Monte Carlo Methods for Uncertainty Quantification

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Lecture outline

Lecture 3: financial SDE applications

- financial models
- approximating SDEs
- weak and strong convergence
- mean square error decomposition
- multilevel Monte Carlo

In computational finance, stochastic differential equations are used to model the behaviour of

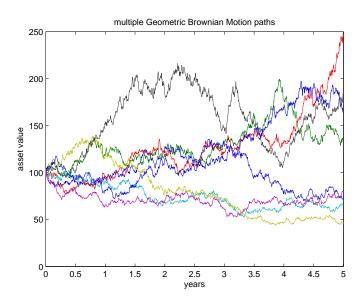
- stocks
- interest rates
- exchange rates
- weather
- electricity/gas demand
- crude oil prices
- ...

Stochastic differential equations are just ordinary differential equations plus an additional random source term.

The stochastic term accounts for the uncertainty of unpredictable day-to-day events.

The aim is **not** to predict exactly what will happen in the future, but to predict the probability of a range of possible things that **might** happen, and compute some averages, or the probability of an excessive loss.

This is really just uncertainty quantification, and they've been doing it for quite a while because they have so much uncertainty.



Examples:

• Geometric Brownian motion (Black-Scholes model for stock prices)

$$dS = r S dt + \sigma S dW$$

Cox-Ingersoll-Ross model (interest rates)

$$dr = \alpha(b - r) dt + \sigma \sqrt{r} dW$$

Heston stochastic volatility model (stock prices)

$$dS = r S dt + \sqrt{V} S dW_1$$

$$dV = \lambda (\sigma^2 - V) dt + \xi \sqrt{V} dW_2$$

with correlation ρ between dW_1 and dW_2

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Generic Problem

Stochastic differential equation with general drift and volatility terms:

$$dS(t) = a(S, t) dt + b(S, t) dW(t)$$

W(t) is a Wiener variable with the properties that for any q < r < s < t, W(t) - W(s) is Normally distributed with mean 0 and variance t - s, independent of W(r) - W(q).

In many finance applications, we want to compute the expected value of an option dependent on the terminal state P(S(T))

Other options depend on the average, minimum or maximum over the whole time interval.

Euler discretisation

Given the generic SDE:

$$dS(t) = a(S) dt + b(S) dW(t), \quad 0 < t < T,$$

the Euler discretisation with timestep h is:

$$\widehat{S}_{n+1} = \widehat{S}_n + a(\widehat{S}_n) h + b(\widehat{S}_n) \Delta W_n$$

where ΔW_n are Normal with mean 0, variance h.

- How good is this approximation?
- How do the errors behave as $h \to 0$?

These are much harder questions when working with SDEs instead of ODEs.

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Weak convergence

For most finance applications, what matters is the **weak** order of convergence, defined by the error in the expected value of the payoff.

For a European option, the weak order is m if

$$\mathbb{E}\left[f(S(T))\right] - \mathbb{E}\left[f(\widehat{S}_N)\right] = O(h^m)$$

The Euler scheme has order 1 weak convergence, so the discretisation "bias" is asymptotically proportional to h.

Strong convergence

In some Monte Carlo applications, what matters is the **strong** order of convergence, defined by the average error in approximating each individual path.

For the generic SDE, the strong order is m if

$$\mathbb{E}\left[\left|S(T)-\widehat{S}_{N}\right|\right]=O(h^{m})$$

The Euler scheme has order 1/2 strong convergence.

The leading order errors are as likely to be positive as negative, and so cancel out – this is why the weak order is higher.

Exotic options

• Lookback option: $P = \left(S(T) - \min_{0 < t < T} S(t)\right)$

Approximation $\widehat{S}_{min} = \min_n \widehat{S}_n$ gives $O(h^{1/2})$ weak convergence

• Barrier option (down-and-out call): $P = \mathbf{1}(\min_{0 < t < T} S(t) > B) \max(0, S(T) - K)$

Approximation using \widehat{S}_{min} gives $O(h^{1/2})$ weak convergence

• Asian option: $P = \max \left(0, T^{-1} \int_0^T S(t) dt - K\right)$

Trapezodial integration gives O(h) weak convergence



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Exotic options

The poor weak convergence for the lookback and barrier options is due to the fact that there is an $O(h^{1/2})$ change in O(S(t)) within each timestep.

It is possible to approximate this (using something called a Brownian Bridge construction) and recover first order weak convergence.

Key point: getting high order convergence is very difficult.

Mean Square Error

Finally, how to decide whether it is better to increase the number of timesteps (reducing the weak error) or the number of paths (reducing the Monte Carlo sampling error)?

If the true option value is

$$V = \mathbb{E}[f]$$

and the discrete approximation is

$$\widehat{V} = \mathbb{E}[\widehat{f}]$$

and the Monte Carlo estimate is

$$\widehat{Y} = \frac{1}{N} \sum_{n=1}^{N} \widehat{f}^{(n)}$$

then ...

Mean Square Error

...the Mean Square Error is

$$\mathbb{E}\left[\left(\widehat{Y} - V\right)^{2}\right] = \mathbb{E}\left[\left(\widehat{Y} - \mathbb{E}[\widehat{f}] + \mathbb{E}[\widehat{f}] - \mathbb{E}[f]\right)^{2}\right]$$

$$= \mathbb{E}\left[\left(\widehat{Y} - \mathbb{E}[\widehat{f}]\right)^{2}\right] + (\mathbb{E}[\widehat{f}] - \mathbb{E}[f])^{2}$$

$$= N^{-1}\mathbb{V}[\widehat{f}] + \left(\mathbb{E}[\widehat{f}] - \mathbb{E}[f]\right)^{2}$$

- first term is due to the variance of estimator
- second term is square of bias due to weak error

Mean Square Error

Given first order weak convergence and M timesteps, the computational cost is proportional to C = NM and the MSE is approximately

$$a N^{-1} + b M^{-2} = a N^{-1} + b C^{-2} N^2.$$

For a fixed computational cost, this is a minimum when

$$N = \left(\frac{aC^2}{2b}\right)^{1/3}, \quad M = \left(\frac{2bC}{a}\right)^{1/3},$$

and the two error terms have a similar magnitude.

Hence the cost to achieve a RMS error of ε requires $M=O(\varepsilon^{-1})$ and $N=O(\varepsilon^{-2})$, so the total cost is $O(\varepsilon^{-3})$.

Multilevel Monte Carlo

When solving finite difference equations coming from approximating PDEs, multigrid combines calculations on a nested sequence of grids to get the accuracy of the finest grid at a much lower computational cost.

Multilevel Monte Carlo uses a similar idea to achieve variance reduction in Monte Carlo path calculations, combining simulations with different numbers of timesteps – same accuracy as finest calculations, but at a much lower computational cost.

Can also be viewed as a recursive control variate strategy.

Consider multiple sets of simulations with different timesteps $h_\ell=2^{-\ell}\ T,\ \ell=0,1,\ldots,L$, and payoff \widehat{P}_ℓ

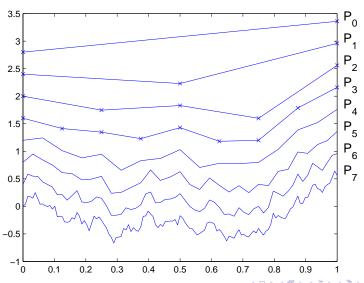
$$\mathbb{E}[\widehat{P}_L] = \mathbb{E}[\widehat{P}_0] + \sum_{\ell=1}^L \mathbb{E}[\widehat{P}_\ell - \widehat{P}_{\ell-1}]$$

Expected value is same – aim is to reduce variance of estimator for a fixed computational cost.

Key point: approximate $\mathbb{E}[\widehat{P}_\ell - \widehat{P}_{\ell-1}]$ using N_ℓ simulations with \widehat{P}_ℓ and $\widehat{P}_{\ell-1}$ obtained using <u>same</u> Brownian path.

$$\widehat{Y}_{\ell} = N_{\ell}^{-1} \sum_{i=1}^{N_{\ell}} \left(\widehat{P}_{\ell}^{(i)} - \widehat{P}_{\ell-1}^{(i)} \right)$$

Discrete Brownian path at different levels



Using independent paths for each level, the variance of the combined estimator is

$$\mathbb{V}\left[\sum_{\ell=0}^L \widehat{Y}_\ell\right] = \sum_{\ell=0}^L N_\ell^{-1} V_\ell, \qquad V_\ell \equiv \mathbb{V}[\widehat{P}_\ell - \widehat{P}_{\ell-1}],$$

and the computational cost is proportional to $\sum_{\ell=0}^L N_\ell \; h_\ell^{-1}$.

Hence, the variance is minimised for a fixed computational cost by choosing N_{ℓ} to be proportional to $\sqrt{V_{\ell} h_{\ell}}$.

The constant of proportionality can be chosen so that the combined variance is $O(\varepsilon^2)$.

For the Euler discretisation and the Lipschitz payoff function

$$\mathbb{V}[\widehat{P}_{\ell} - P] = O(h_{\ell}) \implies \mathbb{V}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] = O(h_{\ell})$$

and the optimal N_ℓ is asymptotically proportional to h_ℓ .

To make the combined variance $O(\varepsilon^2)$ requires

$$N_{\ell} = O(\varepsilon^{-2} L h_{\ell}).$$

To make the bias $O(\varepsilon)$ requires

$$L = \log_2 \varepsilon^{-1} + O(1) \implies h_L = O(\varepsilon).$$

Hence, we obtain an $O(\varepsilon^2)$ MSE for a computational cost which is $O(\varepsilon^{-2}L^2) = O(\varepsilon^{-2}(\log \varepsilon)^2)$.

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Theorem: Let P be a functional of the solution of a stochastic o.d.e., and \widehat{P}_l the discrete approximation using a timestep $h_\ell = M^{-\ell} T$.

If there exist independent estimators \widehat{Y}_{ℓ} based on N_{ℓ} Monte Carlo samples, and positive constants $\alpha \geq \frac{1}{2}, \beta, c_1, c_2, c_3$ such that

i)
$$\left|\mathbb{E}[\widehat{P}_{\ell}-P]\right| \leq c_1 h_{\ell}^{\alpha}$$

ii)
$$\mathbb{E}[\widehat{Y}_{\ell}] = \left\{ egin{array}{ll} \mathbb{E}[\widehat{P}_{0}], & \ell = 0 \\ \mathbb{E}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}], & \ell > 0 \end{array}
ight.$$

- iii) $\mathbb{V}[\widehat{Y}_{\ell}] \leq c_2 N_{\ell}^{-1} h_{\ell}^{\beta}$
- iv) C_{ℓ} , the computational complexity of \widehat{Y}_{ℓ} , is bounded by

$$C_\ell \leq c_3 \ N_\ell \ h_\ell^{-1}$$



then there exists a positive constant c_4 such that for any $\varepsilon < e^{-1}$ there are values L and N_ℓ for which the multi-level estimator

$$\widehat{Y} = \sum_{\ell=0}^{L} \widehat{Y}_{\ell},$$

has Mean Square Error MSE $\equiv \mathbb{E}\left[\left(\widehat{Y} - \mathbb{E}[P]\right)^2\right] < \varepsilon^2$

with a computational complexity C with bound

$$C \leq \left\{ \begin{array}{ll} c_4 \, \varepsilon^{-2}, & \beta > 1, \\ \\ c_4 \, \varepsilon^{-2} (\log \varepsilon)^2, & \beta = 1, \\ \\ c_4 \, \varepsilon^{-2 - (1 - \beta)/\alpha}, & 0 < \beta < 1. \end{array} \right.$$

Convergence Test

Asymptotically,

$$\mathbb{E}[\widehat{P}_L - \widehat{P}_{L-1}] \approx (M-1)\,\mathbb{E}[P - \widehat{P}_L]$$

so this can be used to decide when the bias error is sufficiently small.

In case the correction changes sign at some level, it is safer to use the convergence test

$$\max\left\{M^{-1}\left|\widehat{Y}_{L-1}\right|,\left|\widehat{Y}_{L}\right|\right\}<(M-1)\frac{\varepsilon}{\sqrt{2}}.$$

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Multilevel Algorithm

- \bigcirc start with L=0
- 2 estimate V_L using an initial $N_L = 10^4$ samples
- **3** define optimal N_{ℓ} , $\ell = 0, \ldots, L$
- $oldsymbol{\circ}$ evaluate extra samples as needed for new \mathcal{N}_ℓ
- \odot if $L \ge 2$, test for convergence
- **1** if L < 2 or not converged, set L := L + 1 and go to 2.

Numerical results use M=4, which is almost twice as efficient as M=2.

Geometric Brownian motion:

$$dS = r S dt + \sigma S dW, \qquad 0 < t < 1,$$

$$S(0)=1$$
, $r=0.05$, $\sigma=0.2$

Heston model:

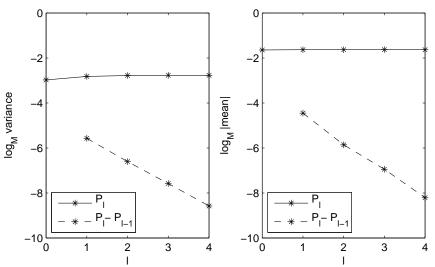
$$\begin{split} \mathrm{d} S &= r \, S \, \mathrm{d} t + \sqrt{V} \, S \, \mathrm{d} W_1, \qquad 0 < t < 1 \\ \mathrm{d} V &= \lambda \left(\sigma^2 - V \right) \mathrm{d} t + \xi \, \sqrt{V} \, \mathrm{d} W_2, \end{split}$$

$$S(0)=1$$
, $V(0)=0.04$, $r=0.05$, $\sigma=0.2$, $\lambda=5$, $\xi=0.25$, $\rho=-0.5$

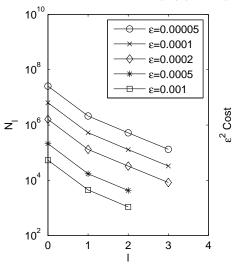
All calculations use M=4, more efficient than M=2.

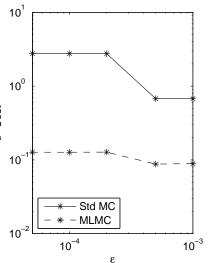


GBM: European call, max(S(1)-1,0)

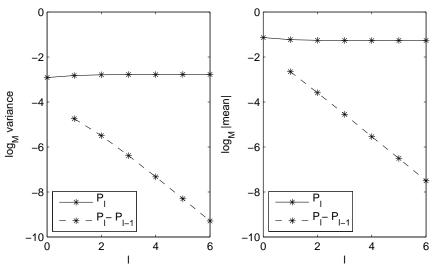


GBM: European call, max(S(1)-1,0)

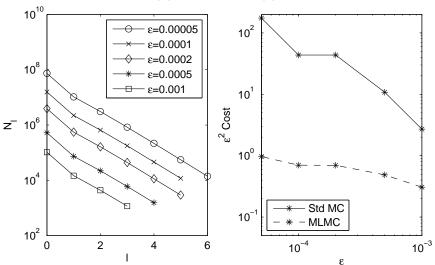




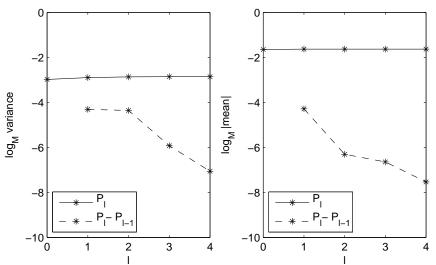
GBM: lookback option, $S(1) - \min_{0 < t < 1} S(t)$



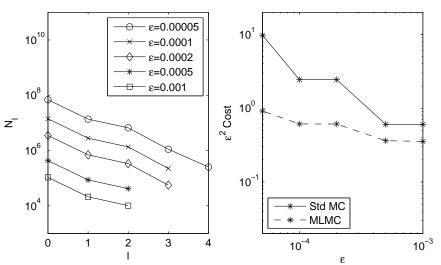
GBM: lookback option, $S(1) - \min_{0 < t < 1} S(t)$



Heston model: European call



Heston model: European call



Extensions

- Milstein discretisation gives better strong convergence, and hence better multilevel performance
- Quasi-Monte Carlo very effective on coarse grids and reduces overall cost to roughly $O(\varepsilon^{-1.5})$ in simplest cases
- multivariate discontinuous payoffs simplest approach is to use "splitting" for multiple simulations of final timestep
- jump-diffusion and Lévy processes (more realistic models than Brownian diffusion)

References

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