

Package ‘spatPomp’

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Type Package

Title Inference for Spatiotemporal Partially Observed Markov Processes

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URL <https://github.com/kidusasfaw/spatPomp>

Description Inference on panel data using spatiotemporal partially-observed Markov process (Spat-POMP) models. The 'spatPomp' package extends 'pomp' to include algorithms taking advantage of the spatial structure in order to assist with handling high dimensional processes. See Asfaw et al. (2024) <[doi:10.48550/arXiv.2101.01157](https://doi.org/10.48550/arXiv.2101.01157)> for further description of the package.

SystemRequirements For Windows users, Rtools (see <https://cran.r-project.org/bin/windows/Rtools/>).

License GPL-3

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BugReports <https://github.com/kidusasfaw/spatPomp/issues/>

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LinkingTo pomp

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Imports foreach, dplyr, tidyr, stringr, abind, rlang, ggplot2

Collate 'spatPomp_class.R' 'abf.R' 'abfir.R' 'arma_benchmark.R'
'get_covariate_names.R' 'as_data_frame.R' 'bm.R' 'bm2.R'
'bm2_kalman_logLik.R' 'bm_kalman_logLik.R' 'bpfilter.R'
'city_data_UK.R' 'iter_filter.R' 'safecall.R' 'pstop.R'
'undefined.R' 'spatPomp.R' 'ibpf.R' 'listie.R' 'conc.R'
'concat.R' 'dunit_measure.R' 'enkf.R' 'eunit_measure.R' 'gbm.R'
'girf.R' 'he10.R' 'he10coordinates.R' 'he10demography.R'
'he10measles.R' 'he10mle.R' 'ienkf.R' 'igirf.R' 'iubf.R'
'loglik.R' 'lorenz.R' 'measles.R' 'measles2.R' 'measlesUK.R'
'munit_measure.R' 'package.R' 'param_formats.R' 'plot.R'
'print.R' 'runit_measure.R' 'simulate.R' 'spatPomp_Csnippet.R'
'spatPomp_workhorses.R' 'unit_names.R' 'vec_dmeasure.R'
'vec_rmeasure.R' 'vunit_measure.R'

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spatPomp-package *Inference for SpatPOMPs (Spatiotemporal Partially Observed Markov Processes)*

Description

The **spatPomp** package provides facilities for inference on panel data using spatiotemporal partially-observed Markov process (SPATPOMP) models. To do so, it relies on and extends a number of facilities that the **pomp** package provides for inference on time series data using partially-observed Markov process (POMP) models.

The **spatPomp** package concerns models consisting of a collection of interacting units. The methods in **spatPomp** may be applicable whether or not these units correspond to spatial locations.

Data analysis using spatPomp

The first step in using **spatPomp** is to encode one’s model(s) and data in objects of class spatPomp. This can be done via a call to the [spatPomp](#) constructor function.

Extending the pomp platform for developing inference tools

spatPomp extends to panel data the general interface to the components of POMP models provided by **pomp**. In doing so, it contributes to the goal of the **pomp** project of facilitating the development of new algorithms in an environment where they can be tested and compared on a growing body of models and datasets.

Documentation

spatPomp is described by Asfaw et al. (2020)

License

spatPomp is provided under the MIT License.

Author(s)

Kidus Asfaw, Joonha Park, Allister Ho, Edward Ionides, Aaron A. King

References

Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package spatPomp. *ArXiv*: 2101.01157. doi:[10.48550/arXiv.2101.01157](https://doi.org/10.48550/arXiv.2101.01157)

See Also

[pomp](#)

 abf

Adapted Bagged Filter (ABF)

Description

An algorithm for estimating the likelihood of a spatiotemporal partially-observed Markov process model. Running `abf` causes the algorithm to run bootstrap replicate jobs which each yield an imperfect adapted simulation. Simulating from the "adapted filter" distribution runs into a curse of dimensionality (COD) problem, which is mitigated by keeping particles in each replicate close to each other through resampling down to one particle per replicate at each observation time point. The adapted simulations are then weighted in a way that mitigates COD by making a weak coupling assumption to get an approximate filter distribution. As a by-product, we also get an estimate of the likelihood of the data.

Usage

```
## S4 method for signature 'spatPomp'
abf(
  object,
  Nrep,
  Np,
  nbhd,
  params,
  tol = 1e-100,
  ...,
  verbose = getOption("verbose", FALSE)
)

## S4 method for signature 'abfd_spatPomp'
abf(object, Nrep, Np, nbhd, tol, ..., verbose = getOption("verbose", FALSE))
```

Arguments

object	A spatPomp object.
Nrep	The number of bootstrap replicates for the adapted simulations.
Np	The number of particles used within each replicate for the adapted simulations.
nbhd	A neighborhood function with three arguments: object, time and unit. The function should return a list of two-element vectors that represent space-time neighbors of (u, n) , which is represented by $c(\text{unit}, \text{time})$. See example below for more details.
params	Parameter vector, required if not available from the spatPomp model object.
tol	If the resampling weight for a particle is zero due to floating-point precision issues, it is set to the value of tol since resampling has to be done.
...	Additional arguments can be used to replace model components.
verbose	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

Upon successful completion, `abf()` returns an object of class ‘`abfd_spatPomp`’ containing the algorithmic parameters used to run `abf()` and the estimated likelihood.

Methods

The following methods are available for such an object:

`logLik` yields an estimate of the log-likelihood of the data under the model.

Author(s)

Kidus Asfaw

References

Ionides, E. L., Asfaw, K., Park, J., and King, A. A. (2021). Bagged filters for partially observed interacting systems. *Journal of the American Statistical Association*, doi:10.1080/01621459.2021.1974867

See Also

likelihood maximization algorithms: `ienkf()`, `igirf`, `iubf`, `ibpf`

Other likelihood evaluation algorithms: `abfir()`, `bpfilter()`, `enkf()`, `girf()`

Examples

```
# Complete examples are provided in the package tests
## Not run:
# Create a simulation of a Brownian motion
b <- bm(U=2, N=5)

# Create a neighborhood function mapping a point in space-time
```

```

# to a list of neighboring points in space-time
bm_nbhd <- function(object, time, unit) {
  nbhd_list = list()
  if(time > 1 && unit > 1){
    nbhd_list = c(nbhd_list, list(c(unit-1, time-1)))
  }
  return(nbhd_list)
}

# Run ABF specified number of Monte Carlo replicates and particles per replicate
abfd_bm <- abf(b, Nrep=2, Np=10, nbhd=bm_nbhd)

# Get the likelihood estimate from ABF
logLik(abfd_bm)

## End(Not run)

```

abfir

Adapted Bagged Filter with Intermediate Resampling (ABF-IR)

Description

An algorithm for estimating the filter distribution and likelihood of a spatiotemporal partially-observed Markov process model. Running `abfir` causes the algorithm to run Monte Carlo replicated jobs which each carry out an adapted simulation using intermediate resampling. Adapted simulation is an easier task than filtering, since particles in each replicate remain close to each other. Intermediate resampling further assists against the curse of dimensionality (COD) problem for importance sampling. The adapted simulations are then weighted in a way that mitigates COD by making a weak coupling assumption to get an approximate filter distribution. As a by-product, we also get an approximation to the likelihood of the data.

Usage

```

## S4 method for signature 'spatPomp'
abfir(
  object,
  Np,
  Nrep,
  nbhd,
  Ninter,
  tol = (1e-100),
  params,
  ...,
  verbose = getOption("verbose", FALSE)
)

## S4 method for signature 'abfird_spatPomp'
abfir(object, Np, Nrep, nbhd, Ninter, tol, params, ...)

```

Arguments

object	A spatPomp object.
Np	The number of particles used within each replicate for the adapted simulations.
Nrep	The number of bootstrap replicates for the adapted simulations.
nbhd	A neighborhood function with three arguments: object, time and unit. The function should return a list of two-element vectors that represent space-time neighbors of (u, n) , which is represented by $c(\text{unit}, \text{time})$. See example below for more details.
Ninter	the number of intermediate resampling time points. By default, this is set equal to the number of units.
tol	If the resampling weight for a particle is zero due to floating-point precision issues, it is set to the value of tol since resampling has to be done.
params	Parameter vector, required if not available from the spatPomp model object.
...	Additional arguments can be used to replace model components.
verbose	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

Upon successful completion, `abfir()` returns an object of class ‘`abfir_spatPomp`’ containing the algorithmic parameters used to run `abfir()` and the estimated likelihood.

Methods

The following methods are available for such an object:

`logLik` yields a biased estimate of the log-likelihood of the data under the model.

Author(s)

Kidus Asfaw

References

Ionides, E. L., Asfaw, K., Park, J., and King, A. A. (2021). Bagged filters for partially observed interacting systems. *Journal of the American Statistical Association*, doi:10.1080/01621459.2021.1974867

See Also

likelihood maximization algorithms: `ienkf()`, `igirf()`, `iubf()`, `ibpf()`

Other likelihood evaluation algorithms: `abf()`, `bpfilter()`, `enkf()`, `girf()`

Examples

```

# Complete examples are provided in the package tests
## Not run:
# Create a simulation of a Brownian motion
b <- bm(U=2, N=5)

# Create a neighborhood function mapping a point in space-time
# to a list of ``neighboring points'' in space-time
bm_nbhd <- function(object, time, unit) {
  nbhd_list = list()
  if(time > 1 && unit > 1){
    nbhd_list = c(nbhd_list, list(c(unit-1, time-1)))
  }
  return(nbhd_list)
}
# Run ABFIR with specified number of Monte Carlo replicates and particles
# per replicate
abfird_bm <- abfir(b,
  Nrep = 2,
  Np=10,
  nbhd = bm_nbhd,
  Ninter = length(unit_names(b)))
# Get the likelihood estimate from ABFIR
logLik(abfird_bm)

## End(Not run)

```

arma_benchmark

Calculated log-ARMA log-likelihood benchmark for spatPomp models

Description

Fits independent log-ARMA models for each unit, and calculates the conditional log-likelihood for each observation, as well as log-likelihood for each unit and total log-likelihood. A simple tool, but one with practical applicability, as demonstrated by King et al (2008) and Wheeler et al (2023). This function is designed for non-negative data, and adds 1 to each observation to avoid $\log(0)$.

Usage

```
arma_benchmark(spo, order = c(2, 0, 1))
```

Arguments

spo A spatPomp object
order A triple (p,d,q) for the ARIMA model fitted to the data. It is intended that d=0

Author(s)

Edward L. Ionides

References

King, A. A., Ionides, E. L., Pascual, M. and Bouma, M. J. (2008). Inapparent infections and cholera dynamics. *Nature* 454 877-880.

Wheeler, J., Rosengart, A. L., Jiang, Z., Tan, K., Treutle, N. and Ionides, E. L. (2023). Informing policy via dynamic models: Cholera in Haiti. *arxiv:2301.08979*.

See Also

Other utilities: [expand_params\(\)](#)

Examples

```
# Complete examples are provided in the package tests
## Not run:
m <- he10(U = 5)
arma_benchmark(m)

## End(Not run)
```

as.data.frame

Coerce to data frame

Description

spatPomp objects can be recast as data frames.

Usage

```
## S3 method for class 'spatPomp'
as.data.frame(x, ...)
```

Arguments

x a spatPomp object.
... additional arguments to be passed to or from methods.

Details

When object is a simple 'spatPomp' object, `as(object, "data.frame")` or `as.data.frame(object)` results in a data frame with the times, units, observables, states (if known), and interpolated covariates (if any).

Value

A 'data.frame' with columns for time, spatial unit and observations.

bm*Brownian motion spatPomp simulator*

Description

Generate a class 'spatPomp' object representing a U-dimensional Brownian motion with spatial correlation decaying geometrically with distance around a circle. The model is defined in continuous time though in this case an Euler approximation is exact at the evaluation times.

Usage

```
bm(U = 5, N = 100, delta_t = 0.1)
```

Arguments

U	A length-one numeric signifying dimension of the process.
N	A length-one numeric signifying the number of observation time steps to evolve the process.
delta_t	Process simulations are performed every delta_t time units whereas observations occur every one time unit

Value

An object of class 'spatPomp' representing a simulation from a U-dimensional Brownian motion

Author(s)

Edward L. Ionides

See Also

Other spatPomp model generators: [bm2\(\)](#), [gbm\(\)](#), [he10\(\)](#), [lorenz\(\)](#), [measles\(\)](#)

Examples

```
# Complete examples are provided in the package tests
## Not run:
b <- bm(U=4, N=20)
# See all the model specifications of the object
spy(b)
# Examples of methodologies applied to this model
# are provided in the tests directory

## End(Not run)
```

bm2	<i>Brownian motion spatPomp generator with shared or unit-specific parameters</i>
-----	---

Description

An extension of `bm` allowing for shared or unit-specific parameters. Generate a class ‘`spatPomp`’ object representing a U -dimensional Brownian motion with spatial correlation decaying geometrically with distance around a circle. The model is defined in continuous time though in this case an Euler approximation is exact at the evaluation times.

Usage

```
bm2(
  U = 5,
  N = 100,
  delta_t = 0.1,
  unit_specific_names = "rho",
  shared_names = NULL,
  unit_params = c(rho = 0.4, sigma = 1, tau = 1, X_0 = 0)
)
```

Arguments

<code>U</code>	A length-one numeric signifying dimension of the process.
<code>N</code>	A length-one numeric signifying the number of observation time steps to evolve the process.
<code>delta_t</code>	Process simulations are performed every <code>delta_t</code> time units whereas observations occur every one time unit
<code>unit_specific_names</code>	determines which parameters take a different value for each unit. Cannot be specified if <code>shared_names</code> is specified. each unit. Other parameters are considered shared between all units.
<code>shared_names</code>	identifies parameters that have common shared value for all units, which by default is all parameters.
<code>unit_params</code>	parameter values used to build the object, copied across each unit for unit-specific parameters

Value

An object of class ‘`spatPomp`’ representing a simulation from a U -dimensional Brownian motion

Author(s)

Edward L. Ionides

See Also

Other spatPomp model generators: [bm\(\)](#), [gbm\(\)](#), [he10\(\)](#), [lorenz\(\)](#), [measles\(\)](#)

Examples

```
# Complete examples are provided in the package tests
## Not run:
b <- bm2(U=4, N=20, shared_names="rho", unit_specific_names=c("sigma", "tau"))
# See all the model specifications of the object
spy(b)
# Examples of methodologies applied to this model
# are provided in the tests directory

## End(Not run)
```

bm2_kalman_logLik	<i>Exact log-likelihood for Brownian motion spatPomp generator with shared or unit-specific parameters</i>
-------------------	--

Description

Computes the exact likelihood for a model constructed using `bm2`, using the Kalman filter. This model is useful for testing methods for models with unit-specific parameters, or method such as `ibpf` which require a unit-specific extension of shared parameters.

Usage

```
bm2_kalman_logLik(bm2_object, params = coef(bm2_object))
```

Arguments

bm2_object	A spatPomp model built using <code>bm2</code> .
params	A parameter vector at which to evaluate the log-likelihood. whereas observations occur every one time unit

Value

A numeric value for the log-likelihood.

Author(s)

Edward L. Ionides

Examples

```
# Further examples are provided in the tests directory
## Not run:
b <- bm2()
bm2_kalman_logLik(b)

## End(Not run)
```

bm_kalman_logLik	<i>Exact log-likelihood for Brownian motion spatPomp generator</i>
------------------	--

Description

Computes the exact likelihood for a model constructed using `bm`, using the Kalman filter. This model is useful for testing methods in a situation where an exact answer is available

Usage

```
bm_kalman_logLik(bm_object, params = coef(bm_object))
```

Arguments

<code>bm_object</code>	A spatPomp model built using <code>bm</code> .
<code>params</code>	A parameter vector at which to evaluate the log-likelihood. whereas observations occur every one time unit

Value

A numeric value for the log-likelihood.

Author(s)

Edward L. Ionides

Examples

```
# Further examples are provided in the tests directory
## Not run:
b <- bm()
bm_kalman_logLik(b)

## End(Not run)
```

 bpfilter

Block particle filter (BPF)

Description

An implementation of the block particle filter algorithm of Rebeschini and van Handel (2015), which is used to estimate the filter distribution of a spatiotemporal partially-observed Markov process. `bpfilter` requires a partition of the spatial units which can be provided by either the `block_size` or the `block_list` argument. The elements of the partition are called blocks. We perform resampling for each block independently based on sample weights within the block. Each resampled block only contains latent states for the spatial components within the block which allows for a “cross-pollination” of particles where the highest weighted segments of each particle are more likely to be resampled and get combined with resampled components of other particles. The method mitigates the curse of dimensionality by resampling locally.

Usage

```
## S4 method for signature 'missing'
bpfilter(object, ...)

## S4 method for signature 'ANY'
bpfilter(object, ...)

## S4 method for signature 'spatPomp'
bpfilter(
  object,
  Np,
  block_size,
  block_list,
  save_states,
  filter_traj,
  ...,
  verbose = getOption("verbose", FALSE)
)

## S4 method for signature 'bpfilterd_spatPomp'
bpfilter(
  object,
  Np,
  block_size,
  block_list,
  save_states,
  filter_traj,
  ...,
  verbose = getOption("verbose", FALSE)
)
```

Arguments

object	A spatPomp object.
...	If a params argument is specified, bpfilter will estimate the likelihood at that parameter set instead of at coef(object).
Np	The number of particles used within each replicate for the adapted simulations.
block_size	The number of spatial units per block. If this is provided, the method subdivides units approximately evenly into blocks with size block_size.
block_list	List that specifies an exact partition of the spatial units. Each partition element, or block, is an integer vector of neighboring units.
save_states	logical. If True, the state-vector for each particle and block is saved.
filter_traj	logical; if TRUE, a filtered trajectory is returned for the state variables and parameters.
verbose	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

Upon successful completion, bpfilter() returns an object of class ‘bpfilterd_spatPomp’ containing the algorithmic parameters used to run bpfilter() and the estimated likelihood.

Details

Only one of block_size or block_list should be specified. If both or neither is provided, an error is triggered.

Methods

The following methods are available for such an object:

[logLik](#) yields an estimate of the log-likelihood of the data under the model.

Author(s)

Kidus Asfaw

References

Rebeschini, P., & Van Handel, R. (2015). Can local particle filters beat the curse of dimensionality?. *The Annals of Applied Probability*, **25(5)**, 2809-2866.

Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package spatPomp. *ArXiv*: 2101.01157. doi:10.48550/arXiv.2101.01157

See Also

likelihood maximization algorithms: [ienkf\(\)](#), [igirf\(\)](#), [iubf\(\)](#), [ibpf\(\)](#)

Other likelihood evaluation algorithms: [abf\(\)](#), [abfir\(\)](#), [enkf\(\)](#), [girf\(\)](#)

Examples

```
# Complete examples are provided in the package tests
## Not run:
# Create a simulation of a Brownian motion
b <- bm(U=4, N=2)

# Run BPF with the specified number of units per block
bpfilterd_b1 <- bpfilter(b, Np = 10, block_size = 2)

# Run BPF with the specified partition
bpfilterd_b2 <- bpfilter(b,
                        Np = 10,
                        block_list = list(c(1,2),c(3,4)) )

# Get a likelihood estimate
logLik(bpfilterd_b2)

## End(Not run)
```

city_data_UK

City data in the United Kingdom

Description

Population and birth information about cities in England and Wales during the measles pre-vaccine era.

Details

Data includes births and population at bi-weekly observations from 40 cities and towns.

Value

a 'data.frame' of the 40 largest cities and towns in the UK and Wales, their latitude, longitude and mean population during the measles pre-vaccine period.

References

Dalziel, Benjamin D. et al. (2016) Persistent chaos of measles epidemics in the prevaccination United States caused by a small change in seasonal transmission patterns. *PLoS Computational Biology*, **12(2)**, e1004655. doi:10.5061/dryad.r4q34

See Also

Other datasets: [measlesUK](#)

concat	<i>Concatenate</i>
--------	--------------------

Description

Concatenate two or more ‘pomp’ objects into a list-like ‘listie’.

Usage

```
## S3 method for class 'SpatPomp'
c(...)
```

Arguments

... elements to be recursively combined into a ‘listie’

Details

concat applied to one or more ‘pomp’ objects or lists of ‘pomp’ objects converts the list into a ‘listie’. In particular, `concat(A,B,C)` is equivalent to `do.call(c,unlist(list(A,B,C)))`.

dunit_measure	<i>dunit_measure</i> dunit_measure evaluates the unit measurement density of a unit's observation given the entire state
---------------	--

Description

dunit_measure dunit_measure evaluates the unit measurement density of a unit's observation given the entire state

Usage

```
## S4 method for signature 'spatPomp'
dunit_measure(object, y, x, unit, time, params, log = TRUE, ...)
```

Arguments

object	An object of class spatPomp
y	A U by 1 matrix of observations for all units
x	A state vector for all units
unit	The unit for which to evaluate the unit measurement density
time	The time for which to evaluate the unit measurement density
params	parameters at which to evaluate the unit measurement density
log	logical; should the density be returned on log scale?
...	additional arguments will be ignored

Value

A class ‘matrix’ with the unit measurement density for spatial unit `unit` corresponding to the corresponding measurement in `y` and states in `x`.

Examples

```
# Complete examples are provided in the package tests
## Not run:
b <- bm(U=3)
s <- states(b)[,1,drop=FALSE]
rownames(s) -> rn
dim(s) <- c(3,1,1)
dimnames(s) <- list(variable=rn, rep=NULL)
p <- coef(b); names(p) -> rnp
dim(p) <- c(length(p),1); dimnames(p) <- list(param=rnp)
o <- obs(b)[,1,drop=FALSE]
dunit_measure(b, y=o, x=s, unit=1, time=1, params=p)

## End(Not run)
```

enkf

Generalized Ensemble Kalman filter (EnKF)

Description

A function to perform filtering using the ensemble Kalman filter of Evensen, G. (1994). This function is generalized to allow for an measurement covariance matrix that varies over time. This is useful if the measurement model varies with the state.

Usage

```
## S4 method for signature 'spatPomp'
enkf(data, Np, ..., verbose = getOption("verbose", FALSE))
```

Arguments

<code>data</code>	A <code>spatPomp</code> object.
<code>Np</code>	The number of Monte Carlo particles used to approximate the filter distribution.
<code>...</code>	Additional arguments can be used to replace model components.
<code>verbose</code>	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

An object of class ‘`enkfd_spatPomp`’ that contains the estimate of the log likelihood (via the `loglik` attribute), algorithmic parameters used to run `enkf()`. Also included are estimated filter means, prediction means and forecasts that are generated during an `enkf()` run.

References

- G. Evensen. Sequential data assimilation with a nonlinear quasi-geostrophic model using Monte Carlo methods to forecast error statistics. *Journal of Geophysical Research: Oceans* **99**, 10143–10162, 1994.
- G. Evensen. *Data assimilation: the ensemble Kalman filter*. Springer-Verlag, 2009.
- J.L. Anderson. An Ensemble Adjustment Kalman Filter for Data Assimilation. *Monthly Weather Review* **129**, 2884–2903, 2001.

See Also

`ienkf()`, `igirf`, `iubf`, `ibpf`

Other likelihood evaluation algorithms: `abf()`, `abfir()`, `bpfilter()`, `girf()`

Examples

```
# Complete examples are provided in the package tests
## Not run:
# Create a simulation of a Brownian motion
b <- bm(U=2, N=5)

# Run EnKF
enkfd_bm <- enkf(b, Np = 20)

# Get a likelihood estimate
logLik(enkfd_bm)

## End(Not run)
```

eunit_measure

eunit_measure

Description

`eunit_measure` evaluates the expectation of a unit's observation given the entire state

Usage

```
## S4 method for signature 'spatPomp'
eunit_measure(object, x, unit, time, params, Np = 1, log = FALSE)
```

Arguments

<code>object</code>	An object of class <code>spatPomp</code>
<code>x</code>	A state vector for all units
<code>unit</code>	The unit for which to evaluate the expectation
<code>time</code>	The time for which to evaluate the expectation

params	parameters at which to evaluate the unit expectation
Np	numeric; defaults to 1 and the user need not change this
log	logical; should the density be returned on log scale?

Value

A class ‘matrix’ with the unit expected observation for spatial unit `unit` corresponding to the corresponding states in `x`.

Examples

```
# Complete examples are provided in the package tests
## Not run:
b <- bm(U=3)
s <- states(b)[,1,drop=FALSE]
rownames(s) -> rn
dim(s) <- c(3,1,1)
dimnames(s) <- list(variable=rn, rep=NULL)
p <- coef(b); names(p) -> rnp
dim(p) <- c(length(p),1); dimnames(p) <- list(param=rnp)
o <- obs(b)[,1,drop=FALSE]
eunit_measure(b, x=s, unit=2, time=1, params=p)

## End(Not run)
```

expand_params

Book-keeping functions for working with expanded parameters

Description

Iterated block particle filters require shared parameters to be expanded into having a value at each unit. `expand_params`, `contract_params` and `mean_by_unit` provide tools for moving between representations. For a unit-specific expansion of a shared parameter, all the values for different units should be the same, and `mean_by_unit` ensures this by taking an average.

Usage

```
expand_params(params, expandedParNames, U)

contract_params(params, expandedParNames, U, average = FALSE)

mean_by_unit(params, expandedParNames, U)
```

Arguments

params	Input parameter vector
expandedParNames	character vector of parameters that are, or should be, expanded. These names should have no numerical suffix 1:U.
U	Number of units
average	Logical value for whether contract_params should average unequal values

Details

These functions assume that expanded parameters have names ending in "1" through "U", where U is the number of units. Contracted parameters, meaning any parameter that is not expanded, should have a name ending in "1". This numerical suffix convention is useful for writing model-building code that allows parameters to be either expanded or contracted.

See Also

Other utilities: [arma_benchmark\(\)](#)

 gbm

Geometric Brownian motion spatPomp simulator

Description

Generate a spatPomp object representing a U-dimensional geometric Brownian motion with spatial correlation decaying geometrically with distance around a circle. The model is defined in continuous time, but an Euler approximation is used for this numerical implementation.

Usage

```
gbm(U = 5, N = 100, delta_t = 0.1, IVP_values = 1, delta_obs = 1)
```

Arguments

U	A length-one numeric signifying dimension of the process.
N	A length-one numeric signifying the number of time steps to evolve the process.
delta_t	process simulations are performed every delta_t time units
IVP_values	initial value parameters for the latent states
delta_obs	observations occur every delta_obs time units

Value

An object of class 'spatPomp' representing a simulation from a U-dimensional geometric Brownian motion

Author(s)

Kidus Asfaw

References

Asfaw, K. T. (2021). Simulation-based Inference for Partially Observed Markov Process Models with Spatial Coupling. University of Michigan Doctoral dissertation. doi:[10.7302/2751](https://doi.org/10.7302/2751)

See Also

Other spatPomp model generators: [bm\(\)](#), [bm2\(\)](#), [he10\(\)](#), [lorenz\(\)](#), [measles\(\)](#)

Examples

```
# Complete examples are provided in the package tests
## Not run:
g <- gbm(U=4, N=20)
# See all the model specifications of the object
spy(g)

## End(Not run)
```

girf

Guided intermediate resampling filter (GIRF)

Description

An implementation of the algorithm of Park and Ionides (2020), following the pseudocode in Asfaw et al. (2020).

Usage

```
## S4 method for signature 'missing'
girf(object, ...)

## S4 method for signature 'ANY'
girf(object, ...)

## S4 method for signature 'spatPomp'
girf(
  object,
  Np,
  Ninter,
  lookahead = 1,
  Nguide,
  kind = c("bootstrap", "moment"),
  tol,
```

```

    ...,
    verbose = getOption("verbose", FALSE)
  )

## S4 method for signature 'girfd_spatPomp'
girf(
  object,
  Np,
  Ninter,
  lookahead,
  Nguide,
  kind = c("bootstrap", "moment"),
  tol,
  ...
)

```

Arguments

object	A spatPomp object.
...	Additional arguments can be used to replace model components.
Np	The number of particles used within each replicate for the adapted simulations.
Ninter	the number of intermediate resampling time points. By default, this is set equal to the number of units.
lookahead	The number of future observations included in the guide function.
Nguide	The number of simulations used to estimate state process uncertainty for each particle.
kind	One of two types of guide function construction. Defaults to 'bootstrap'. See Park and Ionides (2020) for more details.
tol	If all of the guide function evaluations become too small (beyond floating-point precision limits), we set them to this value.
verbose	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

Upon successful completion, `girf()` returns an object of class 'girfd_spatPomp' which contains the algorithmic parameters that were used to run `girf()` and the resulting log likelihood estimate.

Methods

The following methods are available for such an object:

[logLik](#) yields an unbiased estimate of the log-likelihood of the data under the model.

Author(s)

Kidus Asfaw

References

Park, J. and Ionides, E. L. (2020) Inference on high-dimensional implicit dynamic models using a guided intermediate resampling filter. *Statistics and Computing*, doi:10.1007/s11222020099573

Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package spatPomp. *ArXiv*: 2101.01157. doi:10.48550/arXiv.2101.01157

See Also

likelihood maximization algorithms: `ienkf()`, `igirf()`, `iubf()`, `ibpf()`

Other likelihood evaluation algorithms: `abf()`, `abfir()`, `bpfilter()`, `enkf()`

Examples

```
# Complete examples are provided in the package tests
## Not run:
#
# Create a simulation of a Brownian motion
b <- bm(U=2, N=5)

# Run GIRF
girfd_bm <- girf(b,
                 Np = 10,
                 Ninter = length(unit_names(b)),
                 lookahead = 1,
                 Nguides = 10
                )
# Get the likelihood estimate from GIRF
logLik(girfd_bm)

# Compare with the likelihood estimate from particle filter
pfd_bm <- pfilter(b, Np = 10)
logLik(pfd_bm)

## End(Not run)
```

he10

Measles in UK: spatPomp generator with shared or unit-specific parameters

Description

Generate a spatPomp object for measles adding spatial coupling to The model and data from He et al. (2010) with gravity transport as in Park and Ionides (2020). Other transport models may be added in future. The data in the object matches He et al. (2010). The model matches that analysis in the specific case where there is no coupling and all parameters are unit-specific.

Usage

```

he10(
  U = 6,
  dt = 2/365,
  Tmax = 1964,
  expandedParNames = c("alpha", "iota", "R0", "cohort", "amplitude", "gamma", "sigma",
    "sigmaSE", "rho", "psi", "g", "S_0", "E_0", "I_0"),
  basic_params = c(alpha = 1, iota = 0, R0 = 30, cohort = 0, amplitude = 0.5, gamma = 52,
    sigma = 52, mu = 0.02, sigmaSE = 0.15, rho = 0.5, psi = 0.15, g = 400, S_0 = 0.032,
    E_0 = 5e-05, I_0 = 4e-05),
  towns_selected = NULL
)

```

Arguments

U	A length-one numeric signifying the number of cities to be represented in the spatPomp object. Default U=20 gives all the towns studied by He et al., the 10 largest and 10 selected smaller towns.
dt	a numeric (in unit of years) that is used as the Euler time-increment for simulating measles data.
Tmax	Upper time for the window used to construct the object. The lower time is fixed at 1950.0. The default value matches He et al (2010).
expandedParNames	specifies the names of parameters which take unit-specific values. Remaining parameters take a single, shared value for all units.
basic_params	A candidate parameter vector in the basic format, i.e., no unit-specific parameters or unit-related name extensions.
towns_selected	A numeric vector of towns to be modeled. Defaults to 1:U, with cities ranked by decreasing population and 1 being London.

Details

The code for this spatPomp has duplication with measles(), but in future the two models may diverge. The measles() spatPomp is a simplified situation useful for testing some methods. However, measles() does not permit unit-specific parameters, which he10() allows. Also, the structure of this spatPomp is compatible with the spatiotemporal iterated filtering algorithm ibpf(). This requires shared parameters to be represented with a value for each unit, which should be the same for each unit in a valid model instance but may vary between units while optimizing.

Value

An object of class 'spatPomp' representing a U-dimensional spatially coupled measles POMP model.

Relationship to published analysis

The model generator he10() differs from measles() in some details necessitated to reproduce the results of He et al (2010). The measles() model follows the decision of Park and Ionides (2020)

and Ionides et al (2021) to apply the mixing exponent α_u to (I_u/P_u) rather than just to I_u . `he10()` does this for the infections arising from individuals traveling to another town (which don't arise for the panel model of He et al (2010)). However, for infections arising within a city, in order to reproduce the results of He et al (2010), `he10()` uses $(I_u^{\alpha_u}/P_u)$. This is not fully documented in the text of Ionides et al (2022). Models fitted to data have α_u close to 1, so this issue may be negligible in practice.

Another discrepancy between the `he10()` code and the mathematical model written by Ionides et al (2022) arises in whether individuals traveling from u to v use mixing exponent α_u or α_v . Ionides et al (2022) wrote u but the code used implemented v . The implementation in `he10()` matches the implementation of Ionides et al (2022) and so uses v .

It might seem surprising that immigrant infections affect only the first term in the expression for μ_{SE} in Ionides et al (2022), and in the corresponding `he10()` code. This immigration term is needed in the first term to make the model of He et al (2010) a proper sub-model, when coupling is removed by setting the gravitational constant parameter equal to zero. When this constant is allowed to be positive, the role of immigrant infections transmitting to traveling individuals is anticipated to be a negligible, second-order effect which has been omitted from the model.

Note

This function goes through a typical workflow of constructing a typical `spatPomp` object (1-4 below). This allows the user to have a file that replicates the exercise of model building as well as function that creates a typical nonlinear model in epidemiology in case they want to test a new inference methodology. We purposely do not modularize this function because it is not an operational piece of the package and is instead useful as an example.

1. Getting a measurements `data.frame` with columns for times, spatial units and measurements.
2. Getting a covariates `data.frame` with columns for times, spatial units and covariate data.
3. Constructing model components (latent state initializer, latent state transition simulator and measurement model). Depending on the methods used, the user may have to supply a vectorfield to be integrated that represents the deterministic skeleton of the latent process.
4. Bringing all the data and model components together to form a `spatPomp` object via a call to `spatPomp()`.

Author(s)

Edward L. Ionides

References

- Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package `spatPomp`. *ArXiv*: 2101.01157. [doi:10.48550/arXiv.2101.01157](https://doi.org/10.48550/arXiv.2101.01157)
- He, D., Ionides, E. L., and King, A. A. (2010). Plug-and-play inference for disease dynamics: measles in large and small populations as a case study. *Journal of the Royal Society Interface*, 7(43), 271-283. [doi:10.1098/rsif.2009.0151](https://doi.org/10.1098/rsif.2009.0151)
- Ionides, E. L., Asfaw, K., Park, J., and King, A. A. (2021). Bagged filters for partially observed interacting systems. *Journal of the American Statistical Association*, [doi:10.1080/01621459.2021.1974867](https://doi.org/10.1080/01621459.2021.1974867)

Ionides, E. L., Ning, N., and Wheeler, J. (2022). An iterated block particle filter for inference on coupled dynamic systems with shared and unit-specific parameters. *Statistica Sinica*, to appear. doi:10.48550/arXiv.2206.03837

Park, J. and Ionides, E. L. (2020) Inference on high-dimensional implicit dynamic models using a guided intermediate resampling filter. *Statistics and Computing*, doi:10.1007/s11222020099573

See Also

he10coordinates, he10measles, he10mle, he10demography

Other spatPomp model generators: [bm\(\)](#), [bm2\(\)](#), [gbm\(\)](#), [lorenz\(\)](#), [measles\(\)](#)

Examples

```
# Complete examples are provided in the package tests
## Not run:
m <- he10(U = 5)
# See all the model specifications of the object
spy(m)

## End(Not run)
```

he10coordinates	<i>City data in the United Kingdom</i>
-----------------	--

Description

Longitude and latitude for the 20 towns in England and Wales studied by He et al (2010).

Value

a ‘data.frame’ of longitude and latitude for each town.

References

He, D., Ionides, E. L., and King, A. A. (2010). Plug-and-play inference for disease dynamics: measles in large and small populations as a case study. *Journal of the Royal Society Interface*, 7(43), 271-283. doi:10.1098/rsif.2009.0151

See Also

Other datasets he10: [he10demography](#), [he10measles](#), [he10mle](#)

he10demography

Demographic data for 20 towns in the United Kingdom

Description

Population and birth information for some towns in England and Wales during the measles pre-vaccine era.

Details

Data are annual statistics for the 20 towns analyzed by He et al (2010).

Value

a 'data.frame' of with variables town, year, pop and births.

References

He, D., Ionides, E. L., and King, A. A. (2010). Plug-and-play inference for disease dynamics: measles in large and small populations as a case study. *Journal of the Royal Society Interface*, 7(43), 271-283. doi:10.1098/rsif.2009.0151

See Also

Other datasets he10: [he10coordinates](#), [he10measles](#), [he10mle](#)

he10measles

Measles in the United Kingdom

Description

Measles case data from various cities and towns in England and Wales during the pre-vaccine era.

Details

Data are weekly case counts for the 20 towns analyzed by He et al (2010).

Value

a 'data.frame' of reported measles cases for 20 towns, analyzed by He et al (2010).

References

He, D., Ionides, E. L., and King, A. A. (2010). Plug-and-play inference for disease dynamics: measles in large and small populations as a case study. *Journal of the Royal Society Interface*, 7(43), 271-283. doi:10.1098/rsif.2009.0151

See Also

Other datasets he10: [he10coordinates](#), [he10demography](#), [he10mle](#)

 he10mle

Measles in the United Kingdom: MLE from He et al (2010)

Description

Maximum likelihood estimate for fitting a susceptible-exposed-infected- recovered model to the measles case report data analyzed by He et al (2010). The values are similar, but not identical, to those reported by He et al.

Value

a 'data.frame' containing the estimated parameters.

References

He, D., Ionides, E. L., and King, A. A. (2010). Plug-and-play inference for disease dynamics: measles in large and small populations as a case study. *Journal of the Royal Society Interface*, 7(43), 271-283. doi:10.1098/rsif.2009.0151

See Also

Other datasets he10: [he10coordinates](#), [he10demography](#), [he10measles](#)

 ibpf

Iterated block particle filter (IBPF)

Description

An iterated block particle filter, for both shared and unit-specific parameters. We require that the spatPomp has been constructed to have a unit-specific parameter "thetau" for unit u corresponding to an estimated parameter "theta", whether theta is shared or unit-specific. This permits IBPF to implement a spatiotemporal random walk to estimate theta. We require that rw.sd is positive for, and only for, all parameters of the form "thetau" if "theta" is listed in sharedParNames or unitParNames.

Usage

```
## S4 method for signature 'missing'
ibpf(data, ...)

## S4 method for signature 'ANY'
ibpf(data, ...)

## S4 method for signature 'spatPomp'
ibpf(
  data,
  Nbpf,
  Np,
  rw.sd,
  sharedParNames,
  unitParNames,
  cooling.type = "geometric",
  cooling.fraction.50,
  block_size,
  block_list,
  spat_regression,
  ...,
  verbose = getOption("verbose", FALSE)
)

## S4 method for signature 'ibpfd_spatPomp'
ibpf(
  data,
  Nbpf,
  Np,
  rw.sd,
  sharedParNames,
  unitParNames,
  cooling.type = "geometric",
  cooling.fraction.50,
  block_size,
  block_list,
  spat_regression,
  ...,
  verbose = getOption("verbose", FALSE)
)

## S4 method for signature 'bpfilterd_spatPomp'
ibpf(
  data,
  Nbpf,
  Np,
  rw.sd,
  sharedParNames,
```

```

    unitParNames,
    cooling.type = "geometric",
    cooling.fraction.50,
    block_size,
    block_list,
    spat_regression,
    ...,
    verbose = getOption("verbose", FALSE)
)

```

Arguments

<code>data</code>	either a data frame holding the time series data, or an object of class ‘pomp’, i.e., the output of another pomp calculation. Internally, <code>data</code> will be coerced to an array with storage-mode double.
<code>...</code>	If a <code>params</code> argument is specified, <code>bpfilter</code> will estimate the likelihood at that parameter set instead of at <code>coef(object)</code> .
<code>Nbpf</code>	the number of iterations of perturbed BPF.
<code>Np</code>	The number of particles used within each replicate for the adapted simulations.
<code>rw.sd</code>	specification of the magnitude of the random-walk perturbations that will be applied to some or all model parameters. Parameters that are to be estimated should have positive perturbations specified here. The specification is given using the <code>rw.sd</code> function, which creates a list of unevaluated expressions. The latter are evaluated in a context where the model time variable is defined (as time). The expression <code>ivp(s)</code> can be used in this context as shorthand for <code>ifelse(time==time[1],s,0)</code> . Likewise, <code>ivp(s,lag)</code> is equivalent to <code>ifelse(time==time[lag],s,0)</code> . See below for some examples. The perturbations that are applied are normally distributed with the specified s.d. If parameter transformations have been supplied, then the perturbations are applied on the transformed (estimation) scale.
<code>sharedParNames</code>	estimated parameters that are equal for each unit.
<code>unitParNames</code>	estimated parameters that are different for each unit.
<code>cooling.type, cooling.fraction.50</code>	specifications for the cooling schedule, i.e., the manner and rate with which the intensity of the parameter perturbations is reduced with successive filtering iterations. <code>cooling.type</code> specifies the nature of the cooling schedule. See below (under “Specifying the perturbations”) for more detail.
<code>block_size</code>	The number of spatial units per block. If this is provided, the method subdivides units approximately evenly into blocks with size <code>block_size</code> .
<code>block_list</code>	List that specifies an exact partition of the spatial units. Each partition element, or block, is an integer vector of neighboring units.

spat_regression	fraction of each extended parameter regressed toward the unit mean. Not required when all parameters are unit-specific.
verbose	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

Upon successful completion, `ibpf` returns an object of class `'ibpfd_spatPomp'`.

Methods

The following methods are available for such an object:

`coef` gives the Monte Carlo estimate of the maximum likelihood.

Author(s)

Edward L. Ionides

References

Ionides, E. L., Ning, N., and Wheeler, J. (2022). An iterated block particle filter for inference on coupled dynamic systems with shared and unit-specific parameters. *Statistica Sinica*, to appear. [doi:10.48550/arXiv.2206.03837](https://doi.org/10.48550/arXiv.2206.03837)

See Also

likelihood evaluation algorithms: `girf()`, `enkf()`, `bpfilter()`, `abf()`, `abfir()`

Other likelihood maximization algorithms: `ienkf()`, `igirf()`, `iubf()`

Examples

```
# Complete examples are provided in the package tests
## Not run:
# Create a simulation of a Brownian motion, for an extended model with
# unit-specific parameters for all estimated, even if the parameter
# takes the same shared value for each unit.
U <- 4
b2 <- bm2(U=U,N=5,unit_specific_names="rho")

# Run ibpf with two blocks of two units each. estimating rho as a
# shared parameter with all other parameters being fixed.
b2_ibpf <- ibpf(b2,
  sharedParNames="rho",
  unitParNames=NULL,
  Nbpf=5,
  spat_regression=0.1,
  Np=50,
  rw.sd=do.call(rw_sd,setNames(rep(list(0.01),times=U),paste0("rho",1:U))),
  cooling.fraction.50=0.5,
```



```

    block_size=2
  )

  # Get a likelihood estimate
  logLik(b2_ibpf)

  ## End(Not run)

```

 ienkf

Iterated ensemble Kalman filter (IEnKF)

Description

An implementation of a parameter estimation algorithm that uses the ensemble Kalman filter (Evensen, G. (1994)) to perform the filtering step in the parameter-perturbed iterated filtering scheme of Ionides et al. (2015) following the pseudocode in Asfaw, et al. (2020).

Usage

```

## S4 method for signature 'spatPomp'
ienkf(
  data,
  Nenkf = 1,
  rw.sd,
  cooling.type = c("geometric", "hyperbolic"),
  cooling.fraction.50,
  Np,
  ...,
  verbose = getOption("verbose", FALSE)
)

```

Arguments

<code>data</code>	an object of class <code>spatPomp</code>
<code>Nenkf</code>	number of iterations of perturbed EnKF.
<code>rw.sd</code>	specification of the magnitude of the random-walk perturbations that will be applied to some or all model parameters. Parameters that are to be estimated should have positive perturbations specified here. The specification is given using the <code>rw_sd</code> function, which creates a list of unevaluated expressions. The latter are evaluated in a context where the model time variable is defined (as <code>time</code>). The expression <code>ivp(s)</code> can be used in this context as shorthand for <code>ifelse(time==time[1],s,0)</code> . Likewise, <code>ivp(s,lag)</code> is equivalent to <code>ifelse(time==time[lag],s,0)</code> .

See below for some examples.

The perturbations that are applied are normally distributed with the specified s.d. If parameter transformations have been supplied, then the perturbations are applied on the transformed (estimation) scale.

<code>cooling.type</code> , <code>cooling.fraction</code>	<code>.50</code> specifications for the cooling schedule, i.e., the manner and rate with which the intensity of the parameter perturbations is reduced with successive filtering iterations. <code>cooling.type</code> specifies the nature of the cooling schedule. See below (under “Specifying the perturbations”) for more detail.
<code>Np</code>	The number of particles used within each replicate for the adapted simulations.
<code>...</code>	Additional arguments can be used to replace model components.
<code>verbose</code>	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

Upon successful completion, `ienkf` returns an object of class ‘`ienkfd_spatPomp`’. This object contains the convergence record of the iterative algorithm with respect to the likelihood and the parameters of the model (which can be accessed using the `traces` attribute) as well as a final parameter estimate, which can be accessed using the `coef()`.

Methods

The following methods are available for such an object:

`coef` gives the Monte Carlo estimate of the maximum likelihood.

Author(s)

Kidus Asfaw

References

- Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package `spatPomp`. *ArXiv*: 2101.01157. doi:10.48550/arXiv.2101.01157
- Evensen, G. (1994) Sequential data assimilation with a nonlinear quasi-geostrophic model using Monte Carlo methods to forecast error statistics *Journal of Geophysical Research: Oceans* 99:10143–10162
- Evensen, G. (2009) *Data assimilation: the ensemble Kalman filter* Springer-Verlag.
- Anderson, J. L. (2001) An Ensemble Adjustment Kalman Filter for Data Assimilation *Monthly Weather Review* 129:2884–2903

See Also

likelihood evaluation algorithms: `girf()`, `enkf()`, `bpfilter()`, `abf()`, `abfir()`
Other likelihood maximization algorithms: `ibpf()`, `igirf()`, `iubf()`

igirf	<i>Iterated guided intermediate resampling filter (IGIRF)</i>
-------	---

Description

An implementation of a parameter estimation algorithm combining the intermediate resampling scheme of the guided intermediate resampling filter of Park and Ionides (2020) and the parameter perturbation scheme of Ionides et al. (2015) following the pseudocode in Asfaw, et al. (2020).

Usage

```
## S4 method for signature 'missing'
igirf(data, ...)

## S4 method for signature 'ANY'
igirf(data, ...)

## S4 method for signature 'spatPomp'
igirf(
  data,
  Ngirf,
  Np,
  rw.sd,
  cooling.type,
  cooling.fraction.50,
  Ninter,
  lookahead = 1,
  Nguide,
  kind = c("bootstrap", "moment"),
  tol = 1e-100,
  ...,
  verbose = getOption("verbose", FALSE)
)

## S4 method for signature 'igirfd_spatPomp'
igirf(
  data,
  Ngirf,
  Np,
  rw.sd,
  cooling.type,
  cooling.fraction.50,
  Ninter,
  lookahead,
  Nguide,
  kind = c("bootstrap", "moment"),
  tol,
```

```

    ...,
    verbose = getOption("verbose", FALSE)
  )

```

Arguments

<code>data</code>	an object of class <code>spatPomp</code> or <code>igirfd_spatPomp</code>
<code>...</code>	Additional arguments can be used to replace model components.
<code>Ngirf</code>	the number of iterations of parameter-perturbed GIRF.
<code>Np</code>	The number of particles used within each replicate for the adapted simulations.
<code>rw.sd</code>	specification of the magnitude of the random-walk perturbations that will be applied to some or all model parameters. Parameters that are to be estimated should have positive perturbations specified here. The specification is given using the <code>rw_sd</code> function, which creates a list of unevaluated expressions. The latter are evaluated in a context where the model time variable is defined (as <code>time</code>). The expression <code>ivp(s)</code> can be used in this context as shorthand for <code>ifelse(time==time[1],s,0)</code> . Likewise, <code>ivp(s,lag)</code> is equivalent to <code>ifelse(time==time[lag],s,0)</code> . See below for some examples. The perturbations that are applied are normally distributed with the specified s.d. If parameter transformations have been supplied, then the perturbations are applied on the transformed (estimation) scale.
<code>cooling.type, cooling.fraction.50</code>	specifications for the cooling schedule, i.e., the manner and rate with which the intensity of the parameter perturbations is reduced with successive filtering iterations. <code>cooling.type</code> specifies the nature of the cooling schedule. See below (under “Specifying the perturbations”) for more detail.
<code>Ninter</code>	the number of intermediate resampling time points. By default, this is set equal to the number of units.
<code>lookahead</code>	The number of future observations included in the guide function.
<code>Nguide</code>	The number of simulations used to estimate state process uncertainty for each particle.
<code>kind</code>	One of two types of guide function construction. Defaults to 'bootstrap'. See Park and Ionides (2020) for more details.
<code>tol</code>	If all of the guide function evaluations become too small (beyond floating-point precision limits), we set them to this value.
<code>verbose</code>	logical; if TRUE, messages updating the user on progress will be printed to the console.

Value

Upon successful completion, `igirf()` returns an object of class `'igirfd_spatPomp'`. This object contains the convergence record of the iterative algorithm with respect to the likelihood and the parameters of the model (which can be accessed using the `traces` attribute) as well as a final parameter estimate, which can be accessed using the `coef()`. The algorithmic parameters used to run `igirf()` are also included.

Methods

The following methods are available for such an object:

`coef` gives the Monte Carlo maximum likelihood parameter estimate.

Author(s)

Kidus Asfaw

References

Park, J. and Ionides, E. L. (2020) Inference on high-dimensional implicit dynamic models using a guided intermediate resampling filter. *Statistics and Computing*, doi:10.1007/s11222020099573

Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package `spatPomp`. *ArXiv*: 2101.01157. doi:10.48550/arXiv.2101.01157

See Also

likelihood evaluation algorithms: `girf()`, `enkf()`, `bpfilter()`, `abf()`, `abfir()`

Other likelihood maximization algorithms: `ibpf()`, `ienkf()`, `iubf()`

Examples

```
# Complete examples are provided in the package tests
## Not run:
igirf(bm(U=2,N=4),Ngirf=2,
      rw.sd = rw_sd(rho=0.02,X1_0=ivp(0.02)),
      cooling.type="geometric",cooling.fraction.50=0.5,
      Np=10,Ninter=2,lookahead=1,Nguid=5)

## End(Not run)
```

iubf

*Iterated Unadapted Bagged Filter (IUBF)***Description**

An algorithm for estimating the parameters of a spatiotemporal partially-observed Markov process. Running `iubf` causes the algorithm to perform a specified number of iterations of unadapted simulations with parameter perturbation and parameter resamplings. At each iteration, unadapted simulations are performed on a perturbed version of the model, in which the parameters to be estimated are subjected to random perturbations at each observation. After cycling through the data, each replicate's weight is calculated and is used to rank the bootstrap replicates. The highest ranking replicates are recycled into the next iteration. This extra variability introduced through parameter perturbation effectively smooths the likelihood surface and combats particle depletion by introducing diversity into particle population. As the iterations progress, the magnitude of the perturbations is diminished according to a user-specified cooling schedule.

Usage

```
## S4 method for signature 'spatPomp'
iubf(
  object,
  Nubf = 1,
  Nrep_per_param,
  Nparam,
  nbhd,
  prop,
  rw.sd,
  cooling.type = c("geometric", "hyperbolic"),
  cooling.fraction.50,
  tol = (1e-18)^17,
  verbose = getOption("verbose"),
  ...
)
```

Arguments

<code>object</code>	A <code>spatPomp</code> object.
<code>Nubf</code>	The number of iterations to perform
<code>Nrep_per_param</code>	The number of replicates used to estimate the likelihood at a parameter
<code>Nparam</code>	The number of parameters that will undergo the iterated perturbation
<code>nbhd</code>	A neighborhood function with three arguments: <code>object</code> , <code>time</code> and <code>unit</code> . The function should return a list of two-element vectors that represent space-time neighbors of (u, n) , which is represented by <code>c(unit, time)</code> . See example below for more details.
<code>prop</code>	A numeric between 0 and 1. The top <code>prop*100%</code> of the parameters are resampled at each observation

<code>rw.sd</code>	<p>specification of the magnitude of the random-walk perturbations that will be applied to some or all model parameters. Parameters that are to be estimated should have positive perturbations specified here. The specification is given using the <code>rw_sd</code> function, which creates a list of unevaluated expressions. The latter are evaluated in a context where the model time variable is defined (as <code>time</code>). The expression <code>ivp(s)</code> can be used in this context as shorthand for <code>ifelse(time==time[1],s,0)</code>.</p> <p>Likewise, <code>ivp(s,lag)</code> is equivalent to <code>ifelse(time==time[lag],s,0)</code>.</p> <p>See below for some examples.</p> <p>The perturbations that are applied are normally distributed with the specified s.d. If parameter transformations have been supplied, then the perturbations are applied on the transformed (estimation) scale.</p>
<code>cooling.type, cooling.fraction.50</code>	<p>specifications for the cooling schedule, i.e., the manner and rate with which the intensity of the parameter perturbations is reduced with successive filtering iterations. <code>cooling.type</code> specifies the nature of the cooling schedule. See below (under “Specifying the perturbations”) for more detail.</p>
<code>tol</code>	<p>If the resampling weight for a particle is zero due to floating-point precision issues, it is set to the value of <code>tol</code> since resampling has to be done.</p>
<code>verbose</code>	<p>logical; if TRUE, diagnostic messages will be printed to the console.</p>
<code>...</code>	<p>additional arguments are passed to <code>pomp</code>.</p>

Value

Upon successful completion, `iubf()` returns an object of class ‘`iubfd_spatPomp`’. This object contains the convergence record of the iterative algorithm with respect to the likelihood and the parameters of the model (which can be accessed using the `traces` attribute) as well as a final parameter estimate, which can be accessed using the `coef()`. The algorithmic parameters used to run `iubf()` are also included.

Methods

The following methods are available for such an object:

`coef` extracts the point estimate

Author(s)

Kidus Asfaw

References

- Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package `spatPomp`. *ArXiv*: 2101.01157. doi:10.48550/arXiv.2101.01157
- Ionides, E. L., Asfaw, K., Park, J., and King, A. A. (2021). Bagged filters for partially observed interacting systems. *Journal of the American Statistical Association*, doi:10.1080/01621459.2021.1974867

See Also

likelihood evaluation algorithms: `girf()`, `enkf()`, `bpfilter()`, `abf()`, `abfir()`

Other likelihood maximization algorithms: `ibpf()`, `ienkf()`, `igirf()`

`logLik`*Log likelihood*

Description

Extract the estimated log likelihood.

Usage

```
## S4 method for signature 'girfd_spatPomp'  
logLik(object)  
  
## S4 method for signature 'bpfilterd_spatPomp'  
logLik(object)  
  
## S4 method for signature 'abfd_spatPomp'  
logLik(object)  
  
## S4 method for signature 'iubfd_spatPomp'  
logLik(object)  
  
## S4 method for signature 'abfird_spatPomp'  
logLik(object)  
  
## S4 method for signature 'igirfd_spatPomp'  
logLik(object)
```

Arguments

`object` fitted model object

Value

a numeric which is the estimated log likelihood

lorenz	<i>Lorenz '96 spatPomp simulator</i>
--------	--------------------------------------

Description

Generate a spatPomp object representing a U-dimensional stochastic Lorenz '96 process with N measurements made at times $t_n = n * \text{delta_obs}$, simulated using an Euler method with time increment `delta_t`.

Usage

```
lorenz(
  U = 5,
  N = 100,
  delta_t = 0.01,
  delta_obs = 0.5,
  regular_params = c(F = 8, sigma = 1, tau = 1)
)
```

Arguments

U	A length-one numeric signifying the number of spatial units for the process.
N	A length-one numeric signifying the number of observations.
delta_t	A length-one numeric giving the Euler time step for the numerical solution.
delta_obs	A length-one numeric giving the time between observations.
regular_params	A named numeric vector containing the values of the F, sigma and tau parameters. F=8 is a common value that causes chaotic behavior.

Value

An object of class 'spatPomp' representing a simulation from a U-dimensional Lorenz 96 model

Author(s)

Edward L. Ionides

References

Lorenz, E. N. (1996) Predictability: A problem partly solved. *Proceedings of the seminar on predictability*

Ionides, E. L., Asfaw, K., Park, J., and King, A. A. (2021). Bagged filters for partially observed interacting systems. *Journal of the American Statistical Association*, doi:[10.1080/01621459.2021.1974867](https://doi.org/10.1080/01621459.2021.1974867)

See Also

Other spatPomp model generators: `bm()`, `bm2()`, `gbm()`, `he10()`, `measles()`

Examples

```
# Complete examples are provided in the package tests
## Not run:
l <- lorenz(U=5, N=100, delta_t=0.01, delta_obs=1)
# See all the model specifications of the object
spy(l)

## End(Not run)
```

measles

Measles in UK spatPomp generator

Description

Generate a spatPomp object for measles in the top-U most populous cities in England and Wales. The model is adapted from He et al. (2010) with gravity transport following Park and Ionides (2020). The data are from Dalziel et al (2016).

Usage

```
measles(
  U = 6,
  dt = 2/365,
  fixed_ivps = TRUE,
  S_0 = 0.032,
  E_0 = 5e-05,
  I_0 = 4e-05
)
```

Arguments

U	A length-one numeric signifying the number of cities to be represented in the spatPomp object.
dt	a numeric (in unit of years) that is used as the Euler time-increment for simulating measles data.
fixed_ivps	a logical. If TRUE initial value parameters will be declared in the globals slot, shared for each unit, and will not be part of the parameter vector.
S_0	a numeric. If fixed_ivps=TRUE this is the initial proportion of all of the spatial units that are susceptible.
E_0	a numeric. If fixed_ivps=TRUE this is the initial proportion of all of the spatial units that are exposed.
I_0	a numeric. If fixed_ivps=TRUE this is the initial proportion of all of the spatial units that are infected.

Value

An object of class 'spatPomp' representing a U-dimensional spatially coupled measles POMP model.

Relationship to published analysis

This model was used to generate the results of Ionides et al (2021). However, their equation (6) is not exactly correct for the Binomial Gamma infinitesimal model used in the code, as shown by Proposition 5 of Breto and Ionides, 2011. If Poisson Gamma infinitesimal increments were used (Proposition 4 of Breto and Ionides, 2011) then (6) would be correct, but the resulting unbounded increments could break the non-negativity requirement for compartment membership. The same issue arises with the description in Park and Ionides (2020), though that analysis was based on a different model implementation since the spatPomp package was not yet available.

A difference between (6) of Ionides et al (2021) and (2.1) of He et al (2010) is that in (6) the mixing exponent α is applied to (I_u/P_u) rather than just to I_u . In the context of He et al (2010) this changes the parameterization but has negligible effect on the model itself since $P_u(t)$ is approximately constant and so changing its power can be compensated by a corresponding change in the transmission rate, β . In practice, models fitted to data have *alpha* close to 1, so this issue may be moot and this modeling mechanism may not be an effective empirical way to carry out the goal of making allowance for heterogeneous mixing.

The code here includes a cohort effect, c , following He et al (2010), that was not included by Ionides et al (2021). This effect leads to a non-differentiability of expected increments which is problematic for the spatPomp implementation of GIRF. For the results of Ionides et al (2021), this was set to $c = 0$.

The analysis of He et al (2010), and the model generated by `he10()`, use weekly aggregated cases. Weekly reports were not available beyond the 20 cites studied by He et al (2010) so `measles()` relies on the biweekly reports used by Ionides et al (2021) and Ionides & Park (2020).

It turns out to be an important detail of the model by He et al (2010) that a delay is included between birth and entry into the susceptible compartment. He et al (2010) found a 4 year delay fits the data. This value is fixed to be the variable `birth_delay` in the code for `measles()`. The code for Ionides et al (2021) uses a 3 year delay, and the delay is not explained in the abbreviated model description. In `measles()` we have reverted to the 4 year delay identified by He et al (2010).

Note

This function goes through a typical workflow of constructing a typical spatPomp object (1-4 below). This allows the user to have a file that replicates the exercise of model building as well as function that creates a typical nonlinear model in epidemiology in case they want to test a new inference methodology. We purposely do not modularize this function because it is not an operational piece of the package and is instead useful as an example.

1. Getting a measurements `data.frame` with columns for times, spatial units and measurements.
2. Getting a covariates `data.frame` with columns for times, spatial units and covariate data.
3. Constructing model components (latent state initializer, latent state transition simulator and measurement model). Depending on the methods used, the user may have to supply a vectorfield to be integrated that represents the deterministic skeleton of the latent process.
4. Bringing all the data and model components together to form a spatPomp object via a call to `spatPomp()`.

Author(s)

Edward L. Ionides

References

- Ionides, E. L., Asfaw, K., Park, J., and King, A. A. (2021). Bagged filters for partially observed interacting systems. *Journal of the American Statistical Association*, doi:10.1080/01621459.2021.1974867
- Dalziel, Benjamin D. et al. (2016) Persistent chaos of measles epidemics in the prevaccination United States caused by a small change in seasonal transmission patterns. *PLoS Computational Biology*, **12**(2), e1004655. doi:10.5061/dryad.r4q34
- Park, J. and Ionides, E. L. (2020) Inference on high-dimensional implicit dynamic models using a guided intermediate resampling filter. *Statistics and Computing*, doi:10.1007/s11222020099573
- Breto, C. and Ionides, E.L. (2011) Compound Markov counting processes and their applications to modeling infinitesimally over-dispersed systems. *Stochastic Processes and their Applications* **121**, 2571-2591. doi:10.1016/j.spa.2011.07.005

See Also

measles_UK, city_data_UK

Other spatPomp model generators: [bm\(\)](#), [bm2\(\)](#), [gbm\(\)](#), [he10\(\)](#), [lorenz\(\)](#)

Examples

```
# Complete examples are provided in the package tests
## Not run:
m <- measles(U = 5)
# See all the model specifications of the object
spy(m)

## End(Not run)
```

measles2

Measles in UK: spatPomp generator with shared or unit-specific parameters

Description

Generate a spatPomp object for measles in the top-U most populous cities in England and Wales. The model is adapted from He et al. (2010) with gravity transport following Park and Ionides (2019). The structure of this spatPomp is designed to accommodate shared and unit-specific parameters. If carrying out spatiotemporal iterated filtering for shared parameters via `ibpf`, it is necessary to have a unit-specific expansion and so these parameters should be included in `expandedParNames`. This model and data correspond to the biweekly analysis of Park and Ionides (2020) and Ionides et al (2021). There are small differences with the weekly model and data of He et al (2010) and Ionides, Ning and Wheeler (2022).

Usage

```
measles2(
  U = 6,
  dt = 2/365,
  N = 391,
  expandedParNames = c("R0", "c", "A", "muIR", "muEI", "sigmaSE", "rho", "psi", "g",
    "S_0", "E_0", "I_0"),
  contractedParNames = NULL,
  simulated = FALSE,
  basic_params = c(alpha = 0.98, iota = 0.1, R0 = 30, c = 0.3, A = 0.5, muIR = 52, muEI =
    52, muD = 0.02, sigmaSE = 0.15, rho = 0.5, psi = 0.15, g = 400, S_0 = 0.032, E_0 =
    5e-05, I_0 = 4e-05)
)
```

Arguments

U	An integer from 1 to 40 specifying the number of cities to be represented in the spatPomp object.
dt	a numeric (in unit of years) that is used as the Euler time-increment for simulating measles data
N	An integer from 1 to 391 specifying the number of time points.
expandedParNames	specifies parameters that are defined for each unit. This also allows unit perturbations for a parameter with a value shared across units.
contractedParNames	specifies parameters having a shared value across units. Remaining parameters that are neither expanded nor contracted are considered fixed, and will not have a transformation defined for them.
simulated	determines whether to return a simulation from the model or the UK measles data
basic_params	A named vector used to specify shared parameters or unit-specific parameters having common values for each unit.

Value

An object of class 'spatPomp' representing a U-dimensional spatially coupled measles POMP model.

References

He, D., Ionides, E. L., and King, A. A. (2010). Plug-and-play inference for disease dynamics: measles in large and small populations as a case study. *Journal of the Royal Society Interface*, 7(43), 271-283. doi:10.1098/rsif.2009.0151

Park, J. and Ionides, E. L. (2020) Inference on high-dimensional implicit dynamic models using a guided intermediate resampling filter. *Statistics and Computing*, doi:10.1007/s11222020099573

Ionides, E. L., Asfaw, K., Park, J., and King, A. A. (2021). Bagged filters for partially observed interacting systems. *Journal of the American Statistical Association*, doi:10.1080/01621459.2021.1974867

Ionides, E. L., Ning, N., and Wheeler, J. (2022). An iterated block particle filter for inference on coupled dynamic systems with shared and unit-specific parameters. *Statistica Sinica*, to appear. [doi:10.48550/arXiv.2206.03837](https://doi.org/10.48550/arXiv.2206.03837)

Examples

```
# Complete examples are provided in the package tests
## Not run:
m <- measles2(U = 5)
# See all the model specifications of the object
spy(m)

## End(Not run)
```

measlesUK

Measles in the United Kingdom

Description

Measles case data from various cities and towns in England and Wales during the pre-vaccine era.

Details

Data includes bi-weekly case counts as well as births and population from 40 cities and towns.

Value

a ‘data.frame’ of the 40 largest cities and towns in the UK and Wales, their latitude, longitude and bi-weekly measles case counts, population and birthrates.

References

Dalziel, Benjamin D. et al. (2016) Persistent chaos of measles epidemics in the prevaccination United States caused by a small change in seasonal transmission patterns. *PLoS Computational Biology*, **12**(2), e1004655. [doi:10.5061/dryad.r4q34](https://doi.org/10.5061/dryad.r4q34)

See Also

Other datasets: [city_data_UK](#)

munit_measure	<i>munit_measure</i>
---------------	----------------------

Description

munit_measure returns a moment-matched parameter set given an empirically calculated measurement variance and latent states. This is used in girf() and igirf() when they are run with kind='moment'.

Usage

```
## S4 method for signature 'spatPomp'
munit_measure(object, x, vc, unit, time, params, Np = 1)
```

Arguments

object	An object of class spatPomp
x	A state vector for all units
vc	The empirically calculated variance used to perform moment-matching
unit	The unit for which to obtain a moment-matched parameter set
time	The time for which to obtain a moment-matched parameter set
params	parameters to use to obtain a moment-matched parameter set
Np	Number of particle replicates for which to get parameter sets

Value

An array with dimensions `dim(array.params)[1]` by `dim(x)[2]` by `length(unit)` by `length(time)` representing the moment-matched parameter set(s) corresponding to the variance of the measurements, vc, and the states, x.

Author(s)

Kidus Asfaw

Examples

```
# Complete examples are provided in the package tests
## Not run:
b <- bm(U=3)
s <- states(b)[,1,drop=FALSE]
rownames(s) -> rn
dim(s) <- c(3,1,1)
dimnames(s) <- list(variable=rn, rep=NULL)
p <- coef(b); names(p) -> rnp
dim(p) <- c(length(p),1); dimnames(p) <- list(param=rnp)
o <- obs(b)[,1,drop=FALSE]
array.params <- array(p,
```

```

        dim = c(length(p),
                length(unit_names(b)), 1, 1),
        dimnames = list(params = rownames(p)))
vc <- c(4, 9, 16); dim(vc) <- c(length(vc), 1, 1)
munit_measure(b, x=s, vc=vc, Np=1, unit = 1, time=1, params=array.params)

## End(Not run)

```

plot

Plotting spatPomp data

Description

Visualize spatPomp data
 Diagnostic plot for `igirf()`
 Visualize spatPomp data

Usage

```

## S4 method for signature 'igirfd_spatPomp'
plot(x, params = names(coef(x)), ncol = 3)

## S4 method for signature 'spatPomp'
plot(x, type = c("l", "h"), log = F, plot_unit_names = T, ...)

```

Arguments

<code>x</code>	a <code>spatPomp</code> object
<code>params</code>	the names of the parameters for which the user would like to see a trace plot
<code>ncol</code>	the number of columns in the grid plot
<code>type</code>	for visualizing an object of class <code>spatPomp</code> , the user can obtain a grid of line plots by default ('l') or a heat map by supplying argument 'h'.
<code>log</code>	should the data be log-transformed before plotting? This helps in contexts where there are spikes that could take away attention from the dynamics illustrated by the rest of the data.
<code>plot_unit_names</code>	allows suppression of unit names when making a heat map for a large number of units
<code>...</code>	for visualizing an object of class <code>spatPomp</code> , the user can add arguments like <code>nrow</code> to specify the number of rows in the grid.

Value

a ggplot facet plot of class 'gg' and 'ggplot' visualizing the convergence record of running `igirf()` with respect to the likelihood and the parameters of the model.

a ggplot plot of class 'gg' and 'ggplot' visualizing the time series data over multiple spatial units via a tile-plot.

print	<i>Print methods</i>
-------	----------------------

Description

Prints its argument.

Usage

```
## S4 method for signature 'spatPomp'
print(x)
```

Arguments

x a spatPomp object

Value

An object of class 'spatPomp' is returned **invisibly**. The user is notified on the console only the class of the object.

Note

Use `spy()` to see model components of x instead.

runit_measure	<i>runit_measure</i>
---------------	----------------------

Description

`runit_measure` simulates a unit's observation given the entire state

Usage

```
## S4 method for signature 'spatPomp'
runit_measure(object, x, unit, time, params, log = FALSE)
```

Arguments

object	An object of class spatPomp
x	A state vector for all units
unit	The unit for which to simulate an observation
time	The time for which to simulate an observation
params	parameters to use to simulate an observation
log	logical; should the density be returned on log scale?

Value

A matrix with the simulated observation corresponding to state `x` and unit `unit` with parameter set `params`.

Author(s)

Kidus Asfaw

Examples

```
# Complete examples are provided in the package tests
## Not run:
b <- bm(U=3)
s <- states(b)[,1,drop=FALSE]
rownames(s) -> rn
dim(s) <- c(3,1,1)
dimnames(s) <- list(variable=rn, rep=NULL)
p <- coef(b); names(p) -> rnp
dim(p) <- c(length(p),1); dimnames(p) <- list(param=rnp)
o <- obs(b)[,1,drop=FALSE]
runit_measure(b, x=s, unit=2, time=1, params=p)

## End(Not run)
```

simulate

Simulation of a spatiotemporal partially-observed Markov process

Description

simulate generates simulations of the latent and measurement processes.

Usage

```
## S4 method for signature 'spatPomp'
simulate(
  object,
  nsim = 1,
  seed = NULL,
  format = c("spatPomps", "data.frame"),
  include.data = FALSE,
  ...
)
```

Arguments

object	optional; if present, it should be a data frame or a 'pomp' object.
nsim	number of simulations.
seed	optional integer; if set, the pseudorandom number generator (RNG) will be initialized with seed. The RNG will be restored to its original state afterward.
format	the format of the simulated results. If the argument is set to 'spatPomps', the default behavior, then the output is a list of spatPomp objects. Options are 'spatPomps' and 'data.frame'.
include.data	if TRUE, the original data and covariates (if any) are included (with .id = "data"). This option is ignored unless format = "data.frame".
...	additional arguments are passed to pomp .

Value

if format='spatPomps' and nsim=1 an object of class 'spatPomp' representing a simulation from the model in object is returned. If format='spatPomps' and nsim>1 a list of class 'spatPomp' objects is returned. If format='data.frame' then a class 'data.frame' object is returned.

Author(s)

Kidus Asfaw

References

Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package spatPomp. *ArXiv*: 2101.01157. doi:[10.48550/arXiv.2101.01157](https://doi.org/10.48550/arXiv.2101.01157)

Examples

```
# Complete examples are provided in the package tests
## Not run:
# Get a spatPomp object
b <- bm(U=2, N=5)
# Get 2 simulations from same model as data.frame
sims <- simulate(b, nsim=2, format='data.frame')
```

```
## End(Not run)
```

```
spatPomp
```

```
Constructor of the spatPomp object
```

Description

This function constructs a class ‘spatPomp’ object, encoding a spatiotemporal partially observed Markov process (SPATPOMP) model together with a uni- or multi-variate time series on a collection of units. Users will typically develop a POMP model for a single unit before embarking on a coupled SpatPOMP analysis. Consequently, we assume some familiarity with **pomp** and its description by King, Nguyen and Ionides (2016). The spatPomp class inherits from pomp with the additional unit structure being a defining feature of the resulting models and inference algorithms.

Usage

```
spatPomp(
  data,
  units,
  times,
  covar,
  t0,
  ...,
  eunit_measure,
  munit_measure,
  vunit_measure,
  dunit_measure,
  runit_measure,
  rprocess,
  rmeasure,
  dprocess,
  dmeasure,
  skeleton,
  rinit,
  rprior,
  dprior,
  unit_statenames,
  unit_accumvars,
  shared_covarnames,
  globals,
  paramnames,
  params,
  cdir,
  cfile,
  shlib.args,
  PACKAGE,
```

```

    partrans,
    compile = TRUE,
    verbose = getOption("verbose", FALSE)
)

```

Arguments

<code>data</code>	either a dataframe holding the spatiotemporal data, or an object of class ‘ <code>spatPomp</code> ’, i.e., the output of another spatPomp calculation. If dataframe, the user must provide the name of the times column using the <code>times</code> argument and the spatial unit column name using the <code>units</code> argument. The dataframe provided should be sorted in increasing order of time and unit name respectively, i.e. observation 1 in unit A should come before observation 1 in unit B, which should come before observation 2 in unit A.
<code>units</code>	when <code>data</code> is a <code>data.frame</code> this is the name of the column containing the spatial units.
<code>times</code>	the sequence of observation times. <code>times</code> must indicate the column of observation times by name or index. The time vector must be numeric and non-decreasing.
<code>covar</code>	An optional dataframe for supplying covariate information. If provided, there must be two columns that provide the observation time and the observation spatial unit with the same names and arrangement as the <code>data</code> .
<code>t0</code>	The zero-time, i.e., the time of the initial state. This must be no later than the time of the first observation, i.e., <code>t0 <= times[1]</code> .
<code>...</code>	If there are arguments that the user would like to pass to pomp ’s basic constructor function’s <code>...</code> argument, this argument passes them along. Not recommended for this version of spatPomp .
<code>eunit_measure</code>	Evaluator of the expected measurement given the latent states and model parameters. The <code>unit</code> variable is pre-defined, which allows the user to specify differing specifications for each unit using <code>if</code> conditions. Only <code>C</code> snippets are accepted. The <code>C</code> snippet should assign the scalar approximation to the expected measurement to the pre-defined variable <code>ey</code> given the latent state and the parameters. For more information, see the examples section below.
<code>munit_measure</code>	Evaluator of a moment-matched parameter set (like the standard deviation parameter of a normal distribution or the size parameter of a negative binomial distribution) given an empirical variance estimate, the latent states and all model parameters. Only <code>C</code> snippets are accepted. The <code>C</code> snippet should assign the scalar approximation to the measurement variance parameter to the pre-defined variable corresponding to that parameter, which has been predefined with a <code>M_</code> prefix. For instance, if the moment-matched parameter is <code>psi</code> , then the user should assign <code>M_psi</code> to the moment-matched value. For more information, see the examples section below.
<code>vunit_measure</code>	Evaluator of the theoretical measurement variance given the latent states and model parameters. The <code>unit</code> variable is pre-defined, which allows the user to specify differing specifications for each unit using <code>if</code> conditions. Only <code>C</code> snippets are accepted. The <code>C</code> snippet should assign the scalar approximation to the

	measurement variance to the pre-defined variable <code>vc</code> given the latent state and the parameters. For more information, see the examples section below.
<code>dunit_measure</code>	Evaluator of the unit measurement model density given the measurement, the latent states and model parameters. The <code>unit</code> variable is pre-defined, which allows the user to specify differing specifications for each unit using <code>if</code> conditions. Only Csnippets are accepted. The Csnippet should assign the scalar measurement density to the pre-defined variable <code>lik</code> . The user is encouraged to provide a logged density in an <code>if</code> condition that checks whether the pre-defined <code>give_log</code> variable is true. For more information, see the examples section below.
<code>runit_measure</code>	Simulator of the unit measurement model given the latent states and the model parameters. The <code>unit</code> variable is pre-defined, which allows the user to specify differing specifications for each unit using <code>if</code> conditions. Only Csnippets are accepted. The Csnippet should assign the scalar measurement density to the pre-defined which corresponds to the name of the observation for each unit (e.g. <code>cases</code> for the measles spatPomp example). For more information, see the examples section below.
<code>rprocess</code>	simulator of the latent state process, specified using one of the rprocess plugins . Setting <code>rprocess=NULL</code> removes the latent-state simulator. For more information, see rprocess specification for the documentation on these plugins .
<code>rmeasure</code>	simulator of the measurement model, specified either as a C snippet, an R function, or the name of a pre-compiled native routine available in a dynamically loaded library. Setting <code>rmeasure=NULL</code> removes the measurement model simulator. For more information, see rmeasure specification .
<code>dprocess</code>	evaluator of the probability density of transitions of the unobserved state process. Setting <code>dprocess=NULL</code> removes the latent-state density evaluator. For more information, see dprocess specification .
<code>dmeasure</code>	evaluator of the measurement model density, specified either as a C snippet, an R function, or the name of a pre-compiled native routine available in a dynamically loaded library. Setting <code>dmeasure=NULL</code> removes the measurement density evaluator. For more information, see dmeasure specification .
<code>skeleton</code>	optional; the deterministic skeleton of the unobserved state process. Depending on whether the model operates in continuous or discrete time, this is either a vectorfield or a map. Accordingly, this is supplied using either the vectorfield or map fnctions. For more information, see skeleton specification . Setting <code>skeleton=NULL</code> removes the deterministic skeleton.
<code>rinit</code>	simulator of the initial-state distribution. This can be furnished either as a C snippet, an R function, or the name of a pre-compiled native routine available in a dynamically loaded library. Setting <code>rinit=NULL</code> sets the initial-state simulator to its default. For more information, see rinit specification .
<code>rprior</code>	optional; prior distribution sampler, specified either as a C snippet, an R function, or the name of a pre-compiled native routine available in a dynamically loaded library. For more information, see prior specification . Setting <code>rprior=NULL</code> removes the prior distribution sampler.
<code>dprior</code>	optional; prior distribution density evaluator, specified either as a C snippet, an R function, or the name of a pre-compiled native routine available in a dynamically loaded library. For more information, see prior specification . Setting

	dprior=NULL resets the prior distribution to its default, which is a flat improper prior.
unit_statenames	The names of the components of the latent state. E.g. if the user is constructing an joint SIR model over many spatial units, <code>c('S', 'I', 'R')</code> would be passed.
unit_accumvars	a subset of the <code>unit_statenames</code> argument that are accumulator variables. See accumvars for more on the concept of pomp accumulator variables.
shared_covarnames	If <code>covar</code> is supplied, covariates that are shared must still be specified for each unit, i.e., rows with equal values for the same time over all units must be supplied. However, if such covariates exists, supply the names using this argument.
globals	optional character or C snippet; arbitrary C code that will be hard-coded into the shared-object library created when C snippets are provided. If no C snippets are used, <code>globals</code> has no effect.
paramnames	optional character vector; names of model parameters. It is typically only necessary to supply <code>paramnames</code> when C snippets are in use.
params	optional; named numeric vector of parameters. This will be coerced internally to storage mode <code>double</code> .
cdir	optional character variable. <code>cdir</code> specifies the name of the directory within which C snippet code will be compiled. By default, this is in a temporary directory specific to the R session. One can also set this directory using the <code>pomp_cdir</code> global option.
cfile	optional character variable. <code>cfile</code> gives the name of the file (in directory <code>cdir</code>) into which C snippet codes will be written. By default, a random filename is used. If the chosen filename would result in over-writing an existing file, an error is generated.
shlib.args	optional character variables. Command-line arguments to the R CMD SHLIB call that compiles the C snippets. One can, for example, specify libraries against which the C snippets are to be linked. In doing so, take care to make sure the appropriate header files are available to the C snippets, e.g., using the <code>globals</code> argument. See Csnippet for more information.
PACKAGE	optional character; the name (without extension) of the external, dynamically loaded library in which any native routines are to be found. This is only useful if one or more of the model components has been specified using a precompiled dynamically loaded library; it is not used for any component specified using C snippets. <code>PACKAGE</code> can name at most one library.
partrans	optional parameter transformations, constructed using parameter_trans . Many algorithms for parameter estimation search an unconstrained space of parameters. When working with such an algorithm and a model for which the parameters are constrained, it can be useful to transform parameters. One should supply the <code>partrans</code> argument via a call to parameter_trans . For more information, see parameter_trans . Setting <code>partrans=NULL</code> removes the parameter transformations, i.e., sets them to the identity transformation.
compile	logical; if <code>FALSE</code> , compilation of the C snippets will be postponed until they are needed.
verbose	logical; if <code>TRUE</code> , diagnostic messages will be printed to the console.

Details

One implements a SPATPOMP model by specifying some or all of its *basic components*, including:

rinit, the simulator from the distribution of the latent state process at the zero-time;

rprocess, the transition simulator of the latent state process;

dunit_measure, the evaluator of the conditional density at a unit's measurement given the unit's latent state;

eunit_measure, the evaluator of the expectation of a unit's measurement given the unit's latent state;

munit_measure, the evaluator of the moment-matched parameter set given a unit's latent state and some empirical measurement variance;

vunit_measure, the evaluator of the variance of a unit's measurement given the unit's latent state;

runit_measure, the simulator of a unit's measurement conditional on the unit's latent state;

dprocess, the evaluator of the density for transitions of the latent state process;

rmeasure, the simulator of the measurements conditional on the latent state;

dmeasure, the evaluator of the conditional density of the measurements given the latent state;

rprior, the simulator from a prior distribution on the parameters;

dprior, the evaluator of the prior density;

skeleton, which computes the deterministic skeleton of the unobserved state process;

partrans, which performs parameter transformations.

The basic structure and its rationale are described in Asfaw et al. (2020).

Each basic component is supplied via an argument of the same name to `spatPomp()`. The five unit-level model components must be provided via C snippets. The remaining components, whose behaviors are inherited from **pomp** may be furnished using C snippets, R functions, or pre-compiled native routine available in user-provided dynamically loaded libraries.

Value

An object of class 'spatPomp' representing observations and model components from the spatiotemporal POMP model.

Author(s)

Kidus Asfaw, Edward L. Ionides, Aaron A. King

References

Asfaw, K., Park, J., Ho, A., King, A. A., and Ionides, E. L. (2020) Partially observed Markov processes with spatial structure via the R package spatPomp. *ArXiv*: 2101.01157. doi:10.48550/arXiv.2101.01157

King, A. A., Nguyen, D. and Ionides, E. L. (2016) Statistical Inference for Partially Observed Markov Processes via the R Package pomp. *Journal of Statistical Software*, **69**(12), 1–43. doi:10.18637/jss.v069.i12

spatPomp-class	<i>An S4 class to represent a spatiotemporal POMP model and data.</i>
----------------	---

Description

An S4 class to represent a spatiotemporal POMP model and data.

Slots

`unit_names` A vector containing the spatial units of a spatiotemporal POMP.

`unit_statenames` A vector containing the state names such that appending the unit indices to the unit statenames will result in the each unit's corresponding states.

`unit_obsnames` A vector of observation types for a spatial unit.

`eunit_measure` A `pomp_fun` representing the expected measurement for each spatial unit given its states.

`dunit_measure` A `pomp_fun` representing the unit measurement density for each spatial unit.

`runit_measure` A `pomp_fun` representing the unit observation simulator.

spatPomp_Csnippet	<i>C snippets</i>
-------------------	-------------------

Description

`spatPomp_Csnippet()` is used to provide snippets of C code that specify model components. It functions similarly to `Csnippet()` from the **pomp** package; in fact, the output of `spatPomp_Csnippet` is an object of class `Csnippet`. It additionally provides some arguments that allow the user to stay focused on model development in the spatiotemporal context where model size grows.

Usage

```
## S4 method for signature 'character'
spatPomp_Csnippet(
  code,
  method = "",
  unit_statenames,
  unit_obsnames,
  unit_covarnames,
  unit_ivpnames,
  unit_paramnames,
  unit_vfnames
)
```

Arguments

code	encodes a component of a spatiotemporal POMP model using C code
method	a character string matching the name of the 'spatPomp' argument which the code is designed to specify. This argument is ignored unless needed to correctly specify the Csnippet.
unit_statenames	a subset of the unit_statenames slot of the spatPomp object for which we are writing a model. This argument allows the user to get variables that can be indexed conveniently to update states and measurements in a loop. See examples for more details.
unit_obsnames	a subset of the unit_obsnames slot of the spatPomp object for which we are writing a model. This argument allows the user to get variables that can be indexed conveniently to update states and measurements in a loop. See examples for more details.
unit_covarnames	if the model has covariate information for each unit, the names of the covariates for each unit can be supplied to this argument. This allows the user to get variables that can be indexed conveniently to use incorporate the covariate information in a loop. See examples for more details.
unit_ivpnames	This argument is particularly useful when specifying the rinit model component. The paramnames argument to the spatPomp() constructor often has names for initial value parameters for the latent states (e.g. S1_0, S2_0 for the the quantity of susceptibles at unit 1 and unit 2 at the initial time in an SIR model). By supplying unit_ivpnames, we can get variables that can be easily indexed to reference the initial value parameters (in the previous example, unit_ivpnames=c('S') we can get a variable named S_0 that we can index as S_0[0] and S_0[1] to refer to S1_0 and S2_0). See examples for more details.
unit_paramnames	This argument is particularly useful when there are non-initial value parameters that are unit-specific.
unit_vfnames	This argument is particularly useful when specifying the skeleton model component. For all components of the latent state, the user can assume a variable defining the time-derivative is pre-defined (e.g. DS1 and DS2 for the time-derivative of the quantity of the susceptibles at unit 1 and unit 2 in an SIR model). By supplying unit_vfnames, we can get variables that can be easily indexed to reference these variables (in the previous example, setting unit_vfnames=c('S') gets us a variable named DS that we can index as DS[0] and DS[1] to refer to DS1 and DS2). See examples for more details.

Value

An object of class 'Csnippet' which represents a model specification in C code.

Author(s)

Kidus Asfaw

Examples

```

# Set initial states for Brownian motion
bm_rinit <- spatPomp_Csnippet(
  method = "rinit",
  unit_statenames = c("X"),
  unit_ivpnames = c("X"),
  code = "
    for (int u = 0; u < U; u++) {
      X[u]=X_0[u];
    }
"
)
# Skeleton for Brownian motion
bm_skel <- spatPomp_Csnippet(
  method = "skeleton",
  unit_statenames = c("X"),
  unit_vfnames = c("X"),
  code = "
    for (int u = 0 ; u < U ; u++) {
      DX[u] = 0;
    }
"
)

```

unit_names

Unit names of a spatiotemporal model

Description

unit_names outputs the contents of the unit_names slot of a spatPomp object. The order in which the units appear in the output vector determines the order in which latent states and observations for the spatial units are stored.

Usage

```

## S4 method for signature 'spatPomp'
unit_names(x)

```

Arguments

x a spatPomp object

Value

A character vector with the unit names used to create the 'spatPomp' object.

vec_dmeasure *Vector of measurement densities*

Description

Evaluate the unit measurement model density function for each unit. This method is used primarily as part of likelihood evaluation and parameter inference algorithms.

Usage

```
## S4 method for signature 'spatPomp'
vec_dmeasure(object, y, x, units, times, params, log = FALSE, ...)
```

Arguments

object	a spatPomp object
y	numeric; measurements whose densities given the latent states are evaluated
x	numeric; state at which conditional measurement densities are evaluated
units	numeric; units at which measurement densities are evaluated
times	numeric; time at which measurement densities are evaluated
params	numeric; parameter set at which measurement densities is evaluated
log	logical; should the outputted measurement densities be on log scale?
...	additional parameters will be ignored

Value

An array of dimension `length(unit_names(object))` by `dim(x)[2]` by `dim(x)[3]` representing each unit's measurement density assessed for each replicate in `x` for each observation time.

Author(s)

Kidus Asfaw

vec_rmeasure *Vector of simulated measurements*

Description

Simulate from the unit measurement model density function for each unit

Usage

```
## S4 method for signature 'spatPomp'
vec_rmeasure(object, x, times, params, ...)
```

Arguments

object	a spatPomp object
x	numeric; state at which measurements are simulated
times	numeric; time at which measurements are simulated
params	numeric; parameter set at which measurements are simulated
...	additional parameters will be ignored

Value

An array of dimension `length(unit_names(object))` by `dim(x)[2]` by `dim(x)[3]` representing each unit's simulated measurement assessed for each replicate in `x` for each observation time.

Author(s)

Kidus Asfaw

vunit_measure	<i>vunit_measure</i>
---------------	----------------------

Description

`vunit_measure` evaluates the variance of a unit's observation given the entire state

Usage

```
## S4 method for signature 'spatPomp'
vunit_measure(object, x, unit, time, params, Np = 1)
```

Arguments

object	An object of class spatPomp
x	A state vector for all units
unit	The unit for which to evaluate the variance
time	The time for which to evaluate the variance
params	parameters at which to evaluate the unit variance
Np	numeric; defaults to 1 and the user need not change this

Value

A matrix with the unit measurement variance implied by the state, `x`, and the parameter set `params` for unit `unit`.

Author(s)

Kidus Asfaw

Examples

```
# Complete examples are provided in the package tests
## Not run:
b <- bm(U=3)
s <- states(b)[,1,drop=FALSE]
rownames(s) -> rn
dim(s) <- c(3,1,1)
dimnames(s) <- list(variable=rn, rep=NULL)
p <- coef(b); names(p) -> rnp
dim(p) <- c(length(p),1); dimnames(p) <- list(param=rnp)
o <- obs(b)[,1,drop=FALSE]
vunit_measure(b, x=s, unit=2, time=1, params=p)

## End(Not run)
```

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