

# Package ‘fitR’

June 14, 2016

**Title** Tool box for fitting dynamic infectious disease models to time series.

**Version** 0.1

**Description** This package contains helper functions for model fitting and inference

**Depends** R (>= 3.0.2),  
deSolve,  
adaptivetau,ggplot2,reshape2,plyr,tmvtnorm,parallel,doParallel,lubridate,coda,stringr,lattice,dplyr,truncnorm

**License** GPL-2

**LazyData** true

**Roxygen** list(wrap = FALSE)

**RoxygenNote** 5.0.1

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ABCLogPosterior	<i>ABC logged posterior distribution</i>
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### Description

This function evaluates the ABC posterior distribution at theta (using a single simulation trajectory) and returns the result in a suitable format for [mcmcMH](#).

### Usage

```
ABCLogPosterior(epsilon, sum.stats, distanceABC, fitmodel, theta, init.state,
  data)
```

### Arguments

epsilon	numeric vector, ABC tolerances for distances between data and simulations. If a vector of length 1 and the distance function returns a vector of distances, this will be expanded to be same tolerance for all the parameters.
sum.stats	a list of functions to calculate summary statistics. Each of these takes one argument (a trajectory with an "obs" column) and returns a number (the summary statistic given the trajectory)
distanceABC	a function that take three arguments: sum.stats, a list of summary statistics, data.obs (the data trajectory of observations) and model.obs (a model trajectory of observations), and returns the distance between the model run and the data in terms of the summary statistics
fitmodel	a <a href="#">fitmodel</a> object

theta	named numeric vector. Values of the parameters. Names should match <code>fitmodel\$theta.names</code> .
init.state	named numeric vector. Initial values of the state variables. Names should match <code>fitmodel\$state.names</code> .
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".

**Value**

a list of two elements

- `log.density` numeric, logged value of the ABC posterior distribution evaluated at theta
- `trace` named vector with trace information (theta, distance, log.density)

**See Also**

`computeDistanceABC`

---

burnAndThin	<i>Burn and thin MCMC chain</i>
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**Description**

Return a burned and thined trace of the chain.

**Usage**

```
burnAndThin(trace, burn = 0, thin = 0)
```

**Arguments**

trace	either a <code>data.frame</code> or a list of <code>data.frame</code> with all variables in column, as outputed by <code>mcmcMH</code> . Accept also an <code>mcmc</code> or <code>mcmc.list</code> object.
burn	proportion of the chain to burn.
thin	number of samples to discard per sample that is being kept

**Value**

an object with the same format as `trace` (`data.frame` or list of `data.frame` or `mcmc` object or `mcmc.list` object)

---

computeDIC	<i>Compute the DIC</i>
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---

### Description

This function computes the Deviance Information Criterion (DIC) of a `fitmodel` from a MCMC sample.

### Usage

```
computeDIC(trace, fitmodel, init.state, data, margLogLike = dTrajObs, ...)
```

### Arguments

<code>trace</code>	either a <code>data.frame</code> or <code>mcmc</code> object. Must contain one column with the posterior <code>log.likelihood</code> .
<code>fitmodel</code>	a <code>fitmodel</code> object
<code>init.state</code>	named numeric vector. Initial values of the state variables. Names should match <code>fitmodel\$state.names</code> .
<code>data</code>	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
<code>margLogLike</code>	R-function to compute the marginal log-likelihood of theta.
<code>...</code>	further arguments to be passed to <code>margLogLike</code>

### Value

a list of 5 elements:

- DIC value of the DIC
- `theta_bar` mean posterior of theta
- `log_like_theta_bar` log-likelihood of `theta_bar`
- `D_theta_bar` deviance of `theta_bar`
- `p_D` effective number of parameters

---

computeDistanceABC	<i>Compute the distance between a model and data for ABC</i>
--------------------	--

---

### Description

Compute the distance (using `distance.ABC`) between the observed time series and a simulated time series of observations obtained by running the model with parameters `theta`.

### Usage

```
computeDistanceABC(sum.stats, distanceABC, fitmodel, theta, init.state, data)
```

**Arguments**

<code>sum.stats</code>	a list of functions to calculate summary statistics. Each of these takes one argument (a trajectory with an "obs" column) and returns a number (the summary statistic given the trajectory)
<code>distanceABC</code>	a function that take three arguments: <code>sum.stats</code> , a list of summary statistics, <code>data.obs</code> (the data trajectory of observations) and <code>model.obs</code> (a model trajectory of observations), and returns the distance between the model run and the data in terms of the summary statistics
<code>fitmodel</code>	a <code>fitmodel</code> object
<code>theta</code>	named numeric vector. Values of the parameters. Names should match <code>fitmodel\$theta.names</code> .
<code>init.state</code>	named numeric vector. Initial values of the state variables. Names should match <code>fitmodel\$state.names</code> .
<code>data</code>	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".

**Value**

a sampled distance between

---

`distanceOscillation`     *Distance weighted by number of oscillations*

---

**Description**

This positive distance is the mean squared differences between  $x$  and the  $y$ , divided by the square of the number of times the  $x$  oscillates around the  $y$  (see note below for illustration).

**Usage**

```
distanceOscillation(x, y)
```

**Arguments**

`x`, `y`                numeric vectors of the same length.

**Note**

To illustrate this distance, suppose we observed a time series  $y = c(1, 3, 5, 7, 5, 3, 1)$  and we have two simulated time series  $x1 = (3, 5, 7, 9, 7, 5, 3)$  and  $x2 = (3, 5, 3, 5, 7, 5, 3)$ ;  $x1$  is consistently above  $y$  and  $x2$  oscillates around  $y$ . While the squared differences are the same, we obtain  $d(y, x1) = 4$  and  $d(y, x2) = 1.3$ .

---

dTrajObs

*Log-likelihood of a trajectory for a deterministic model*


---

### Description

Compute the trajectory log-likelihood of theta for a deterministic model defined in a [fitmodel](#) object by summing the point log-likelihoods.

### Usage

```
dTrajObs(fitmodel, theta, init.state, data, log = FALSE)
```

### Arguments

fitmodel	a <a href="#">fitmodel</a> object
theta	named numeric vector. Values of the parameters. Names should match fitmodel\$theta.names.
init.state	named numeric vector. Initial values of the state variables. Names should match fitmodel\$state.names.
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".

### Value

numeric value of the log-likelihood

### See Also

[rTrajObs](#)

---

export2Tracer

*Export trace in Tracer format*


---

### Description

Print trace in a file that can be read by the software Tracer.

### Usage

```
export2Tracer(trace, file)
```

### Arguments

trace	a data.frame with one column per estimated parameter, as returned by <a href="#">burnAndThin</a>
file	either a character string naming a file or a <a href="#">connection</a> open for writing. "" indicates output to the console.

### Note

Tracer is a program for analysing the trace files generated by Bayesian MCMC runs. It can be downloaded at <http://tree.bio.ed.ac.uk/software/tracer>.

**See Also**

burnAndThin

---

fitmodel	<i>Constructor of fitmodel object</i>
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**Description**

A `fitmodel` object is a list that stores some variables and functions that will be useful to simulate and fit your model during the course.

**Usage**

```
fitmodel(name = NULL, state.names = NULL, theta.names = NULL,
         simulate = NULL, rPointObs = NULL, dprior = NULL, dPointObs = NULL)
```

**Arguments**

- |             |   |
|-------------|---|
| name        | character. Name of the model (required).  |
| state.names | character vector. Names of the state variables i.e. <code>c("S", "I", "R")</code> (required).   |
| theta.names | character vector. Names of the parameters i.e. <code>c("R0", "infectious.period")</code> (required).  |
| simulate    | <p>R-function to simulate forward the model (required). This function takes 3 arguments:</p> <ul style="list-style-type: none"> <li>• <code>theta</code> named numeric vector. Values of the parameters. Names should match <code>theta.names</code>.</li> <li>• <code>init.state</code> named numeric vector. Initial values of the state variables. Names should match <code>state.names</code>.</li> <li>• <code>times</code> numeric vector. Time sequence for which the state of the model is wanted; the first value of <code>times</code> must be the initial time, i.e. the time of <code>init.state</code>.</li> </ul> <p>and returns a <code>data.frame</code> containing the simulated trajectories that is the values of the state variables (1 per column) at each observation time (1 per row). The first column is time.</p> |
| rPointObs   | <p>R-function that generates a (randomly sampled) observation point from a model point, using an observation model (optional). It thus acts as an inverse of <code>dPointObs</code> (see below). This function takes 2 arguments</p> <ul style="list-style-type: none"> <li>• <code>model.point</code> named numeric vector. State of the model at a given point in time.</li> <li>• <code>theta</code> named numeric vector. Values of the parameters. Names should match <code>theta.names</code>.</li> </ul> <p>and returns an observation point</p>   |
| dprior      | <p>R-function that evaluates the prior density of the parameters at a given <code>theta</code> (optional). The function should take 2 arguments:</p> <ul style="list-style-type: none"> <li>• <code>theta</code> named numeric vector. Values of the parameters. Names should match <code>theta.names</code>.</li> </ul>  |

- `log` boolean. determines whether the logarithm of the prior density should be returned.

and returns the (logged, if requested) value of the prior density distribution.

`dPointObs` R-function that evaluates the likelihood of one data point given the state of the model at the same time point. This function takes 4 arguments:

- `data.point` named numeric vector. Observation time and observed data point.
- `model.point` named numeric vector containing the state of the model at the observation time point.
- `theta` named numeric vector. Parameter values. Useful since parameters are usually needed to compute the likelihood (i.e. reporting rate).
- `log` boolean. determines whether the logarithm of the likelihood should be returned.

and returns the (log-)likelihood. (optional)

### Value

a `fitmodel` object that is a list of 7 elements:

- `name` character, name of the model
- `state.names` vector, names of the state variables.
- `theta.names` vector, names of the parameters.
- `simulate` R-function to simulate forward the model; usage: `simulate(theta, init.state, times)`.
- `rPointObs` R-function to generate simulated observations; usage: `rPointObs(model.point, theta)`.
- `dprior` R-function to evaluate the log-prior of the parameter values; usage: `dprior(theta)`.
- `dPointObs` R-function to evaluate the log-likelihood of one data point; usage: `dPointObs(data.point, model.point, theta, log)`.

### See Also

[testFitmodel](#)

### Examples

```
## create a simple deterministic SIR model with constant population size

SIR_name <- "SIR with constant population size"
SIR_state.names <- c("S", "I", "R")
SIR_theta.names <- c("R0", "D_inf")

SIR_simulateDeterministic <- function(theta, init.state, times) {

  SIR_ode <- function(time, state, parameters) {

    ## parameters
    beta <- parameters[["R0"]] / parameters[["D_inf"]]
    nu <- 1 / parameters[["D_inf"]]

    ## states
    S <- state[["S"]]
    I <- state[["I"]]
    R <- state[["R"]]
```



```

    N <- S + I + R

    dS <- -beta * S * I/N
    dI <- beta * S * I/N - nu * I
    dR <- nu * I

    return(list(c(dS, dI, dR)))
  }

  trajectory <- data.frame(ode(y = init.state,
                              times = times,
                              func = SIR_ode,
                              parms = theta,
                              method = "ode45"))

  return(trajectory)
}

## function to compute log-prior
SIR_prior <- function(theta, log = FALSE) {

  ## uniform prior on R0: U[1,100]
  log.prior.R0 <- dunif(theta[["R0"]], min = 1, max = 100, log = TRUE)
  ## uniform prior on infectious period: U[0,30]
  log.prior.D <- dunif(theta[["D_inf"]], min = 0, max = 30, log = TRUE)

  log.sum <- log.prior.R0 + log.prior.D

  return(ifelse(log, log.sum, exp(log.sum)))
}

## function to compute the likelihood of one data point
SIR_pointLike <- function(data.point, model.point, theta, log = FALSE){

  ## the prevalence is observed through a Poisson process
  return(dpois(x = data.point[["obs"]],
               lambda = model.point[["I"]],
               log = log))
}

## function to generate observation from a model simulation
SIR_genObsPoint <- function(model.point, theta){

  ## the prevalence is observed through a Poisson process
  obs.point <- rpois(n = 1, lambda = model.point[["I"]])

  return(c(obs = obs.point))
}

## create deterministic SIR fitmodel
SIR <- fitmodel(
  name = SIR_name,
  state.names = SIR_state.names,
  theta.names = SIR_theta.names,
  simulate = SIR_simulateDeterministic,
  dprior = SIR_prior,

```

```

rPointObs = SIR_genObsPoint,
dPointObs = SIR_pointLike)

## test
## theta <- c(R0 = 3, D_inf = 2)
## init.state <- c(S = 999, I = 1, R = 0)
## data(epi)

```

---

fitR

---

*fitR*


---

### Description

fitR

---

FluTdC1971

---

*Time-series of the 1971 influenza epidemic in Tristan-da-Cunha*


---

### Description

A dataset containing the daily incidence recorded during the 1971 influenza A/H3N2 two-wave epidemic on the island of Tristan-da-Cunha.

### Format

A data frame with 59 rows and 3 variables

### Details

- date calendar date of the record
- time day of record since beginning of epidemic
- Inc daily count incidence of influenza-like-illness

### Source

<http://eutils.ncbi.nlm.nih.gov/entrez/eutils/elink.fcgi?dbfrom=pubmed&id=4511951&retmode=ref&cmd=prlinks>

---

logPosterior	<i>Posterior distribution for a fitmodel</i>
--------------	--

---

### Description

This function evaluates the posterior distribution at theta and returns the result in a suitable format for [mcmcMH](#).

### Usage

```
logPosterior(fitmodel, theta, init.state, data, margLogLike = dTrajObs, ...)
```

### Arguments

fitmodel	a <a href="#">fitmodel</a> object
theta	named numeric vector. Values of the parameters. Names should match fitmodel\$theta.names.
init.state	named numeric vector. Initial values of the state variables. Names should match fitmodel\$state.names.
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
margLogLike	R-function to compute the marginal log-likelihood of theta.
...	further arguments to be passed to margLogLike

### Value

a list of two elements

- log.density numeric, logged value of the posterior density evaluated at theta
- trace named vector with trace information (theta, log.prior, marg.log.like, log.posterior)

### See Also

[dTrajObs](#), [margLogLikeSto](#)

---

logPosteriorWrapper	<i>A wrapper for logPosterior</i>
---------------------	-----------------------------------

---

### Description

A wrapper for [logPosterior](#) that returns a function that can be used as a target for [mcmcMH](#)

### Usage

```
logPosteriorWrapper(fitmodel, init.state, data, margLogLike, ...)
```

**Arguments**

fitmodel	a <a href="#">fitmodel</a> object
init.state	named numeric vector. Initial values of the state variables. Names should match fitmodel\$state.names.
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
margLogLike	R-function to compute the marginal log-likelihood of theta.
...	further arguments to be passed to margLogLike

**Value**

a R-function with one argument called theta.

---

margLogLikeSto	<i>Marginal log-likelihood for a stochastic model</i>
----------------	---

---

**Description**

Compute a Monte-Carlo estimate of the log-likelihood of theta for a stochastic model defined in a [fitmodel](#) object, using [particleFilter](#)

**Usage**

```
margLogLikeSto(fitmodel, theta, init.state, data, n.particles, n.cores = 1)
```

**Arguments**

fitmodel	a <a href="#">fitmodel</a> object
theta	named numeric vector. Values of the parameters. Names should match fitmodel\$theta.names.
init.state	named numeric vector. Initial values of the state variables. Names should match fitmodel\$state.names.
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
n.particles	number of particles
n.cores	number of cores on which propogation of the particles is parallelised. By default no parallelisation (n.cores=1). If NULL, set to the value returned by <a href="#">detectCores</a> .

**Value**

Monte-Carlo estimate of the marginal log-likelihood of theta

**See Also**

[particleFilter](#)

mcmcMH

*Metropolis-Hasting MCMC***Description**

Run `n.iterations` of a Metropolis-Hasting MCMC to sample from the target distribution using a gaussian proposal kernel. Two optional optimizations are also implemented: truncated gaussian proposal (to match the support of the target distribution, i.e. boundary of the parameters) and adaptative gaussian proposal (to match the size and the shape of the target distribution).

**Usage**

```
mcmcMH(target, init.theta, proposal.sd = NULL, n.iterations, covmat = NULL,
        limits = list(lower = NULL, upper = NULL), adapt.size.start = NULL,
        adapt.size.cooling = 0.99, adapt.shape.start = NULL,
        print.info.every = n.iterations/100, verbose = FALSE,
        max.scaling.sd = 50, acceptance.rate.weight = NULL,
        acceptance.window = NULL)
```

**Arguments**

<code>target</code>	R-function that takes a single argument: <code>theta</code> (named numeric vector of parameter values) and returns a list of 2 elements: <ul style="list-style-type: none"> <li>• <code>log.density</code> the logged value of the target density, evaluated at <code>theta</code>.</li> <li>• <code>trace</code> a named numeric vector of values to be printed in the <code>trace.data.frame</code> returned by <code>mcmcMH</code>.</li> </ul>
<code>init.theta</code>	named vector of initial parameter values to start the chain.
<code>proposal.sd</code>	vector of standard deviations. If this is given and <code>covmat</code> is not, a diagonal matrix will be built from this to use as covariance matrix of the multivariate Gaussian proposal distribution. By default, this is set to <code>init.theta/10</code> .
<code>n.iterations</code>	number of iterations to run the MCMC chain.
<code>covmat</code>	named numeric covariance matrix of the multivariate Gaussian proposal distribution. Must have named rows and columns with at least all estimated <code>theta</code> . If <code>proposal.sd</code> is given, this is ignored.
<code>limits</code>	limits for the - potentially truncated - multi-variate normal proposal distribution of the MCMC. Contains 2 elements: <ul style="list-style-type: none"> <li>• <code>lower</code> named numeric vector. Lower truncation points in each dimension of the Gaussian proposal distribution. By default they are set to <code>-Inf</code>.</li> <li>• <code>upper</code> named numeric vector. Upper truncation points in each dimension of the Gaussian proposal distribution. By default they are set to <code>Inf</code>.</li> </ul>
<code>adapt.size.start</code>	number of iterations to run before adapting the size of the proposal covariance matrix (see note below). Set to 0 (default) if size is not to be adapted.
<code>adapt.size.cooling</code>	cooling factor for the scaling factor of the covariance matrix during size adaptation (see note below).
<code>adapt.shape.start</code>	number of accepted jumps before adapting the shape of the proposal covariance matrix (see note below). Set to 0 (default) if shape is not to be adapted

<code>print.info.every</code>	frequency of information on the chain: acceptance rate and state of the chain. Default value to <code>n.iterations/100</code> . Set to NULL to avoid any info.
<code>verbose</code>	logical. If TRUE, information are printed.
<code>max.scaling.sd</code>	numeric. Maximum value for the scaling factor of the covariance matrix. Avoid too high values for the scaling factor, which might happen due to the exponential update scheme. In this case, the covariance matrix becomes too wide and the sampling from the truncated proposal kernel becomes highly inefficient
<code>acceptance.rate.weight</code>	if this is non-NULL, the acceptance rate is calculated as a moving average. This makes sure the acceptance rate is a reflection of recent proposals, and is useful for adaptive MCMC to prevent overshoot. If this is set to a value between 0 and 1, the acceptance rate will be calculated as $a = a_{\text{old}}(1 - w) + a_{\text{new}}w$ , where $a_{\text{old}}$ is the (moving) average of the acceptance rate at the previous iteration, $a_{\text{new}}$ is 1 if the current proposal is accepted and 0 if it is rejected, and $w$ is the value of <code>acceptance.rate.weight</code> . Values of $w$ close to 1 give a lot of weight to recent proposals, while values closer to 0 give more weight to older proposals. A reasonable starting value is in the order of $1e-2$ .
<code>acceptance.window</code>	if given, how many acceptances to store

### Value

a list with 3 elements:

- `trace` a data.frame. Each row contains a state of the chain (as returned by `target`).
- `acceptance.rate` acceptance rate of the MCMC chain.
- `covmat.empirical` empirical covariance matrix of the target sample.

### Note

The size of the proposal covariance matrix is adapted using the following formulae:

$$\Sigma_{n+1} = \sigma_n * \Sigma_n$$

with  $\sigma_n = \sigma_{n-1} * \exp(\alpha^n * (acc - 0.234))$ , where  $\alpha$  is equal to `adapt.size.cooling` and  $acc$  is the acceptance rate of the chain.

The shape of the proposal covariance matrix is adapted using the following formulae:

$$\Sigma_{n+1} = 2.38^2/d * \Sigma_n$$

with  $\Sigma_n$  the empirical covariance matrix and  $d$  is the number of estimated parameters in the model.

### References

Roberts GO, Rosenthal JS. Examples of adaptive MCMC. Journal of Computational and Graphical Statistics. Taylor & Francis; 2009;18(2):349-67.

measles

*Time-series of a measles outbreak***Description**

A dataset containing the weekly incidence recorded during a recent outbreak of measles in Europe

**Format**

A data frame with 37 rows and 2 variables

**Details**

- time week of the record
- Inc weekly recorded incidence of measles

particleFilter

*Run a particle filter for fitmodel object***Description**

The particle filter returns an estimate of the marginal log-likelihood  $L = p(y(t_{1:T})|\theta)$  as well as the set of filtered trajectories and their respective weights at the last observation time  $\omega(t_T) = p(y(t_T)|\theta)$ .

**Usage**

```
particleFilter(fitmodel, theta, init.state, data, n.particles,
  progress = FALSE, n.cores = 1)
```

**Arguments**

fitmodel	a <a href="#">fitmodel</a> object
theta	named numeric vector. Values of the parameters. Names should match <code>fitmodel\$theta.names</code> .
init.state	named numeric vector. Initial values of the state variables. Names should match <code>fitmodel\$state.names</code> .
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
n.particles	number of particles
progress	if TRUE progression of the filter is displayed in the console.
n.cores	number of cores on which propogation of the particles is parallelised. By default no parallelisation ( <code>n.cores=1</code> ). If NULL, set to the value returned by <a href="#">detectCores</a> .

**Value**

A list of 3 elements:

- dPointObs the marginal log-likelihood of the theta.
- traj a list of size n.particles with all filtered trajectories.
- traj.weight a vector of size n.particles with the normalised weight of the filtered trajectories.

**Note**

An unbiased state sample  $x(t_{0:T}) \sim p(X(t_{0:T})|\theta, y(t_{0:T}))$  can be obtained by sampling the set of trajectories traj with probability traj.weight.

**See Also**

plotSMC

---

plotESSBurn

---

*Plot Effective Sample Size (ESS) against burn-in*


---

**Description**

Takes an mcmc trace and tests the ESS at different values of burn-in

**Usage**

```
plotESSBurn(trace, longest.burn.in = ifelse(is.data.frame(trace) |
  is.mcmc(trace), nrow(trace), nrow(trace[[1]]))/2,
  step.size = round(longest.burn.in/50))
```

**Arguments**

trace	either a data.frame or a list of data.frame with all variables in column, as outputed by <code>mcmcMH</code> . Accept also mcmc or mcmc.list objects.
longest.burn.in	The longest burn-in to test. Defaults to half the length of the trace
step.size	The size of the steps of burn-in to test. Defaults to 1/50th of longest.burn.in

**Value**

a plot of the ESS against burn.in



plotFit

*Plot fit of model to data***Description**

This function simulates the model under `theta`, generates observation and plot them against the data. Since simulation and observation processes can be stochastic, `n.replicates` can be plotted.

**Usage**

```
plotFit(fitmodel, theta, init.state, data, n.replicates = 1, summary = TRUE,
        alpha = min(1, 10/n.replicates), all.vars = FALSE, non.extinct = NULL,
        plot = TRUE)
```

**Arguments**

<code>fitmodel</code>	a <a href="#">fitmodel</a> object
<code>theta</code>	named numeric vector. Values of the parameters. Names should match <code>fitmodel\$theta.names</code> .
<code>init.state</code>	named numeric vector. Initial values of the state variables. Names should match <code>fitmodel\$state.names</code> .
<code>data</code>	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
<code>n.replicates</code>	numeric, number of replicated simulations.
<code>summary</code>	logical. If TRUE, the mean, median as well as the 50th and 95th percentile of the trajectories are plotted (default). If FALSE, all individual trajectories are plotted (transparency can be set with <code>alpha</code> ).
<code>alpha</code>	transparency of the trajectories (between 0 and 1).
<code>all.vars</code>	logical, if FALSE only the observations are plotted. Otherwise, all state variables are plotted.
<code>non.extinct</code>	character vector. Names of the infected states which must be non-zero so the epidemic is still ongoing. When the names of these states are provided, the extinction probability is plotted by computing the proportion of faded-out epidemics over time. An epidemic has faded-out when all the infected states (whose names are provided) are equal to 0. This is only relevant for stochastic models. In addition, if <code>summary == TRUE</code> , the summaries of the trajectories conditioned on non-extinction are shown. Default to NULL.
<code>plot</code>	if TRUE the plot is displayed, and returned otherwise.

**Value**

if `plot == FALSE`, a list of 2 elements is returned:

- `simulations` data.frame of `n.replicates` simulated observations.
- `plot` the plot of the fit.

---

plotHPDregion2D	<i>2D highest posterior density region</i>
-----------------	--

---

**Description**

Given a sample from a multivariate posterior distribution, plot the bivariate region of highest marginal posterior density (HPD) for two variables with defined levels.

**Usage**

```
plotHPDregion2D(trace, vars, prob = c(0.95, 0.75, 0.5, 0.25, 0.1),
  xlab = NULL, ylab = NULL, plot = TRUE)
```

**Arguments**

trace	either a <code>data.frame</code> or <code>mcmc</code> object.
vars	which variables to plot: numeric or character vector
prob	probability level
xlab	x axis label
ylab	y axis label
plot	if TRUE the plot is displayed, and returned otherwise.

**Note**

HPD levels are computed using the function [HPDregionplot](#) from the package `emdbook`.

---

plotPosteriorDensity	<i>Plot MCMC posterior densities</i>
----------------------	--------------------------------------

---

**Description**

Plot the posterior density.

**Usage**

```
plotPosteriorDensity(trace, prior = NULL, colour = NULL, plot = TRUE)
```

**Arguments**

trace	either a <code>data.frame</code> or a list of <code>data.frame</code> with all variables in column, as outputed by <a href="#">mcmcMH</a> . Accept also an <code>mcmc</code> , a <code>mcmc.list</code> object or a list of <code>mcmc.list</code> .
prior	a <code>data.frame</code> containing the prior density. It must have the three following columns: <ul style="list-style-type: none"> <li>• theta names of the parameters</li> <li>• x value of the parameters</li> <li>• density density of the prior at x</li> </ul>
colour	named vector of two characters and containing colour names for posterior and prior distributions. Vector names must be posterior and prior.
plot	if TRUE the plot is displayed, and returned otherwise.

**See Also**

burnAndThin

---

plotPosteriorFit      *Plot MCMC posterior fit*


---

**Description**

Plot posterior distribution of observation generated under model's posterior parameter distribution.

**Usage**

```
plotPosteriorFit(trace, fitmodel, init.state, data,
  posterior.summary = c("sample", "median", "mean", "max"), summary = TRUE,
  sample.size = 100, non.extinct = NULL, alpha = min(1, 10/sample.size),
  plot = TRUE, all.vars = FALSE, init.date = NULL)
```

**Arguments**

trace	a data.frame with one column per estimated parameter, as returned by <a href="#">burnAndThin</a>
fitmodel	a <a href="#">fitmodel</a> object
init.state	named numeric vector. Initial values of the state variables. Names should match fitmodel\$state.names.
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
posterior.summary	character. Set to "sample" to plot trajectories from a sample of the posterior (default). Set to "median", "mean" or "max" to plot trajectories corresponding to the median, mean and maximum of the posterior density.
summary	logical, if TRUE trajectories are summarised by their mean, median, 50% and 95% quantile distributions. Otherwise, the trajectories are plotted.
sample.size	number of theta sampled from posterior distribution (if posterior.summary == "sample"). Otherwise, number of replicated simulations.
non.extinct	character vector. Names of the infected states which must be non-zero so the epidemic is still ongoing. When the names of these states are provided, the extinction probability is plotted by computing the proportion of faded-out epidemics over time. An epidemic has faded-out when all the infected states (whose names are provided) are equal to 0. This is only relevant for stochastic models. In addition, if summary == TRUE, the summaries of the trajectories conditioned on non-extinction are shown. Default to NULL.
alpha	transparency of the trajectories (between 0 and 1).
plot	if TRUE the plot is displayed, and returned otherwise.
all.vars	logical, if FALSE only the observations are plotted. Otherwise, all state variables are plotted.
init.date	character. Date of the first point of the time series (default to NULL). If provided, the x-axis will be in calendar format. NB: currently only works if the unit of time is the day.

**Value**

If `plot == FALSE`, a list of 2 elements is returned:

- `theta` the `theta(s)` used for plotting (either a vector or a `data.frame`)
- `traj` a `data.frame` with the trajectories (and observations) sampled from the posterior distribution.
- `plot` the plot of the fit displayed.

---

plotSMC	<i>Plot result of SMC</i>
---------	---------------------------

---

**Description**

Plot the observation generated by the filtered trajectories together with the data.

**Usage**

```
plotSMC(smc, fitmodel, theta, data = NULL, summary = TRUE, alpha = 1,
        all.vars = FALSE, plot = TRUE)
```

**Arguments**

<code>smc</code>	output of <a href="#">particleFilter</a>
<code>fitmodel</code>	a <a href="#">fitmodel</a> object
<code>theta</code>	named numeric vector. Values of the parameters. Names should match <code>fitmodel\$theta.names</code> .
<code>data</code>	data frame. Observation times and observed data. The time column must be named as given by <code>time.column</code> , whereas the name of the data column should match one of <code>fitmodel\$state.names</code> .
<code>summary</code>	logical. If TRUE, the mean, median as well as the 50th and 95th percentile of the trajectories are plotted (default). If FALSE, all individual trajectories are plotted (transparency can be set with <code>alpha</code> ).
<code>alpha</code>	transparency of the trajectories (between 0 and 1).
<code>all.vars</code>	logical, if FALSE only the observations are plotted. Otherwise, all state variables are plotted.
<code>plot</code>	if TRUE the plot is displayed, and returned otherwise.

**See Also**

[particleFilter](#)

---

plotTrace	<i>Plot MCMC trace</i>
-----------	------------------------

---

**Description**

Plot the traces of all estimated variables.

**Usage**

```
plotTrace(trace, estimated.only = FALSE)
```

**Arguments**

`trace` a `data.frame` with one column per estimated parameter, as returned by [burnAndThin](#)  
`estimated.only` logical, if TRUE only estimated parameters are displayed.

**See Also**

[burnAndThin](#)

---

plotTraj	<i>Plot model trajectories</i>
----------	--------------------------------

---

**Description**

This function use faceting to plot all trajectories in a data frame. Convenient to see results of several simulations, or data. Also, if data is present, then an additional plot is displayed with data and potentially observation generated.

**Usage**

```
plotTraj(traj = NULL, state.names = NULL, data = NULL,
         time.column = "time", lines.data = FALSE, summary = TRUE,
         replicate.column = "replicate", non.extinct = NULL, alpha = 1,
         plot = TRUE, colour = "red", init.date = NULL, same = FALSE)
```

**Arguments**

`traj` `data.frame`, output of `fitmodel$simulate` or `simulateModelReplicates`.  
`state.names` character vector. Names of the state variables to plot. Names must match `fitmodel$state.names`. If NULL (default) all state variables are plotted.  
`data` data frame. Observation times and observed data. The time column must be named as given by `time.column`, whereas the name of the data column should match one of `fitmodel$state.names`.  
`time.column` character vector. The column in the data that indicates time  
`lines.data` logical. If TRUE, the data will be plotted as lines  
`summary` logical. If TRUE, the mean, median as well as the 50th and 95th percentile of the trajectories are plotted (default). If FALSE, all individual trajectories are plotted (transparency can be set with `alpha`).

<code>replicate.column</code>	character Vector. The column in the data that indicates the replicate (if multiple replicates are to be plotted, i.e. if <code>summary</code> is <code>FALSE</code> )
<code>non.extinct</code>	character vector. Names of the infected states which must be non-zero so the epidemic is still ongoing. When the names of these states are provided, the extinction probability is plotted by computing the proportion of faded-out epidemics over time. An epidemic has faded-out when all the infected states (whose names are provided) are equal to 0. This is only relevant for stochastic models. In addition, if <code>summary == TRUE</code> , the summaries of the trajectories conditioned on non-extinction are shown. Default to <code>NULL</code> .
<code>alpha</code>	transparency of the trajectories (between 0 and 1).
<code>plot</code>	if <code>TRUE</code> the plot is displayed, and returned otherwise.
<code>colour</code>	character vector. If a character, will use that colour to plot trajectories. If "all", use all available colours. If <code>NULL</code> , don't set the colour.
<code>init.date</code>	character. Date of the first point of the time series (default to <code>NULL</code> ). If provided, the x-axis will be in calendar format. NB: currently only works if the unit of time is the day.

**See Also**

[simulateModelReplicates](#)

---

<code>printNamedVector</code>	<i>Print named vector</i>
-------------------------------	---------------------------

---

**Description**

Print named vector with format specified by `fmt` (2 decimal places by default).

**Usage**

```
printNamedVector(x, fmt = "%.2f", sep = " | ")
```

**Arguments**

<code>x</code>	named vector
<code>fmt</code>	a character vector of format strings, each of up to 8192 bytes.
<code>sep</code>	a character string to separate the terms. Not <a href="#">NA_character_</a> .

**See Also**

[sprintf](#)

rTrajObs

*Generate an observation trajectory for a fitmodel***Description**

This function simulates a model defined in a [fitmodel](#) object and generates observations at each time point. It returns the trajectory with an additions obs column.

**Usage**

```
rTrajObs(fitmodel, theta, init.state, times)
```

**Arguments**

fitmodel	a <a href="#">fitmodel</a> object
theta	named numeric vector. Values of the parameters. Names should match fitmodel\$theta.names.
init.state	named numeric vector. Initial values of the state variables. Names should match fitmodel\$state.names.
times	the times at which to generate observations

**Value**

data.frame

**See Also**

[dTrajObs](#)

SEIT2L\_deter

*The deterministic SEIT2L model with constant population size***Description**

The deterministic SEIT2L model with constant population size, uniform prior and Poisson observation with reporting rate.

**Format**

A [fitmodel](#) object, that is a list with the following elements:

**Details**

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.

- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing `SEIT2L_deter$simulate` for instance. There are some comments included.

---

SEIT2L\_pomp

*The SEIT2L model for pomp*


---

## Description

The SEIT2L model with constant population size, uniform prior and Poisson observation with reporting rate, as a [pomp](#) object.

## Format

A [pomp](#) object.

## Details

Look at the documentation of [pomp](#) for more details about each of these elements.

## Examples

```
## load pomp package
require('pomp')

## load Tristan da Cunha data
data(FluTdC1971)

## load SEITL_pomp object
data(SEITL_pomp)

## define deterministic skeleton
SEIT2L.skel.c <- '
    double trans[6];

    double beta = R0 / D_inf;
    double epsilon = 1 / D_lat;
    double nu = 1 / D_inf;
    double tau = 1 / D_imm;

    double N = S + E + I + T1 + T2 + L;

    trans[0] = beta * I / N * S;
    trans[1] = epsilon * E;
    trans[2] = nu * I;
    trans[3] = 2 * tau * T1;
    trans[4] = 2 * alpha * tau * T2;
    trans[5] = 2 * (1 - alpha) * tau * T2;

    DS = -trans[0] + trans[5];
    DE = trans[0] - trans[1];
    DI = trans[1] - trans[2];
    DT1 = trans[2] - trans[3];
```



```

    DT2 = trans[3] - trans[4] - trans[5];
    DL = trans[4];
    DInc = trans[1];
,

## define stochastic model, for use with euler.sim, see ?euler.sim
SEIT2L.sim.c <- '
    double rate[6];
    double dN[6];

    double beta = R0 / D_inf;
    double epsilon = 1 / D_lat;
    double nu = 1 / D_inf;
    double tau = 1 / D_imm;

    double N = S + E + I + T1 + T2 + L;

    rate[0] = beta * I / N;
    rate[1] = epsilon;
    rate[2] = nu;
    rate[3] = 2 * tau;
    rate[4] = 2 * alpha * tau;
    rate[5] = 2 * (1 - alpha) * tau;

    reulermultinom(1, S, &rate[0], dt, &dN[0]);
    reulermultinom(1, E, &rate[1], dt, &dN[1]);
    reulermultinom(1, I, &rate[2], dt, &dN[2]);
    reulermultinom(1, T1, &rate[3], dt, &dN[3]);
    reulermultinom(2, T2, &rate[4], dt, &dN[4]);

    S += -dN[0] + dN[5];
    E += dN[0] - dN[1];
    I += dN[1] - dN[2];
    T1 += dN[2] - dN[3];
    T2 += dN[3] - dN[4] - dN[5];
    L += dN[4];
    Inc += dN[1];
,

## construct pomp object
SEIT2L_pomp <- pomp(data = FluTdc1971[, c("time", "obs")],
    skeleton = Csnippet(SEIT2L.skel.c),
    skeleton.type = "vectorfield",
    rprocess = euler.sim(step.fun = Csnippet(SEIT2L.sim.c),
        delta.t = 0.1),
    rmeasure = Csnippet(SEITL.rmeas.c),
    dmeasure = Csnippet(SEITL.dmeas.c),
    toEstimationScale = Csnippet(SEITL.logtrans.c),
    fromEstimationScale = Csnippet(SEITL.exptrans.c),
    times = "time",
    t0 = 1,
    zeronames = "Inc",
    paramnames = c("R0", "D_inf", "D_lat", "D_imm", "alpha", "rho"),
    statenames = c("S", "E", "I", "T1", "T2", "L", "Inc"),
    obsnames = c("obs"))

```

SEIT2L\_stoch

*The stochastic SEIT2L model with constant population size***Description**

The stochastic SEIT2L model with constant population size, uniform prior and Poisson observation with reporting rate.

**Format**

A [fitmodel](#) object, that is a list with the following elements:

**Details**

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.
- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing SEIT2L\_stoch\$simulate for instance. There are some comments included.

SEIT4L\_deter

*The deterministic SEIT4L model with constant population size***Description**

The deterministic SEIT4L model with constant population size, uniform prior and Poisson observation with reporting rate.

**Format**

A [fitmodel](#) object, that is a list with the following elements:

**Details**

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.
- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing SEIT4L\_deter\$simulate for instance. There are some comments included.

SEIT4L\_pomp

*The SEIT4L model for pomp***Description**

The SEIT4L model with constant population size, uniform prior and Poisson observation with reporting rate, as a [pomp](#) object.

**Format**

A [pomp](#) object.

**Details**

Look at the documentation of [pomp](#) for more details about each of these elements.

**Examples**

```
## load pomp package
require('pomp')

## load Tristan da Cunha data
data(FluTdC1971)

## load SEITL_pomp object
example(SEITL_pomp)

## define deterministic skeleton
SEIT4L.skel.c <- '
    double trans[8];

    double beta = R0 / D_inf;
    double epsilon = 1 / D_lat;
    double nu = 1 / D_inf;
    double tau = 1 / D_imm;

    double N = S + E + I + T1 + T2 + L;

    trans[0] = beta * I / N * S;
    trans[1] = epsilon * E;
    trans[2] = nu * I;
    trans[3] = 4 * tau * T1;
    trans[4] = 4 * tau * T2;
    trans[5] = 4 * tau * T3;
    trans[6] = 4 * alpha * tau * T4;
    trans[7] = 4 * (1 - alpha) * tau * T4;

    DS = -trans[0] + trans[7];
    DE = trans[0] - trans[1];
    DI = trans[1] - trans[2];
    DT1 = trans[2] - trans[3];
    DT2 = trans[3] - trans[4];
    DT3 = trans[4] - trans[5];
    DT4 = trans[5] - trans[6];
```

```

    DL = trans[6];
    DInc = trans[1];
,

## define stochastic model, for use with euler.sim, see ?euler.sim
SEIT4L.sim.c <- '
    double rate[8];
    double dN[8];

    double beta = R0 / D_inf;
    double epsilon = 1 / D_lat;
    double nu = 1 / D_inf;
    double tau = 1 / D_imm;

    double N = S + E + I + T1 + T2 + L;

    rate[0] = beta * I / N;
    rate[1] = epsilon;
    rate[2] = nu;
    rate[3] = 4 * tau;
    rate[4] = 4 * tau;
    rate[5] = 4 * tau;
    rate[6] = 4 * alpha * tau;
    rate[7] = 4 * (1 - alpha) * tau;

    reulermultinom(1, S, &rate[0], dt, &dN[0]);
    reulermultinom(1, E, &rate[1], dt, &dN[1]);
    reulermultinom(1, I, &rate[2], dt, &dN[2]);
    reulermultinom(1, T1, &rate[3], dt, &dN[3]);
    reulermultinom(1, T2, &rate[4], dt, &dN[4]);
    reulermultinom(1, T3, &rate[5], dt, &dN[5]);
    reulermultinom(2, T4, &rate[6], dt, &dN[6]);

    S += -dN[0] + dN[7];
    E += dN[0] - dN[1];
    I += dN[1] - dN[2];
    T1 += dN[2] - dN[3];
    T2 += dN[3] - dN[4];
    T3 += dN[4] - dN[5];
    T4 += dN[5] - dN[6] - dN[7];
    L += dN[6];
    Inc += dN[1];
,

## construct pomp object
SEIT4L_pomp <- pomp(data = FluTdc1971[, c("time", "obs")],
    skeleton = Csnippet(SEIT4L.skel.c),
    skeleton.type = "vectorfield",
    rprocess = euler.sim(step.fun = Csnippet(SEIT4L.sim.c),
        delta.t = 0.1),
    rmeasure = Csnippet(SEITL.rmeas.c),
    dmeasure = Csnippet(SEITL.dmeas.c),
    toEstimationScale = Csnippet(SEITL.logtrans.c),
    fromEstimationScale = Csnippet(SEITL.exptrans.c),
    times = "time",
    t0 = 1,
    zeronames = "Inc",

```

```
paramnames = c("R0", "D_inf", "D_lat", "D_imm", "alpha", "rho"),
statenames = c("S", "E", "I", "T1", "T2", "T3", "T4", "L", "Inc"),
obsnames = c("obs"))
```

SEIT4L\_stoch

*The stochastic SEIT4L model with constant population size***Description**

The stochastic SEIT4L model with constant population size, uniform prior and Poisson observation with reporting rate.

**Format**

A [fitmodel](#) object, that is a list with the following elements:

**Details**

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.
- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing `SEIT4L_stoch$simulate` for instance. There are some comments included.

SEITL\_deter

*The deterministic SEITL model with constant population size***Description**

The deterministic SEITL model with constant population size, uniform prior and Poisson observation with reporting rate.

**Format**

A [fitmodel](#) object, that is a list with the following elements:

## Details

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.
- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing `SEITL_deter$simulate` for instance. There are some comments included.

---

SEITL\_distanceOscillation

*ABC distance with oscillations*

---

## Description

This positive distance is the mean squared differences between the simulation and the observation, divided by the square of the number of times the simulation oscillates around the observation.

## Usage

```
SEITL_distanceOscillation(simu.traj.obs, data)
```

## Arguments

`simu.traj.obs` data.frame of simulated trajectory with observation, as returned by [rTrajObs](#).  
`data` data.frame of times and observations. Must have two columns: time and Inc.

## See Also

`distanceOscillation`

## Examples

```
## Not run:
# Suppose we observed a time series:
data <- data.frame(time=1:7, Inc=c(1,3,5,7,5,3,1))
# and we have two simulated time series:
traj1 <- data.frame(time=1:7, observation=c(3,5,7,9,7,5,3))
traj2 <- data.frame(time=1:7, observation=c(3,5,3,5,7,5,3))
# traj1 is consistently above data and traj2 oscillates around data:
plot(data$time, data$Inc, t='l', ylim=c(0,10))
lines(traj1$time, traj1$observation, col="red")
lines(traj2$time, traj2$observation, col="blue")
# While the squared differences are the same, we obtain a smaller distance for traj2:
d1 <- SEITL_distanceOscillation(traj1, data)
# d1 = 4
d2 <- SEITL_distanceOscillation(traj2, data)
```

```
# d2 = 1.3

## End(Not run)
```

---

SEITL\_pomp

---

*The SEITL model for pomp*


---

## Description

The SEITL model with constant population size, uniform prior and Poisson observation with reporting rate, as a [pomp](#) object.

## Format

A [pomp](#) object.

## Details

Look at the documentation of [pomp](#) for more details about each of these elements.

## Examples

```
## load pomp package
require('pomp')

## load Tristan da Cunha data
data(FluTdC1971)

## define deterministic skeleton
SEITL.skel.c <- '
  double trans[5];

  double beta = R0 / D_inf;
  double epsilon = 1 / D_lat;
  double nu = 1 / D_inf;
  double tau = 1 / D_imm;

  double N = S + E + I + T + L;

  trans[0] = beta * I / N * S;
  trans[1] = epsilon * E;
  trans[2] = nu * I;
  trans[3] = alpha * tau * T;
  trans[4] = (1 - alpha) * tau * T;

  DS = -trans[0] + trans[4];
  DE = trans[0] - trans[1];
  DI = trans[1] - trans[2];
  DT = trans[2] - trans[3] - trans[4];
  DL = trans[3];
  DInc = trans[1];
,

## define stochastic model, for use with euler.sim, see ?euler.sim
```

```

SEITL.sim.c <- '
  double rate[5];
  double dN[5];

  double beta = R0 / D_inf;
  double epsilon = 1 / D_lat;
  double nu = 1 / D_inf;
  double tau = 1 / D_imm;

  double N = S + E + I + T + L;

  rate[0] = beta * I / N;
  rate[1] = epsilon;
  rate[2] = nu;
  rate[3] = alpha * tau;
  rate[4] = (1 - alpha) * tau;

  reulermultinom(1, S, &rate[0], dt, &dN[0]);
  reulermultinom(1, E, &rate[1], dt, &dN[1]);
  reulermultinom(1, I, &rate[2], dt, &dN[2]);
  reulermultinom(2, T, &rate[3], dt, &dN[3]);

  S += -dN[0] + dN[4];
  E += dN[0] - dN[1];
  I += dN[1] - dN[2];
  T += dN[2] - dN[3] - dN[4];
  L += dN[3];
  Inc += dN[1];
,

## define sampling random point observations
SEITL.rmeas.c <- '
  obs = rpois(rho * Inc > 0 ? rho * Inc : 0);
,

## define point observation probability density
SEITL.dmeas.c <- '
  lik = dpois(obs, rho * Inc > 0 ? rho * Inc : 0, give_log);
,

## define prior density
SEITL.dprior.c <- '
  lik = dunif(R0, 1, 50, 1) +
        dunif(D_lat, 0, 10, 1) +
        dunif(D_inf, 0, 15, 1) +
        dunif(D_imm, 0, 50, 1) +
        dunif(alpha, 0, 1, 1) +
        dunif(rho, 0, 1, 1);

  lik = give_log ? lik : exp(lik);
,

SEITL.logtrans.c <- '
  TR0 = log(R0);
  TD_inf = log(D_inf);
  TD_lat = log(D_lat);
  TD_imm = log(D_imm);

```



```

    Talpha = log(alpha);
    Trho = log(rho);
  ,

SEITL.exptrans.c <- '
  TR0 = exp(R0);
  TD_inf = exp(D_inf);
  TD_lat = exp(D_lat);
  TD_imm = exp(D_imm);
  Talpha = exp(alpha);
  Trho = exp(rho);
'

## construct pomp object
SEITL_pomp <- pomp(data = FluTdC1971[, c("time", "obs")],
  skeleton = Csnippet(SEITL.skel.c),
  skeleton.type = "vectorfield",
  rprocess = euler.sim(step.fun = Csnippet(SEITL.sim.c),
    delta.t = 0.1),
  rmeasure = Csnippet(SEITL.rmeas.c),
  dmeasure = Csnippet(SEITL.dmeas.c),
  dprior = Csnippet(SEITL.dprior.c),
  toEstimationScale = Csnippet(SEITL.logtrans.c),
  fromEstimationScale = Csnippet(SEITL.exptrans.c),
  times = "time",
  t0 = 1,
  zeronames = "Inc",
  paramnames = c("R0", "D_inf", "D_lat", "D_imm", "alpha", "rho"),
  statenames = c("S", "E", "I", "T", "L", "Inc"),
  obsnames = c("obs"))

```

SEITL\_stoch

*The stochastic SEITL model with constant population size***Description**

The stochastic SEITL model with constant population size, uniform prior and Poisson observation with reporting rate.

**Format**

A `fitmodel` object, that is a list with the following elements:

**Details**

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.

- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing `SEITL_stoch$simulate` for instance. There are some comments included.

---

```
simulateFinalStateAtExtinction
```

*Simulate model until extinction*

---

### Description

Return final state at extinction

### Usage

```
simulateFinalStateAtExtinction(fitmodel, theta, init.state, extinct = NULL,
  time.init = 0, time.step = 100, n = 100, observation = FALSE,
  n.cores = 1)
```

### Arguments

<code>fitmodel</code>	a <a href="#">fitmodel</a> object
<code>theta</code>	named numeric vector. Values of the parameters. Names should match <code>fitmodel\$theta.names</code> .
<code>init.state</code>	named numeric vector. Initial values of the state variables. Names should match <code>fitmodel\$state.names</code> .
<code>extinct</code>	character vector. Simulations stop when all these state are extinct.
<code>time.init</code>	numeric. Start time of simulation.
<code>time.step</code>	numeric. Time step at which extinction is checked
<code>n</code>	number of replicated simulations.
<code>observation</code>	logical, if TRUE simulated observation are generated by <code>rTrajObs</code> .
<code>n.cores</code>	number of cores on which propagation of the particles is parallelised. By default no parallelisation ( <code>n.cores=1</code> ). If NULL, set to the value returned by <a href="#">detectCores</a> .

---

```
simulateModelReplicates
```

*Simulate several replicate of the model*

---

### Description

Simulate several replicate of a `fitmodel` using its function `simulate`

### Usage

```
simulateModelReplicates(fitmodel, theta, init.state, times, n,
  observation = FALSE)
```

**Arguments**

<code>fitmodel</code>	a <a href="#">fitmodel</a> object
<code>theta</code>	named numeric vector. Values of the parameters. Names should match <code>fitmodel\$theta.names</code> .
<code>init.state</code>	named numeric vector. Initial values of the state variables. Names should match <code>fitmodel\$state.names</code> .
<code>times</code>	vector of times at which you want to observe the states of the model.
<code>n</code>	number of replicated simulations.
<code>observation</code>	logical, if TRUE simulated observation are generated by <code>rTrajObs</code> .

**Value**

a data.frame of dimension `[nlength(times)]x[length(init.state)+2]` with column names equal to `c("replicate", "time", names(init.state))`.

---

```
simulateModelStochastic
```

*Simulate forward a stochastic model*

---

**Description**

This function uses the function [ssa.adaptivetau](#) to simulate the model and returns the trajectories in a valid format for the class [fitmodel](#).

**Usage**

```
simulateModelStochastic(theta, init.state, times, transitions, rateFunc)
```

**Arguments**

<code>theta</code>	named vector of model parameters.
<code>init.state</code>	named vector of initial state of the model.
<code>times</code>	time sequence for which state of the model is wanted; the first value of times must be the initial time.
<code>transitions</code>	One of two possible data types: <ul style="list-style-type: none"> <li>• A list with length equal to the number of transitions. Each element of the list should be a vector specifying a transition (i.e., which state(s) change and by how much). Each entry in the vector needs a name (specifying which state variable to change, either by name or index) and a value (specifying the amount by which this variable will change).</li> <li>• A two-dimensional matrix of integers specifying how each state variable (rows) should be changed for a given transition (columns). Generally this will be a sparse matrix of primarily 1s and -1s, which can make this structure inefficient.</li> </ul>

See the example below for details as well as [ssa.maketrans](#) or the vignette accompanying this package.

`rateFunc` R function that returns instantaneous transition rates for each transition in the form a real-valued one-dimensional vector with length equal to the number of transitions. The order of these rates must match the order in which transitions were specified in the `transitions` parameter above. This function must accept the following arguments:

- vector of current values for all state variables (in order used in the `init.values` argument above)
- parameters as supplied in argument to `ssa.adaptivetau`
- single real number giving the current time (all simulations start at  $t=0$ )

### Value

a data.frame of dimension `length(times) × (length(init.state)+1)` with column names equal to `c("time", names(init.state))`.

---

SIR

*A simple deterministic SIR model with constant population size*

---

### Description

A simple deterministic SIR model with constant population size, uniform prior and Poisson observation.

### Format

A `fitmodel` object, that is a list with the following elements:

### Details

- `name` character.
- `state.names` character vector.
- `theta.names` character vector.
- `simulate` R-function.
- `rPointObs` R-function.
- `dprior` R-function.
- `dPointObs` R-function.

Look at the documentation of `fitmodel` for more details about each of these elements. You can look at the code of the R-functions by typing `SIR$simulate` for instance. There are some comments included.

---

SIR_exp	<i>A simple deterministic SIR model with constant population size and parameters on the exponential scale</i>
---------	---

---

### Description

A simple deterministic SIR model with constant population size, uniform prior and Poisson observation. The parameters are transformed using an exponential transformation.

### Format

A `fitmodel` object, that is a list with the following elements:

### Details

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.
- dPointObs R-function.

Look at the documentation of `fitmodel` for more details about each of these elements. You can look at the code of the R-functions by typing `SIR_exp$simulate` for instance. There are some comments included.

---

SIR_reporting	<i>A simple deterministic SIR model with constant population size and reporting rate</i>
---------------	--

---

### Description

A simple deterministic SIR model with constant population size, uniform prior and Poisson observation with reporting rate.

### Format

A `fitmodel` object, that is a list with the following elements:

### Details

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.

- dprior R-function.
- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing `SIR_reporting$simulate` for instance. There are some comments included.

---

SIR\_stoch

*A simple stochastic SIR model with constant population size*

---

### Description

A simple stochastic SIR model with constant population size, uniform prior and Poisson observation.

### Format

A [fitmodel](#) object, that is a list with the following elements:

### Details

- name character.
- state.names character vector.
- theta.names character vector.
- simulate R-function.
- rPointObs R-function.
- dprior R-function.
- dPointObs R-function.

Look at the documentation of [fitmodel](#) for more details about each of these elements. You can look at the code of the R-functions by typing `SIR_stoch$simulate` for instance. There are some comments included.

---

testFitmodel

*Test a fitmodel*

---

### Description

This function performs a serie of checks on the [fitmodel](#) provided by the user in order to make sure that it will be compatible both with the functions coded during the course and the functions available in the `fitR` package. The latters can be used as a correction.

### Usage

```
testFitmodel(fitmodel, theta, init.state, data = NULL, verbose = TRUE)
```

**Arguments**

fitmodel	a <a href="#">fitmodel</a> object
theta	named numeric vector. Values of the parameters. Names should match fitmodel\$theta.names.
init.state	named numeric vector. Initial values of the state variables. Names should match fitmodel\$state.names.
data	data frame. Observation times and observed data. The time column must be named "time" and the observation column must be named "obs".
verbose	if TRUE, print details of the test performed to check validity of the arguments

**See Also**[fitmodel](#)**Examples**

```
## create a simple deterministic SIR model with constant population size

SIR_name <- "SIR with constant population size"
SIR_state.names <- c("S", "I", "R")
SIR_theta.names <- c("R0", "D_inf")

SIR_simulateDeterministic <- function(theta, init.state, times) {

  SIR_ode <- function(time, state, parameters) {

    ## parameters
    beta <- parameters[["R0"]] / parameters[["D_inf"]]
    nu <- 1 / parameters[["D_inf"]]

    ## states
    S <- state[["S"]]
    I <- state[["I"]]
    R <- state[["R"]]

    N <- S + I + R

    dS <- -beta * S * I/N
    dI <- beta * S * I/N - nu * I
    dR <- nu * I

    return(list(c(dS, dI, dR)))
  }

  trajectory <- data.frame(ode(y = init.state,
                              times = times,
                              func = SIR_ode,
                              parms = theta,
                              method = "ode45"))

  return(trajectory)
}

## function to compute log-prior
SIR_prior <- function(theta, log = FALSE) {
```

```

## uniform prior on R0: U[1,100]
log.prior.R0 <- dunif(theta[["R0"]], min = 1, max = 100, log = TRUE)
## uniform prior on infectious period: U[0,30]
log.prior.D <- dunif(theta[["D_inf"]], min = 0, max = 30, log = TRUE)

log.sum <- log.prior.R0 + log.prior.D

return(ifelse(log, log.sum, exp(log.sum)))
}

## function to compute the likelihood of one data point
SIR_pointLike <- function(data.point, model.point, theta, log = FALSE){

  ## the prevalence is observed through a Poisson process
  return(dpois(x = data.point[["obs"]],
               lambda = model.point[["I"]],
               log = log))
}

## function to generate observation from a model simulation
SIR_genObsPoint <- function(model.point, theta){

  ## the prevalence is observed through a Poisson process
  obs.point <- rpois(n = 1, lambda = model.point[["I"]])

  return(c(obs = obs.point))
}

## create deterministic SIR fitmodel
SIR <- fitmodel(
  name = SIR_name,
  state.names = SIR_state.names,
  theta.names = SIR_theta.names,
  simulate = SIR_simulateDeterministic,
  dprior = SIR_prior,
  rPointObs = SIR_genObsPoint,
  dPointObs = SIR_pointLike)

## test
## theta <- c(R0 = 3, D_inf = 2)
## init.state <- c(S = 999, I = 1, R = 0)
## data(epi)

```

---

updateCovmat

*Update covariance matrix*


---

## Description

Update covariance matrix using a stable one-pass algorithm. This is much more efficient than using [cov](#) on the full data.

## Usage

```
updateCovmat(covmat, theta.mean, theta, i)
```



**Arguments**

<code>covmat</code>	covariance matrix at iteration $i-1$ . Must be numeric, symmetrical and named.
<code>theta.mean</code>	mean vector at iteration $i-1$ . Must be numeric and named.
<code>theta</code>	vector of new value at iteration $i$ . Must be numeric and named.
<code>i</code>	current iteration.

**Value**

A list of two elements

- `covmat` update covariance matrix
- `theta.mean` updated mean vector

**References**

<http://en.wikipedia.org/wiki/Algorithms%5Ffor%5Fcalculating%5Fvariance#Covariance>

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