& Numerical solutions to SDEs

Recall our SDE

dX(t) = a(X(t)) dt + b(X(t)) dw(t), X(0)=Xo,

for some drift and dispersion functions a and b.

In this section, we aim to simulate a skell-ton (trajectory) of the SDE at discrete times. More specifically, we would like to sample X(t,), X(tz)... X(tT) for any times of the trajectory times of the trajectory times.

of times. For notation clarity, we may some times use shorthands

 $X_{k}:=X(t_{k})$, $X_{i}(t_{k})$, the i-th element of vector $X(t_{k})\in\mathbb{R}^{d}$, $X_{i,k}:=X_{i}(t_{k})$.

We will also use $\Delta t_R := t_R - t_{R-1}$ and $\Delta W_R := W(t_R) - W(t_{R-1})$ $\sim N(0, \Delta t_R)$.

The simulation of SDEs is easy if we explicitly know their solutions. However, this is possible only for a few isolated cases, for example, the geometric Brownian motion (see Exercise 1, Assignment 4) and linear SDEs. It is also possible to use the airsanou-based method by Beskos, 2005 to make exact Samples, but the method is restricted to very limited cont classes of SITES. Hence, in this section, we formulate a few most commonly used numerical schemes to approximate the SDEs Solutions.

The Euler-Marayama is argubly the simplest approximate scheme. The idea is:

$$X(t_{k-1}) = X(t_{k-1}) + \begin{cases} t_k & \alpha(X(s)) ds + \int t_k & b(X(s)) & \alpha(X(s)) &$$

Essentially, the Euler-Monayama method approximates

the integrals above by rectangles. This works reasonably
provided that the discretisation time the them is sufficiently

Small. Furthermore, if the dispersion coefficient is

a constant b instead of a function of X, vizz, then the

Itô integral approximation is exact, viz,

the b dw(s) = b DWk.

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This suggests that the EM method is better when dealing with constant dispersion than that of X-dependent ones.

Let's sumarise the algorithm.

Algorithm 9. Euler-Mamyama

Juput: times ti, tz, ... tī output: X(ti), X(tz), _ X(ti) approximate samples

- 1. Draw X60 n P(xo)
- 2. For k=1,2,... T do Draw Dwk ~ N(0, tk-tk-1) X(tk) = X(tk-1) + a(X(tk-1)) (tk-tk-1) + b(X(tk-1)) Dwk.

Refum XItil, Xltz)... Xltz)

Example 10. Modified Duffing-van der Pol Simulation. Consider a two-dimensional SDE:

$$d\left[\begin{array}{c}X_{1}(t)\\X_{2}(t)\end{array}\right]=\left[\begin{array}{c}X_{1}(t)\\X_{1}(t)\end{array}\left(\begin{array}{c}X_{1}(t)\end{array}\right)-X_{2}(t)\end{array}\right]dt+\left[\begin{array}{c}O\\X_{1}(t)\end{array}\right]dw(t),$$

Set d=2, and the initial values $X_1(0)=-3$ and $X_2(0)=0$. Simulate a path of X at times $t_1=0.01$, $t_2=0.02$, ... $t_{1000}=10$. Please See 'Lec3_van_der_Pol. ipynb' for how this is done in Euler--/Mamyama.

Albeit the simplicity of EM, the method can be a quite crucke estimator. More precisely, suppose that twice the drift and the dispersion functions are continuously differentiable and their derivatives are bounded, we can show that the L2-approximation error is such that $\pm [11 \times (t_R) - \frac{1}{2}]^{\frac{1}{2}} \le C(t_R - t_R)^{\beta}$, Lem approximated constant

for $\beta = \frac{1}{2}$ (see, Lord et al., 2014, PP. 338 for details).

also see kloeden and Platen, 1995, ch. 10 for a complete proof.

This means that the approximation emor goes to zero as $\Delta t_R \to 0$ at a speed of $\beta = \frac{1}{2}$. It is then of interests to ask: Can we develop a better approximate scheme in the way that the convergence speed $\beta > \frac{1}{2}$ is faster?

To develop such a high-order simulator, we can leverage Itô's formula. The idea is similar to using Taylor expansion for ODEs.

(intentionally left blank)

Recall the SDE Xttr) = X(tr-1)+ (tr a(X(S)) ds + (tr b(X(S)) c/W(S)), and we let XEIRd, a:IRd > IRd, b:IRdxw > IRd, WEIRM. WE Use notation Xi(th) to denote the i-th element of the vector X(th). Let us apply Itô's formula on ai and bij: (1x(x(s)) = ai(x(tn-1)) + (1xai(x(z))) a(x(z)) + = tr(T(x(z)) Hxai(x(z))) dz $T'(x) := b(x) b(x)^T$ any 5 > tre-1 + (& Qi (X(Z))) h(X(Z)) dW(Z), bij (X(S)) = bij (X(te)) + ((x bij (X(2))) a(X(2)) + = tr(T(X(2)) Hxbij (X(2))) d2 + (((x bij (x (2))) b (x (2)) d w (2).

Now substitute ai(x(s)) and bij(x(s)) back to the SIDE, we get:

If we discard the terms O-O, the we recover the Enter-Manyan scheme. To do better than Enter-Manyama, we can keep some of the terms the appy Ito's formula again on them. As an example, applying Ito's formula on O gives:

(\(\forall \) \(\text{X(Z1)} \) \(\text{Tb(X(Z1)} \) = (\(\text{Xbij} (\text{X(tr-1)}) \) \(\text{D(X(tr-1)}) \) \(\text{Tb} \) \(

which we can substitute back to 19 and discard (3) giving:

recover the celebrated Milstein's method; Xi(tr) x Xi(tr-1) + Qi(X(tr-1))(tr-tr-1) + = (bij (X(tr-1)) DKj, k + (Vx bij (X(tr-1))) b(X(tr-1)) fter (dk/2) The Mistein's method, as we can see from the approximation above, is an improvement of Eulor-Maruyam with an additional (\(\text{bij}(\text{X(the-i)})\) \(\text{b(the-i)} \) \(\text{the fler} \) \(\text{dW(2) dWj(5)}, \) which leverages the gradient information of the dispersion function. When b is a constant, the gradient is zero, the get it reduces to Euler-Manyama. However, the iterated Ito integral the for dw(2) dw(5) is a trouble maker, and is nasty to simulate. Essentially, we have to simulate a bunch of the (s) dw2(2) dw3(s) for all i,j=1,2,...w., which does not scale well in the dimension of the Brownian motion W. FYI: this is highly related to Hermite pbly nomial, see kloden and

platen, 1995.

Fortunately, if the dispersion function b admits special structures, for example, diagonal, or, unidimensional, then the iterated Ito integral is cheap to compute. To see this, suppose that b(x(tk,1)) is alway diagonal, then big(·) = o for all i+j, hence we only need to deal with the start the start dwi(2) dwi(s) for i=1,2, -w. Morover, doing some algebras gives:

So the Milstein's method for this the diagonal dispersion functions simplifies to:

We can rewrite it in a more compact vector format:

O: To prove it, use Itô's formula on \$(1011) = wit).

where

$$\overline{b}(X(t|_{R-1})):=\frac{-c(b_{11}(X(t|_{R-1}))}{c(X_1)}$$

$$\overline{c(x_1)}$$

$$c(x_2)$$

$$c(x_2)$$

$$c(x_3)$$

$$c(x_4)$$

$$c(x_4)$$

Let's sumarise it Milstein's method.

```
Algorithm 11. Milstein's method for diagonal dispersion
Input: Times to, tz, __ tr
output: Approximate &(t), X(tz), ... X(t)
  1. Draw X(0) ~ B. P(X.)
  2, For k=1,2, ... T do
           Draw swk n N(0, tr-tr-1)
            X(tr) = X(tr-1) + a(x(tr-1))(tr-tr-1) + b(x(tr-1)) Alle
                += b(x(tr.1)) b(x(tr.1) ( DWk - (tr-tr.1))
Return X(t,), X(tr), ~ X(tr)
```

Example 12. dx1t)=-2x(t)dt + x/1-x(t)2dW(t), X0=0 Simulate a trajectory at times to=v.vl, tz=v.v2, ... to=5 using Milstein's method. then compare to the true solution XLt) = Sin(XWLt), See, Kloecken and Platen. 1995.
PP. 121. See 'tes' Lec3_ emvs_milstein. ipyub'.

The convergence of Milstein's method is such that #[11X(tr)-X(tr)/2] EC (tr-tr)

for B=1 which is greater that Euler-Manyam's 2.

(under a similar assamption on the SDE coefficients)

It is possible to do better further than Milstein's method, however, as we can see from the derivation of Milstein's method, getting higher ord schemes would be super nasty. The main challenge is really the iterated It's integrals. If the readers are interested, please see kloeden and Platen, 1995 for how to derive high order discretisation schemes, and

they als:	o show how to simply the schemes if the SIDE
coefficien	ts have special structures (e.g., diagonal or commutative,
	B=1.5
latus h	oring up a method called order 1.5 strong Taylor
Let W	I be consider the devivation is seen in
	or constant dispersion. The derivation is seen in
kloeden au	nd Platen, 1995.
Algorithe	M 13. Order 1.5 strong Taylor Scheme.
	Times t, tr, to
super.	approximate Xtt.), Xtt.)
	Draw &(0) ~ P(X.)
2,	For k=1,2, T do
	Draw two independent U., Uz uN(O, In)
	Ada = ofth-the U.
	$\Delta \beta_{R} = \frac{1}{2} (t_{R} - t_{R-1})^{\frac{3}{2}} (U_{1} + \frac{U_{2}}{\sqrt{3}})$ use broadcasting einsum
	For i=1,2, of do
	er=[e,k. ez,k ez,k ed,k]
	X(tr) = X(tr.1) + Q(X(tr.1)) (tr-tr.1) + b Ddr
	+ (Ja(Xthe)) a(Xthe) + 2 ex) (th-th)2 + Ja(Xthe) b BR
K - 1	
Dutpet:	Ztti), Ztti) Jacobian of a.

Scanssian-increment based simulations

We find that the aforementioned Euler-Mamya ma and order 1.5 strong Taylor methods are Gaussian based. More precisely, we can generalise than and the Same framework:

X(the) & #[X(the) | X(the)] + Cov[X(the) | X(the)) &k,

&k ~ N(0, Id),

Cholesky clecomposition

Euler-Maruyama gives

Euler-Maruyama gives

E[Xltn) | Xltn) | Xltn) + a(Xltn) | str,

Car[Xltn) | Xltn) | Car[Xltn) | Car[Xltn) | str.

Order 1.5 strong Taylor gives:

E[XIthe) 1 x (the) 2 X(the) + a(XIthe) str

+(Ja(XIthe)) a(XIthe)) + 2 er) str

Cov[X(the) | x(the)) & Dota + Ja(X(the)) Ta(X(the)) \frac{1}{2} \frac{1}{2}

+ (TJa(X(the-1))T+ Ja(X(the-1))T) sth?

Algorithm It. Gaussian-increment based sampling.

Input: Times t, tz, _ t7

Output: Approximate Xlti), Xlti), Xlti).

1. Draw X(0) ~ P(X0)

2. For k=1, 2, ... T do

Draw Lk ~ N(0, Id)

X(tk) = \(\frac{1}{2}\) Xlti), Xlti) \(\frac{1}{2}\) Xltin) \(\frac{1}{2}\) Xltin) \(\frac{1}{2}\) Xltin) \(\frac{1}{2}\) Xltin)

Reform \(\frac{1}{2}\) Liti), Xlti), \(\frac{1}{2}\) \(\frac{1}{2}\)

'Lec3_em_vs_ito15_vs_time. ipynb' implements the Gaussian-based Simulation.

The Milstein's method, however, cannot be put into this generalisation because it has DWk which is not Gaussian.

The upside of such a generalisation is that we can to some extent avoid the complicated iterations of Itô's formula to derive the high order schemes. More specifically, in the next section, we will show that it is can be easier and automatic to approximate the statistical quantities, such as E[XCGN [XCGN]] and ON[XCGN] [XCGN]. Compared to approximate the painful iterated Itô integrals.

Another upsicle is that we can easily compare the methods by letting by letting under the same framework 50 we can tet the methods Share the Same randomness (i.e., h., hr., _ 1/47).

Example # 15

dx(+)=-0x(+)d+ dw(+), 0>0.

This is a linear SDE called Ornstein-Uhlenbech process, we can compute its conditional mean and Variance in closed-form:

E[Xltr) 1 Xltr) = e O(tr-tr)
Xltry)

We can compare Euler-Maruyama and the order 1.5 method, See 'lec3-em-vs-itols-vs-tme.ipyub. In the Script, there is another method Called TME which we shall detail in the next Section.

The downside of the Gaussian-based simulation is that it can be a poor estimator when the transition is quite non-Gaussian. Even if we can approximate EIX(GR) IX(GR-1) and Cov[X(GR-1)] exactly, the estimator can still be poor; for some models it can even be worse than Milstein's method, especially when the dispersion term is non-constant.