current" and ind.pred missing, the last series is taken
M2pred	optional, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector")
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleBayesTrajectory"
sample.length	length of samples to be drawn
grid	fineness degree of approximation
plot.prediction	if TRUE, result are plotted

### Examples

```

mu <- 2; Omega <- 0.4; phi <- matrix(rnorm(21, mu, sqrt(Omega)))
model <- set.to.class("mixedDiffusion",
  parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1),
  b.fun = function(phi, t, x) phi*x, sT.fun = function(t, x) x)
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est_mixdiff <- estimate(model, t, data[1:20,], 100) # nMCMC should be much larger
plot(est_mixdiff)
## Not run:
pred_mixdiff <- predict(est_mixdiff, b.fun.mat = function(phi, t, y) phi[,1]*y)
lines(t, data[21,], lwd = 2)
mean(apply(pred_mixdiff$Y, 2, quantile, 0.025) <= data[21, ] &
  apply(pred_mixdiff$Y, 2, quantile, 0.975) >= data[21, ])
mean(sapply(1:20, function(i){
  mean(apply(pred_mixdiff$Y, 2, quantile, 0.025) <= data[i, ] &
    apply(pred_mixdiff$Y, 2, quantile, 0.975) >= data[i, ]))})
pred_mixdiff2 <- predict(est_mixdiff, b.fun.mat = function(phi, t, y) phi[,1]*y,
  which.series = "current")
pred_mixdiff3 <- predict(est_mixdiff, b.fun.mat = function(phi, t, y) phi[,1]*y,
  which.series = "current", y.start = data[20, 51], t = t[51:101])

## End(Not run)
pred_mixdiff <- predict(est_mixdiff, Euler.interval = TRUE,
  b.fun.mat = function(phi, t, y) phi[,1]*y); lines(t, data[21,], lwd = 2)
# one step Euler approximation
pred_mixdiff <- predict(est_mixdiff, pred.alg = "simpleTrajectory", sample.length = 100)
for(i in 1:100) lines(t, pred_mixdiff$Y[i,], col = "grey")
pred_mixdiff <- predict(est_mixdiff, pred.alg = "simpleBayesTrajectory")

# OU
## Not run:
b.fun <- function(phi, t, y) phi[1]-phi[2]*y; y0.fun <- function(phi, t) phi[3]
mu <- c(10, 1, 0.5); Omega <- c(0.9, 0.01, 0.01)
phi <- sapply(1:3, function(i) rnorm(21, mu[i], sqrt(Omega[i])))

```

```

cl <- set.to.class("mixedDiffusion",
  parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1),
  y0.fun = y0.fun, b.fun = b.fun, sT.fun = function(t, x) 1)
t <- seq(0, 1, by = 0.01)
data <- simulate(cl, t = t, plot.series = TRUE)
est <- estimate(cl, t, data[1:20,], 2000)
plot(est)
pred_mixdiff <- predict(est, t = seq(0, 1, length = 21), b.fun.mat = function(phi, t, y) phi[,1]-phi[,2]*y)
lines(t, data[21,], lwd = 2)
mean(apply(pred_mixdiff$Y, 2, quantile, 0.025) <= data[21, seq(1, length(t), length = 21)] &
  apply(pred_mixdiff$Y, 2, quantile, 0.975) >= data[21, seq(1, length(t), length = 21)]) &
  mean(sapply(1:20, function(i){
    mean(apply(pred_mixdiff$Y, 2, quantile, 0.025) <= data[i, seq(1, length(t), length = 21)] &
      apply(pred_mixdiff$Y, 2, quantile, 0.975) >= data[i, seq(1, length(t), length = 21)]))}))

## End(Not run)

```

---

predict,est.mixedRegression-method

*Prediction for mixed regression model*

---

## Description

Bayesian prediction of the regression model  $y_i = f(\phi_j, t_i) + \epsilon_i, \phi_j \sim N(\mu, \Omega)$ .

## Usage

```

## S4 method for signature 'est.mixedRegression'
predict(object, t, only.interval = TRUE,
  level = 0.05, burnIn, thinning, fun.mat, which.series = c("new",
    "current"), ind.pred, M2pred = 10, cand.length = 1000, sample.length,
  grid, plot.prediction = TRUE)

```

## Arguments

object	class object of MCMC samples: "est.mixedRegression"
t	vector of time points to make predictions for
only.interval	if TRUE: only calculation of prediction intervals
level	level of the prediction intervals
burnIn	burn-in period
thinning	thinning rate
fun.mat	matrix-wise definition of drift function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
ind.pred	index of series to be predicted, optional, if which.series = "current" and ind.pred missing, the last series is taken
M2pred	optional, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector")

sample.length    length of samples to be drawn  
 grid             fineness degree of approximation  
 plot.prediction    if TRUE, result are plotted

### Examples

```
mu <- c(10, 5); Omega <- c(0.9, 0.01)
phi <- cbind(rnorm(21, mu[1], sqrt(Omega[1])), rnorm(21, mu[2], sqrt(Omega[2])))
cl <- set.to.class("mixedRegression",
  parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1),
  fun = function(phi, t) phi[1]*t + phi[2], sT.fun = function(t) 1)
t <- seq(0, 1, by = 0.01)
data <- simulate(cl, t = t, plot.series = TRUE)
est <- estimate(cl, t, data[1:20,], 2000)
plot(est)
pred <- predict(est, fun.mat = function(phi, t) phi[,1]*t + phi[,2])
points(t, data[21,], pch = 20)
```

---

predict.est.NHPP-method

*Prediction for Poisson process*

---

### Description

Bayesian prediction of a nonhomogeneous Poisson process.

### Usage

```
## S4 method for signature 'est.NHPP'
predict(object, variable = c("eventTimes",
  "PoissonProcess"), t, burnIn, thinning, Lambda.mat, which.series = c("new",
  "current"), Tstart, M2pred = 10, rangeN = c(0, 5), cand.length = 1000,
  pred.alg = c("Trajectory", "Distribution", "simpleTrajectory",
  "simpleBayesTrajectory"), sample.length, grid = 1e-05,
  plot.prediction = TRUE)
```

### Arguments

object	class object of MCMC samples: "est.NHPP"
variable	if prediction of event times ("eventTimes") or of Poisson process variables ("PoissonProcess")
t	vector of time points to make predictions for (only for variable = "PoissonProcess")
burnIn	burn-in period
thinning	thinning rate
Lambda.mat	matrix-wise definition of drift function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")

Tstart	optional, if missing, first (which.series = "new") or last observation variable ("current") is taken
M2pred	optional, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
rangeN	vector of candidate area for differences of N, only if pred.alg = "Distribution" and variable = "PoissonProcess"
cand.length	length of candidate samples (if method = "vector")
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleBayesTrajectory"
sample.length	length of samples to be drawn
grid	fineness degress of approximation
plot.prediction	if TRUE, result are plotted

### Examples

```

model <- set.to.class("NHPP", parameter = list(xi = c(5, 1/2)),
                    Lambda = function(t, xi) (t/xi[2])^xi[1])
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
est_NHPP <- estimate(model, t, data$Times, 1000) # nMCMC should be much larger!
plot(est_NHPP)
pred <- predict(est_NHPP, Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1], variable = "PoissonProcess", pred.a
## Not run:
pred_NHPP <- predict(est_NHPP, Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1])
pred_NHPP <- predict(est_NHPP, variable = "PoissonProcess", Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1])
pred_NHPP2 <- predict(est_NHPP, which.series = "current", Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1])
pred_NHPP3 <- predict(est_NHPP, variable = "PoissonProcess", which.series = "current", Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1])
pred_NHPP4 <- predict(est_NHPP, pred.alg = "simpleTrajectory", M2pred = length(data$Times))

## End(Not run)
pred_NHPP <- predict(est_NHPP, variable = "PoissonProcess", pred.alg = "simpleTrajectory",
                    M2pred = length(data$Times))
pred_NHPP <- predict(est_NHPP, variable = "PoissonProcess", pred.alg = "simpleBayesTrajectory",
                    M2pred = length(data$Times), sample.length = 100)

```

---

predict,est.Regression-method

*Prediction for regression model*

---

### Description

Bayesian prediction of regression model  $y_i = f(\phi, t_i) + \epsilon_i$ .

### Usage

```

## S4 method for signature 'est.Regression'
predict(object, t, only.interval = TRUE,
        level = 0.05, burnIn, thinning, fun.mat, which.series = c("new",
        "current"), M2pred = 10, cand.length = 1000, method = c("vector",
        "free"), sampling.alg = c("BinSearch", "RejSamp"), sample.length, grid,
        plot.prediction = TRUE)

```



**Arguments**

object	class object of MCMC samples: "est.Regression"
t	vector of time points to make predictions for
only.interval	if TRUE: only calculation of prediction intervals
level	level of the prediction intervals
burnIn	burn-in period
thinning	thinning rate
fun.mat	matrix-wise definition of drift function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
M2pred	optional, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector")
method	vectorial ("vector") or not ("free")
sampling.alg	sampling algorithm, inversion method ("BinSearch") or rejection sampling ("RejSamp")
sample.length	length of samples to be drawn
grid	fineness degree of approximation
plot.prediction	if TRUE, result are plotted

**Examples**

```

t <- seq(0,1, by = 0.01)
cl <- set.to.class("Regression", fun = function(phi, t) phi[1]*t + phi[2],
                  parameter = list(phi = c(1,2), gamma2 = 0.1))
data <- simulate(cl, t = t, plot.series = TRUE)
est <- estimate(cl, t, data, 1000)
plot(est)
pred <- predict(est, fun.mat = function(phi, t) phi[,1]*t + phi[,2])
## Not run:
pred2 <- predict(est, fun.mat = function(phi, t) phi[,1]*t + phi[,2], only.interval = FALSE)
plot(density(pred2[,10]))

## End(Not run)

```

---

predict,est.reg\_hiddenNHPP-method

*Prediction for regression model dependent on Poisson process*

---

**Description**

Bayesian prediction of a regression model  $y_i = f(t_i, N_i, \theta) + \epsilon_i$ .

**Usage**

```
## S4 method for signature 'est.reg_hiddenNHPP'
predict(object, t, only.interval = TRUE,
  level = 0.05, burnIn, thinning, Lambda.mat, fun.mat,
  which.series = c("new", "current"), M2pred = 10, cand.length = 1000,
  pred.alg = c("Distribution", "simpleTrajectory", "simpleBayesTrajectory"),
  sample.length, grid = 1e-05, plot.prediction = TRUE)
```

**Arguments**

object	class object of MCMC samples: "est.reg_hiddenNHPP"
t	vector of time points to make predictions for
only.interval	if TRUE: only calculation of prediction intervals
level	level of the prediction intervals
burnIn	burn-in period
thinning	thinning rate
Lambda.mat	matrix-wise definition of intensity rate function (makes it faster)
fun.mat	matrix-wise definition of regression function (makes it faster)
which.series	which series to be predicted, new one ("new") or further development of current one ("current")
M2pred	default 10, if current series to predicted and t missing, M2pred variables will be predicted with dt of observation time points
cand.length	length of candidate samples (if method = "vector"), for jump diffusion
pred.alg	prediction algorithm, "Distribution", "Trajectory", "simpleTrajectory" or "simpleTrajectory"
sample.length	length of samples to be drawn
grid	fineness degress of approximation, for Poisson process
plot.prediction	if TRUE, result are plotted

**Examples**

```
t <- seq(0,1, by = 0.01)
cl <- set.to.class("reg_hiddenNHPP", fun = function(t, N, theta) theta[1]*t + theta[2]*N,
  parameter = list(theta = c(1,2), gamma2 = 0.1, xi = c(3, 1/4)),
  Lambda = function(t, xi) (t/xi[2])^xi[1])
data <- simulate(cl, t = t, plot.series = TRUE)
est <- estimate(cl, t, data, 1000)
plot(est)
## Not run:
pred <- predict(est, Lambda.mat = function(t, xi) (t/xi[,2])^xi[,1],
  fun.mat = function(t, N, theta) theta[,1]*t + theta[,2]*N)

## End(Not run)
pred <- predict(est, pred.alg = "simpleTrajectory", sample.length = 100)
```

---

prediction.intervals    *Prediction interval function*

---

### Description

Bayesian prediction of the parameter of a stochastic process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t + h(\eta, t, Y_t)dN_t$ .

### Usage

```
prediction.intervals(samples, Fun, x0, level = 0.05, candArea, grid = 0.001)
```

### Arguments

samples	MCMC samples
Fun	cumulative distribution function
x0	starting point
level	level of prediction intervals
candArea	candidate area
grid	fineness degree

### Value

prediction interval

---

predPhi	<i>Prediction Of The Random Effects In Mixed Stochastic Differential Equations</i>
---------	------------------------------------------------------------------------------------

---

### Description

Prediction of the random effects  $\phi \sim N(\mu, \Omega)$

### Usage

```
predPhi(samples, cand)
```

### Arguments

samples	output of estSDE or estREg
cand	candidates for phi (matrix with p columns)

### Value

matrix phi

---

predReg	<i>Bayesian prediction in mixed nonlinear regression models</i>
---------	-----------------------------------------------------------------

---

**Description**

Bayesian prediction in the mixed nonlinear regression model  $y_{ij} = f(\phi_j, t_{ij}) + \epsilon_{ij}, \epsilon_{ij} \sim N(0, \gamma^2 * s^2(t_{ij}), \phi_j) \sim N(\mu, \Omega)$ .

**Usage**

```
predReg(t, samples, fODE, sVar, cand, len = 1000, mod = "Gompertz")
```

**Arguments**

t	vector of times which are predicted
samples	list of samples from the posterior
fODE	regression function
sVar	variance function
cand	vector of candidates for trajectory sampling
len	number of samples from the predictive distribution
mod	model out of Gompertz, Richards, logistic, Weibull, only used instead of fODE

**Value**

matrix of predictions in t

---

proposal	<i>Sampling from proposal density for strictly positive parameters</i>
----------	------------------------------------------------------------------------

---

**Description**

Used in Metropolis Hastings algorithms.

**Usage**

```
proposal(parOld, propSd)
```

**Arguments**

parOld	the parameter from the last iteration step
propSd	proposal standard deviation

**Value**

candidate for MH ratio

**Examples**

```
plot(replicate(100, proposal(1, 0.1)), type = "l")
```

---

proposalRatio	<i>Sampling from proposal density for strictly positive parameters</i>
---------------	------------------------------------------------------------------------

---

**Description**

Used in Metropolis Hastings algorithms.

**Usage**

```
proposalRatio(parOld, parNew, propSd)
```

**Arguments**

parOld	the parameter from the last iteration step
parNew	drawn candidate
propSd	proposal standard deviation

**Value**

MH ratio

**Examples**

```
cand <- proposal(1, 0.01)
proposalRatio(1, cand, 0.01)
```

---

RejSampling	<i>Rejection Sampling Algorithm</i>
-------------	-------------------------------------

---

**Description**

Rejection Sampling

**Usage**

```
RejSampling(Fun, dens, len, cand, grid = 0.001, method = c("vector",
"free"))
```

**Arguments**

Fun	cumulative distribution function
dens	density
len	number of samples
cand	candidate area
grid	fineness degree
method	vectorial ("vector") or not ("free")

**Value**

vector of samples

**Examples**

```
plot(density(RejSampling(Fun = function(x) pnorm(x, 5, 1), dens = function(x) dnorm(x, 5, 1), len = 500, cand = seq(2, 9, by = 1)),
lines(density(RejSampling(function(x) pnorm(x, 5, 1), function(x) dnorm(x, 5, 1), 500, cand = seq(2, 9, by = 1)),
curve(dnorm(x, 5, 1), from = 2, to = 8, add = TRUE, col = 3)
```

---

scoreRule	<i>Calculation of interval score</i>
-----------	--------------------------------------

---

**Description**

Scoring rule of Raftery and Gneiting (??).

**Usage**

```
scoreRule(l, u, x, alpha = 0.05)
```

**Arguments**

l	lower bound
u	upper bound
x	true value
alpha	level

**Value**

interval score

---

set.to.class	<i>Builds classes</i>
--------------	-----------------------

---

**Description**

Defines classes

**Usage**

```
set.to.class(class.name = c("jumpDiffusion", "Merton", "Diffusion",
"mixedDiffusion", "hiddenDiffusion", "hiddenmixedDiffusion", "reg_hiddenNHPP",
"NHPP", "Regression", "mixedRegression"), parameter, prior, start, b.fun,
s.fun, h.fun, sT.fun, y0.fun, fun, Lambda, priorRatio)
```

**Arguments**

class.name	name of model class
parameter	list of parameter values
prior	optional list of prior parameters
start	optional list of starting values
b.fun	drift function b
s.fun	variance function s
h.fun	jump high function h
sT.fun	variance function $\tilde{s}$
y0.fun	function for the starting point, if dependent on parameter
fun	regression function
Lambda	intensity rate of Poisson process
priorRatio	list of functions for prior ratios, only for jumpDiffusion, is missing: non-informative estimation

**Value**

class

**Examples**

```
set.to.class("jumpDiffusion")
cl_jd <- set.to.class("jumpDiffusion", parameter = list(theta = 0.1, phi = 0.01, gamma2 = 0.1, xi = 3))
summary(class.to.list(cl_jd))
```

---

simN

*Simulation of counting process*


---

**Description**

Simulation of counting process and event times.

**Usage**

```
simN(t, xi, len, start = c(0, 0), Lambda, int = c("Weibull", "Exp"))
```

**Arguments**

t	vector of times
xi	parameter vector $\xi$
len	number of samples to be drawn
start	vector: start[1] starting point time, start[2] starting point for Poisson process
Lambda	intensity rate function
int	if no Lambda: one out of "Weibull" or "Exp" for intensity function

**Value**

N	Poisson process
Times	event times

---

simulate,Diffusion-method

*Simulation of diffusion process*

---

### Description

Simulation of a stochastic process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ .

### Usage

```
## S4 method for signature 'Diffusion'
simulate(object, nsim = 1, seed = NULL, t, y0,
         mw = 1, plot.series = TRUE)
```

### Arguments

object	class object of parameters: "Diffusion"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
y0	starting point of the process
mw	mesh width for finer Euler approximation
plot.series	logical(1), if TRUE, simulated series are depicted grafically

### Examples

```
model <- set.to.class("Diffusion", parameter = list(phi = 0.5, gamma2 = 0.01))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5, plot.series = TRUE)
```

---

simulate,hiddenDiffusion-method

*Simulation of diffusion process*

---

### Description

Simulation of a hidden stochastic process  $Z_i = Y_{t_i} + \epsilon_i, dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ .

### Usage

```
## S4 method for signature 'hiddenDiffusion'
simulate(object, nsim = 1, seed = NULL, t,
         mw = 10, plot.series = TRUE)
```



**Arguments**

object	class object of parameters: "hiddenDiffusion"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
mw	mesh width for finer Euler approximation
plot.series	logical(1), if TRUE, simulated series are depicted grafically

**Examples**

```
model <- set.to.class("hiddenDiffusion", parameter = list(phi = 0.5, gamma2 = 0.01, sigma2 = 0.1))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
```

---

```
simulate,hiddenmixedDiffusion-method
```

*Simulation of hidden mixed diffusion process*

---

**Description**

Simulation of a stochastic process  $Z_{ij} = Y_{t_{ij}} + \epsilon_{ij}$ ,  $dY_t = b(\phi_j, t, Y_t)dt + s(\gamma, t, Y_t)dW_t$ ,  $\phi_j \sim N(\mu, \Omega)$ .

**Usage**

```
## S4 method for signature 'hiddenmixedDiffusion'
simulate(object, nsim = 1, seed = NULL, t,
         mw = 10, plot.series = TRUE)
```

**Arguments**

object	class object of parameters: "hiddenmixedDiffusion"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
mw	mesh width for finer Euler approximation
plot.series	logical(1), if TRUE, simulated series are depicted grafically

**Examples**

```
mu <- c(5, 1); Omega <- c(0.9, 0.04); phi <- cbind(rnorm(21, mu[1], sqrt(Omega[1])), rnorm(21, mu[2], sqrt(Omega[2])))
y0.fun <- function(phi, t) phi[2]
model <- set.to.class("hiddenmixedDiffusion", y0.fun = y0.fun, b.fun = function(phi, t, y) phi[1], parameter = list(mu = mu, Omega = Omega))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t)
```

---

simulate,jumpDiffusion-method  
*Simulation of jump diffusion process*

---

### Description

Simulation of jump diffusion process  $dY_t = b(\phi, t, Y_t)dt + s(\gamma, t, Y_t)dW_t + h(\eta, t, Y_t)dN_t$ .

### Usage

```
## S4 method for signature 'jumpDiffusion'
simulate(object, nsim = 1, seed = NULL, t, y0,
         mw = 1, plot.series = TRUE)
```

### Arguments

object	class object of parameters: "jumpDiffusion"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
y0	starting point of process
mw	mesh width for finer Euler approximation
plot.series	logical(1), if TRUE, simulated series are depicted grafically

### Examples

```
model <- set.to.class("jumpDiffusion", parameter = list(theta = 0.1, phi = 0.05, gamma2 = 0.1, xi = c(3, 1/4)))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5)
```

---

simulate,Merton-method  
*Simulation of jump diffusion process*

---

### Description

Simulation of jump diffusion process  $Y_t = y_0 \exp(\phi t - \gamma^2/2t + \gamma W_t + \log(1 + \theta)N_t)$ .

### Usage

```
## S4 method for signature 'Merton'
simulate(object, nsim = 1, seed = NULL, t, y0,
         plot.series = TRUE)
```

**Arguments**

object	class object of parameters: "Merton"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
y0	starting point of process
plot.series	logical(1), if TRUE, simulated series are depicted grafically

**Examples**

```
model <- set.to.class("Merton", parameter = list(thetaT = 0.1, phi = 0.05, gamma2 = 0.1, xi = 10))
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, y0 = 0.5)
```

---

simulate,mixedDiffusion-method

*Simulation of diffusion process*

---

**Description**

Simulation of a stochastic process  $dY_t = b(\phi_j, t, Y_t)dt + s(\gamma, t, Y_t)dW_t, \phi_j \sim N(\mu, \Omega)$ .

**Usage**

```
## S4 method for signature 'mixedDiffusion'
simulate(object, nsim = 1, seed = NULL, t,
         mw = 1, plot.series = TRUE)
```

**Arguments**

object	class object of parameters: "mixedDiffusion"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
mw	mesh width for finer Euler approximation
plot.series	logical(1), if TRUE, simulated series are depicted grafically

**Examples**

```
mu <- 2; Omega <- 0.4; phi <- matrix(rnorm(21, mu, sqrt(Omega)))
model <- set.to.class("mixedDiffusion", parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1), y0.
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
```

---

 simulate,mixedRegression-method

*Simulation of mixed regression model*


---

### Description

Simulation of regression model  $y_i = f(\phi_j, t_i) + \epsilon_i, \phi_j \sim N(\mu, \Omega)$ .

### Usage

```
## S4 method for signature 'mixedRegression'
simulate(object, nsim = 1, seed = NULL, t,
  plot.series = TRUE)
```

### Arguments

object	class object of parameters: "mixedRegression"
nsim	number of response vectors to simulate. Defaults = 1.
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
plot.series	logical(1), if TRUE, simulated series are depicted grafically

### Examples

```
mu <- 2; Omega <- 0.4; phi <- matrix(rnorm(21, mu, sqrt(Omega)))
model <- set.to.class("mixedRegression", parameter = list(phi = phi, mu = mu, Omega = Omega, gamma2 = 0.1), fu
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
```

---

 simulate,NHPP-method    *Simulation of Poisson process*


---

### Description

Simulation of Poisson process.

### Usage

```
## S4 method for signature 'NHPP'
simulate(object, nsim = 1, seed = NULL, t,
  plot.series = TRUE)
```

### Arguments

object	class object of parameters: "NHPP"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
plot.series	logical(1), if TRUE, simulated series are depicted grafically

**Examples**

```
model <- set.to.class("NHPP", parameter = list(xi = c(5, 1/2)), Lambda = function(t, xi) (t/xi[2])^xi[1])
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t)
```

---

simulate,Regression-method

*Simulation of regression model*

---

**Description**

Simulation of the regression model  $y_i = f(\phi, t_i) + \epsilon_i$ .

**Usage**

```
## S4 method for signature 'Regression'
simulate(object, nsim = 1, seed = NULL, t,
         plot.series = TRUE)
```

**Arguments**

object	class object of parameters: "Diffusion"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
plot.series	logical(1), if TRUE, simulated series are depicted grafically

**Examples**

```
model <- set.to.class("Regression", parameter = list(phi = 5, gamma2 = 0.1), fun = function(phi, t) phi*t)
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t, plot.series = TRUE)
```

---

simulate,reg\_hiddenNHPP-method

*Simulation of regression model dependent on Poisson process*

---

**Description**

Simulation of of the regression model  $y_i = f(t_i, N_i, \theta) + \epsilon_i$ .

**Usage**

```
## S4 method for signature 'reg_hiddenNHPP'
simulate(object, nsim = 1, seed = NULL, t,
         plot.series = TRUE)
```

Arguments

object	class object of parameters: "reg_hiddenNHPP"
nsim	number of response vectors to simulate. Defaults to 1
seed	optional: seed number for random number generator
t	vector of time points to make predictions for
plot.series	logical(1), if TRUE, simulated series are depicted grafically

Examples

```
model <- set.to.class("reg_hiddenNHPP", fun = function(t, N, theta) theta[1]*t + theta[2]*N, parameter = list(
t <- seq(0, 1, by = 0.01)
data <- simulate(model, t = t)
```

---

simY	<i>Simulation of Jump diffusion process</i>
------	---------------------------------------------

---

Description

Simulation of Jump diffusion process.

Usage

```
simY(t, phi, thetaT, gamma2, start, N)
```

Arguments

t	vector of times
phi	parameter $\phi$
thetaT	parameter $\tilde{\theta}$
gamma2	parameter $\gamma^2$
start	starting point of process $y_0$
N	Poisson process variables in t

Value

matrix or vector

---

sim_JD_Euler	<i>Simulation of Jump diffusion process</i>
--------------	---------------------------------------------

---

**Description**

Simulation of Jump diffusion process.

**Usage**

```
sim_JD_Euler(t, phi, theta, gamma2, b.fun, s.fun, h.fun, start, N, mw = 1)
```

**Arguments**

t	vector of times
phi	parameter $\phi$
theta	parameter $\theta$
gamma2	parameter $\gamma^2$
b.fun	drift function
s.fun	variance function
h.fun	jump high function
start	starting point $y_0$
N	Poisson process variables in t
mw	mesh width for finer Euler approximation

**Value**

matrix or vector

---

sim_reg_hiddenNHPP	<i>Simulation of regression model including the NHPP</i>
--------------------	----------------------------------------------------------

---

**Description**

Simulation.

**Usage**

```
sim_reg_hiddenNHPP(t, N, fun, theta, gamma2)
```

**Arguments**

t	vector of times
N	vector of Poisson process
fun	regression function
theta	parameter $\theta$
gamma2	parameter $\gamma^2$

**Value**

matrix or vector

---

TimestoN*Transformation of vector of event times to counting process*

---

**Description**

Transformation of vector of event times to counting process.

**Usage**

```
TimestoN(times, t)
```

**Arguments**

times	vector of event times
t	times of counting process

**Value**

vector of counting process observations in t

**Examples**

```
t <- seq(0, 1, by = 0.01)
times <- simN(t, c(5, 0.5), len = 1)$Times
process <- TimestoN(times, t)
```



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