
SDE Documentation

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Contents

1	Layout	1
2	Method	3
3	Location of the documentation	5
3.1	SDE	5
3.2	Module Schauder	6
3.3	Module Diffusion	9
3.4	Module Randm	11
4	Indices and tables	13

The main module

SDE Simulate diffusion processes in one or more dimension. Especially simulate vector linear processes / Ornstein-Uhlenbeck processes Monte Carlo sample diffusion bridges, diffusion processes conditioned to hit a point v at a prescribed time T Functions for transition density, mean and covariance of linear processes Perform Monte Carlo estimates of transition densities of general diffusion processes

with a Submodule

SDE.Schauder To nonparametrically estimate the drift of a diffusion with unit diffusion coefficient using a Schauder wavelet “basis”

The package contains the additional modules:

Diffusion Alternative API to generate Ito processes and diffusions

Randm Random symmetric, positive definite, stable matrix for testing purposes.

CHAPTER 2

Method

The procedure in `SDE.Schauder` is a Julia implementation of nonparametric Bayesian inference for “continuously” observed one dimensional diffusion processes with unit diffusion coefficient. The drift is modeled as linear combination of hierarchical Faber–Schauder basis functions with a Gaussian prior on the coefficients. This incorporates a Brownian motion like prior on the drift function. The posterior is then computed using Gaussian conjugacy.

This is work in progress.

Location of the documentation

<https://sdejl.readthedocs.org>

Contents:

SDE

Miscellaneous

SDE.**syl** (a, b, c)

Solves the Sylvester equation $AX + XB = C$, where C is symmetric and A and $-B$ have no common eigenvalues using (inefficient) algebraic approach via the Kronecker product, see http://en.wikipedia.org/wiki/Sylvester_equation

Stochastic Processes

SDE.**mu** (t, x, T, P)

Expectation $E(t, x)(X_T)$

SDE.**K** (t, T, P)

Covariance matrix $Cov(X_T - x_t)$

SDE.**H** (t, T, P)

Negative Hessian of $\log p(t, x; T, v)$ as a function of x .

SDE.**r** (t, x, T, v, P)

Returns $r(t, x) = \text{grad}_x \log p(t, x; T, v)$ where p is the transition density of the process P .

SDE.**bstar** ($t, x, T, v, P::MvPro$)

Returns the drift function of a vector linear process bridge which end at time T in point v .

SDE.**bcirc** ($t, x, T, v, Pt::Union(MvLinPro, MvAffPro), P::MvPro$)

Drift for guided proposal derived from a vector linear process bridge which end at time T in point v .

SDE.lp (t, x, T, y, P)
Returns $\log p(t, x; T, y)$, the log transition density of the process P

SDE.samplep (t, x, T, P)
Samples from the transition density of the process P .

SDE.exact (u, tt, P)
Simulate process P starting in u on a discrete grid tt from its transition probability.

SDE.ll (X, P)
Compute log likelihood evaluated in B , β and Lyapunov matrix λ for a observed linear process on a discrete grid dt from its transition density.

SDE.lp (s, x, t, y, P)
Returns $\log p(t, x; T, y)$, the log transition density

SDE.euler ($u, W::CTPath, P::CTPro$)
Multivariate euler scheme for U , starting in u using the same time grid as the underlying Wiener process W .

SDE.llikeliXcirc ($t, T, Xcirc, b, a, B, \beta, \lambda$)
Loglikelihood (log weights) of $Xcirc$ with respect to $Xstar$.

 t, T – timespan $Xcirc$ – bridge proposal (drift $Bcirc$ and diffusion coefficient σ) b, σ – diffusion coefficient σ target B, β – drift $b(x) = Bx + \beta$ of $Xtilde$ λ – solution of the lyapunov equation for $Xtilde$

SDE.tofs ($s, tmin, T$)
SDE.soft ($t, tmin, T$)
Time change mapping s in $[0, T=t_2 - t_1]$ (U-time) to t in $[t_1, t_2]$ (X-time), and inverse.

SDE.Vs (s, T, v, B, β)
SDE.dotVs (s, T, v, B, β)
Time changed V and time changed time derivative of V for generation of U

SDE.XofU ($UU, tmin, T, v, P$)
 U is the scaled and time changed process

 $U(s) = \exp(s/2) * (v(s) - X(tofs(s)))$

 $XofU$ transforms entire process U sampled at time points ss to X at tt .

SDE.stable (Y, d, ep)
Return real stable d -dim matrix with real eigenvalues smaller than $-ep$ parametrized with a vector of length $d*d$,

For maximum likelihood estimation we need to search the maximum over all stable matrices. These are matrices with eigenvalues with strictly negative real parts. We obtain a $d \times d$ stable matrix as difference of a antisymmetric matrix and a positive definite matrix.

Module Schauder

Introduction

In the following $\hat{x}$ is the piecewise linear function taking values $(0, 0)$, $(0.5, 1)$, $(1, 0)$ on the interval $[0, 1]$ and 0 elsewhere.

The Schauder basis of level $L > 0$ in the interval $[a, b]$ can be defined recursively from $n = 2^{L-1}$ classical finite elements $\psi_i(x)$ on the grid

$$a + (1:n) / (n+1) * (b-a).$$

Assume that f is expressed as linear combination

$$f(x) = \sum_{i=1}^n c_i \psi_i(x)$$

with

$$\psi_{2j-1}(x) = \text{hat}(nx - j + 1) \text{ for } j = 1 \dots 2^{L-1}$$

and

$$\psi_{2j}(x) = \text{hat}(nx - j + 1/2) \text{ for } j = 1 \dots 2^{L-1} - 1$$

Note that these coefficients are easy to find for the finite element basis, just

```
function fe_transf(f, a, b, L)
    n = 2^L-1
    return map(f, a + (1:n) / (n+1) * (b-a))
end
```

Then the coefficients of the same function with respect to the Schauder basis

$$f(x) = \sum_{i=1}^n c_i \phi_i(x)$$

where for $L = 2$

$$\phi_2(x) = 2\text{hat}(x)$$

$$\phi_1(x) = \psi_1(x) = \text{hat}(2x)$$

$$\phi_3(x) = \psi_3(x) = \text{hat}(2x - 1)$$

can be computed directly, but also using the recursion

$$\phi_2(x) = \psi_1(x) + 2\psi_2(x) + \psi_2(x)$$

This can be implemented inplace (see `pickup()`), and is used throughout.

```
for l in 1:L-1
    b = sub(c, 2^l:2^l:n)
    b[:] *= 0.5
    a = sub(c, 2^(l-1):2^l:n-2^(l-1))
    a[:] -= b[:]
    a = sub(c, 2^l+2^(l-1):2^l:n)
    a[:] -= b[1:length(a)]
end
```

Reference

`Schauder.pickup!(x)`

Inplace computation of 2^L-1 Schauder-Faber coefficients from 2^L-1 overlapping finite-element coefficients x .

– inverse of `Schauder.drop`

– $L = \text{level}(x)$

`Schauder.drop!(x)`

Inplace computation of 2^L-1 finite element coefficients from 2^L-1 Faber schauder coefficients x .

– inverse of `Schauder.pickup`

`Schauder.finger_permute(x)`

Reorders vector x or matrix A according to the reordering of the elements of a Faber-Schauder-basis from left to right, from bottom to top.

`Schauder.finger_pm(L, K)`

Returns the permutation used in `finger_permute`. Memoized reordering of faber schauder elements from low level to high level. The last K elements/rows are left untouched.

`Schauder.level(x)`

Gives the no. of levels of the biggest Schauder basis with less then length(x) elements. `level(x) = ilogb(size(x,1)+1)`

`Schauder.level(x, K)`

Gives the no. of levels `l` of the biggest Schauder basis with less then `length(x)` elements and the number of additional elements `n-2^l+1`.

`Schauder.vectoroflevels(L, K)`

Gives a vector with the level of the hierarchical elements.

`Schauder.hat(x)`

Hat function. Piecewise linear functions with values $(-\infty, 0)$, $(0, 0)$, $(0.5, 1)$, $(1, 0)$, $(\infty, 0)$. - x vector or number

Introduction

The procedure is as follows. Consider the diffusion process $(x_t: 0 \leq t \leq T)$ given by

$$dx_t = b(x_t)dt + dw_t$$

where the drift b is expressed as linear combination

$$f(x) = \sum_{i=1}^n c_i \phi_i(x)$$

(see [Module Schauder](#)) and prior distribution on the coefficients

$$c_i \sim N(0, \xi_i)$$

Then the posterior distribution of b given observations x_t is given by

$$c_i | x_s \sim N(W^{-1}\mu, W^{-1}) \quad W = \Sigma + (\text{diag}(\xi))^{-1},$$

with the $n \times n$ -matrix

$$\Sigma_{ij} = \int_0^T \phi_i(x_t) \phi_j(x_t) dt$$

and the n -vector

$$\mu_i = \int_0^T \phi_i(x_t) dx_t.$$

Using the recursion detailed in [Module Schauder](#), one rather computes

$$\Sigma'_{ij} = \int_0^T \psi_i(x_t) \psi_j(x_t) dt$$

and the n -vector

$$\mu'_i = \int_0^T \psi_i(x_t) dx_t$$

and uses `pickup_mu!(mu)` and `pickup_Sigma!(Sigma)` to obtain μ and Σ .

Optional additional basis functions

One can extend the basis by additional functions, implemented are variants. B1 includes a constant, B2 two linear functions

B1 $\phi_1 \dots \phi_n, c$

B2 $\phi_1 \dots \phi_n, \max(1-x, 0), \max(x, 0)$

To compute μ , use

```
mu = pickup_mu!(fe_mu(y, L, 0))
mu = fe_muB1(mu, y);
```

or

```
mu = pickup_mu!(fe_mu(y, L, 0))
mu = fe_muB2(mu, y);
```

Reference

Functions taking y without parameter $[a, b]$ expect y to be shifted into the interval $[0, 1]$.

Schauder.pickup_mu!(μ)
computes μ from μ'

Schauder.drop_mu!(μ)
Computes μ' from μ .

Schauder.pickup_Sigma!(Σ)
Transforms Σ' into Σ .

Schauder.drop_Sigma!(Σ)
Transforms Σ into Σ' .

Schauder.fe_mu(y, L, K)
Computes μ' from the observations y using 2^{L-1} basis elements and returns a vector with K trailing zeros (in case one ones to customize the basis).

Schauder.fe_muB1(μ, y)
Append $\mu_{n+1} = \int_0^T \phi_{n+1} dx_t$ with $\phi_{n+1} = 1$.

Schauder.fe_muB2(μ, y)
Append $\mu_{n+1} = \int_0^T \phi_{n+1} dx_t$ with $\phi_{n+1} = \max(1 - x, 0)$ and $\mu_{n+2} = \int_0^T \phi_{n+2} dx_t$ with $\phi_{n+2} = \max(x, 0)$

Schauder.fe_Sigma(y, dt, L)
Computes the matrix Σ' from the observations y uniformly spaced at distance dt using 2^{L-1} basis elements.

Schauder.bayes_drift($x, dt, a, b, L, xirem, beta, B$)
Performs estimation of drift on observations x in $[a, b]$ spaced at distance dt using the Schauder basis of level L and level wise coefficients decaying at rate $beta$. A Brownian motion like prior is obtained for $beta = 0.5$. The K remaining optional basiselements have variance $xirem$.

The result is returned as `[designp coeff se]` where `coeff` are coefficients of finite elements with maximum at the designpoints `designp` and standard error `se`.

Observations outside $[a, b]$ may influence the result through $\phi_{n+1}, \dots, \phi_{n+K}$

Module Diffusion

Introduction

The functions in this module operate on three conceptual different objects, (although they are currently just represented as vectors and arrays.)

Stochastic processes, denoted x, y, w are arrays of values which are sampled at distance dt, ds , where dt, ds are either scalar or Vectors `length(dt)=size(W)[end]`. Stochastic differentials are denoted dx, dw etc., and are first differences of stochastic processes. Finally, t can denote the total time or correspond to a vector of `size(W)[end]` sampling time points.

Note the following convention: In analogy with the definition of the Ito integral,

$$\text{intxdw}[i] = x[i](w[i+1]-w[i]) \quad (== x[i]dw[i])$$

and

$$\text{length}(w) = \text{length}(dw) + 1$$

Reference

`Diffusion.brown1(u, t, n::Integer)`

Compute n equally spaced samples of 1d Brownian motion in the interval $[0, t]$, starting from point u

`Diffusion.brown(u, t, d::Integer, n::Integer)`

Simulate n equally spaced samples of d -dimensional Brownian motion in the interval $[0, t]$, starting from point u

`Diffusion.dw1(t, n::Integer)`

`Diffusion.dw(t, d::Integer, n::Integer)`

Simulate a 1-dimensional (d -dimensional) Wiener differential with n values in the the interval $[0, t]$, starting from point u

`Diffusion.dw(dt::Vector, d::Integer)`

Simulate a d -dimensional Wiener differential sampled at time points with distances given by the vector dt

`Diffusion.ito(y, dx)`

`Diffusion.ito(dx)`

`Diffusion.cumsum0(dx)`

Integrate a valued stochastic process with respect to a stochastic differential. R, R^2 (d rows, n columns), R^3 .

`ito(dx)` is a shortcut for `ito(ones(size(dx)[end], dx))`. So `ito(dx)` is just a `cumsum0` function which is a inverse to `dx = diff([0, x1, x2, x3, ...])`.

`..(y, dx)`

`Diffusion.ydx(y, dx)`

`y .. dx` returns the stochastic differential `ydx` defined by the property

$$\text{ito}(ydx) == \text{ito}(y, dx)$$

`Diffusion.bb(u, v, t, n)`

Simulates n equidistant samples of a Brownian bridge from point u to v in time t

`Diffusion.dwcond1(v, t, n)`

Simulates n equidistant samples of a “bridge noise”: that is a Wiener differential dw conditioned on $W(t) = v$

`Diffusion.aug(dw, dt, n)`

`Diffusion.aug(dt, n)`

Take Wiener differential sampled at dt and return Wiener differential subsampled n times between each observation with new length `length(dw)*n`. `aug(dt, n)` computes the corresponding subsample of times.

`Diffusion.quvar(x)`

Computes quadratic variation of x .

`Diffusion.bracket(x)`

`Diffusion.bracket` (x, y)

Computes quadratic variation process of x (of x and y).

`Diffusion.euler` ($t0, u, b, sigma, dt, dw$)

`Diffusion.euler` ($t0, u, b, sigma, dt$)

Simulates a 1-dimensional diffusion process using the Euler-Maruyama approximation with drift $b(t, x)$ and diffusion coefficient $\sigma(t, x)$ starting in $(t0, u)$ using dt and given Wiener differential dw .

Module Randm

Introduction

Random matrices for testing purposes. I did not figure out the actual distributions the matrices are drawn from.

Reference

`Randm.randposdef` (d)

Random positive definite matrix of dimension d .

`Randm.randstable` (d)

Random stable matrix (matrix with eigenvalues with negative real part) with dimension d .

`Randm.randunitary` (d)

Random unitary matrix of dimension d .

`Randm.randorth` (d)

Orthogonal matrix drawn according to the Haar measure on the group of orthogonal matrices.

`Randm.randnormal` (d)

Random normal matrix.

CHAPTER 4

Indices and tables

- `genindex`
- `modindex`
- `search`

Symbols

..() (in module Diffusion), 10

A

aug() (in module Diffusion), 10

B

bayes_drift() (in module Schauder), 9

bb() (in module Diffusion), 10

bcirc() (in module SDE), 5

bracket() (in module Diffusion), 10

brown() (in module Diffusion), 10

brown1() (in module Diffusion), 10

bstar() (in module SDE), 5

C

cumsum0() (in module Diffusion), 10

D

dotVs() (in module SDE), 6

drop

 () (in module Schauder), 7

drop_mu

 () (in module Schauder), 9

drop_Sigma

 () (in module Schauder), 9

dW() (in module Diffusion), 10

dW1() (in module Diffusion), 10

dWcond1() (in module Diffusion), 10

E

euler() (in module Diffusion), 11

euler() (in module SDE), 6

exact() (in module SDE), 6

F

fe_mu() (in module Schauder), 9

fe_muB1() (in module Schauder), 9

fe_muB2() (in module Schauder), 9

fe_Sigma() (in module Schauder), 9

finger_permute() (in module Schauder), 7

finger_pm() (in module Schauder), 8

H

H() (in module SDE), 5

hat() (in module Schauder), 8

I

ito() (in module Diffusion), 10

K

K() (in module SDE), 5

L

level() (in module Schauder), 8

ll() (in module SDE), 6

llikeliXcirc() (in module SDE), 6

lp() (in module SDE), 5, 6

M

mu() (in module SDE), 5

P

pickup

 () (in module Schauder), 7

pickup_mu

 () (in module Schauder), 9

pickup_Sigma

 () (in module Schauder), 9

Q

quvar() (in module Diffusion), 10

R

r() (in module SDE), 5

randnormal() (in module Randm), 11

randorth() (in module Randm), 11

randposdef() (in module Randm), 11

`randstable()` (in module `Randm`), 11
`randunitary()` (in module `Randm`), 11

S

`samplep()` (in module `SDE`), 6
`soft()` (in module `SDE`), 6
`stable()` (in module `SDE`), 6
`syl()` (in module `SDE`), 5

T

`tofs()` (in module `SDE`), 6

V

`vectoroflevels()` (in module `Schauder`), 8
`Vs()` (in module `SDE`), 6

X

`XofU()` (in module `SDE`), 6

Y

`ydx()` (in module `Diffusion`), 10