

Multilevel Monte Carlo methods for financial applications

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Introduction

Outline:

- ▶ basic ideas
- ▶ three finance applications:
 - ▶ SDEs for option pricing
 - ▶ reduced precision computing and approximate random variables
 - ▶ nested simulations, such as CVaR

In doing this I hope to emphasise:

- ▶ the simplicity of the idea
- ▶ its flexibility – it's not prescriptive, more a generic approach
- ▶ sources of further information

Objective

To achieve a root-mean-square accuracy of ε , Monte Carlo simulation requires $O(\varepsilon^{-2})$ samples.

In many cases the cost of each sample also depends on ε , so the overall cost is often $O(\varepsilon^{-3})$ or worse.

The aim is to reduce the total cost to $O(\varepsilon^{-2})$, and reduce the cost even it is already $O(\varepsilon^{-2})$.

Control variates

Control variates are a well-established technique for reducing variance, and hence computational cost

If we want to estimate $\mathbb{E}[P(\omega)]$, and we know $\mathbb{E}[Q(\omega)]$ for some $Q(\omega)$ which is well correlated to $P(\omega)$, then

$$\mathbb{E}[P] = \underbrace{\lambda \mathbb{E}[Q]}_{\text{known}} + \mathbb{E}[P - \lambda Q]$$

so we can instead use Monte Carlo to estimate $\mathbb{E}[P - \lambda Q]$ and choose λ to minimise the variance, giving

$$\mathbb{V}[P - \lambda Q] = (1 - \rho^2) \mathbb{V}[P]$$

where ρ is the correlation coefficient.

Two-level Monte Carlo

If we want to estimate $\mathbb{E}[P]$ but it is much cheaper to simulate $\tilde{P} \approx P$, then since

$$\mathbb{E}[P] = \mathbb{E}[\tilde{P}] + \mathbb{E}[P - \tilde{P}]$$

we can use the estimator

$$N_0^{-1} \sum_{n=1}^{N_0} \tilde{P}^{(0,n)} + N_1^{-1} \sum_{n=1}^{N_1} \left(P^{(1,n)} - \tilde{P}^{(1,n)} \right)$$

Similar to a control variate except that

- ▶ we don't know analytic value of $\mathbb{E}[\tilde{P}]$, so need to estimate it
- ▶ there is no multiplicative factor λ

Benefit: if $P - \tilde{P}$ is small, its variance will be small, so won't need many samples to accurately estimate $\mathbb{E}[P - \tilde{P}]$, so cost will be reduced greatly.

Two-level Monte Carlo

If we define

- ▶ C_0, V_0 cost and variance of one sample of \tilde{P}
- ▶ C_1, V_1 cost and variance of one sample of $P - \tilde{P}$

then the total cost and variance of this estimator is

$$C_{tot} = N_0 C_0 + N_1 C_1 \quad \implies \quad V_{tot} = V_0/N_0 + V_1/N_1$$

Treating N_0, N_1 as real variables, using a Lagrange multiplier to minimise the cost subject to a fixed variance gives

$$\frac{\partial}{\partial N_\ell} (C_{tot} + \mu^2 V_{tot}) = 0, \quad N_\ell = \mu \sqrt{V_\ell / C_\ell}$$

Choosing μ s.t. $V_{tot} = \varepsilon^2$ gives

$$C_{tot} = \varepsilon^{-2} (\sqrt{V_0 C_0} + \sqrt{V_1 C_1})^2$$

Multilevel Monte Carlo

Natural generalisation: given a sequence $\hat{P}_0, \hat{P}_1, \dots, \hat{P}_L$

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

we can use the estimator

$$\hat{Y} = N_0^{-1} \sum_{n=1}^{N_0} \hat{P}_0^{(0,n)} + \sum_{\ell=1}^L \left\{ N_\ell^{-1} \sum_{n=1}^{N_\ell} \left(\hat{P}_\ell^{(\ell,n)} - \hat{P}_{\ell-1}^{(\ell,n)} \right) \right\}$$

with independent estimation for each level of correction

Multilevel Monte Carlo

If we define

- ▶ C_0, V_0 to be cost and variance of \widehat{P}_0
- ▶ C_ℓ, V_ℓ to be cost and variance of $\widehat{P}_\ell - \widehat{P}_{\ell-1}$

then the total cost is $\sum_{\ell=0}^L N_\ell C_\ell$ and the variance is $\sum_{\ell=0}^L N_\ell^{-1} V_\ell$.

Minimise the cost for a fixed variance

$$\frac{\partial}{\partial N_\ell} \sum_{k=0}^L (N_k C_k + \mu^2 N_k^{-1} V_k) = 0$$

gives

$$N_\ell = \mu \sqrt{V_\ell / C_\ell} \quad \implies \quad N_\ell C_\ell = \mu \sqrt{V_\ell C_\ell}$$

Multilevel Monte Carlo

Setting the total variance equal to ε^2 gives

$$\mu = \varepsilon^{-2} \left(\sum_{\ell=0}^L \sqrt{V_\ell C_\ell} \right)$$

and hence, the total cost is

$$\sum_{\ell=0}^L N_\ell C_\ell = \varepsilon^{-2} \left(\sum_{\ell=0}^L \sqrt{V_\ell C_\ell} \right)^2$$

in contrast to the standard cost which is approximately $\varepsilon^{-2} V_0 C_L$.

The MLMC cost savings are therefore approximately:

- ▶ V_L/V_0 , if $\sqrt{V_\ell C_\ell}$ increases with level
- ▶ C_0/C_L , if $\sqrt{V_\ell C_\ell}$ decreases with level

Multilevel Monte Carlo

If $\widehat{P}_0, \widehat{P}_1, \dots \rightarrow P$, then the Mean Square Error has the decomposition

$$\begin{aligned}\mathbb{E} \left[(\widehat{Y} - \mathbb{E}[P])^2 \right] &= \mathbb{V}[\widehat{Y}] + \left(\mathbb{E}[\widehat{Y}] - \mathbb{E}[P] \right)^2 \\ &= \sum_{\ell=0}^L V_{\ell}/N_{\ell} + \left(\mathbb{E}[\widehat{P}_L] - \mathbb{E}[P] \right)^2\end{aligned}$$

so can choose L so that $\left| \mathbb{E}[\widehat{P}_L] - \mathbb{E}[P] \right| < \varepsilon/\sqrt{2}$

and then choose N_{ℓ} so that $\sum_{\ell=0}^L V_{\ell}/N_{\ell} < \varepsilon^2/2$

MLMC Theorem

(Slight generalisation of version in my original 2008 *Operations Research* paper, "Multilevel Monte Carlo path simulation")

If there exist independent estimators \hat{Y}_ℓ based on N_ℓ Monte Carlo samples, each costing C_ℓ , and positive constants $\alpha, \beta, \gamma, c_1, c_2, c_3$ such that $\alpha \geq \frac{1}{2} \min(\beta, \gamma)$ and

$$\text{i) } \left| \mathbb{E}[\hat{P}_\ell - P] \right| \leq c_1 2^{-\alpha \ell}$$

$$\text{ii) } \mathbb{E}[\hat{Y}_\ell] = \begin{cases} \mathbb{E}[\hat{P}_0], & \ell = 0 \\ \mathbb{E}[\hat{P}_\ell - \hat{P}_{\ell-1}], & \ell > 0 \end{cases}$$

$$\text{iii) } \mathbb{V}[\hat{Y}_\ell] \leq c_2 N_\ell^{-1} 2^{-\beta \ell}$$

$$\text{iv) } \mathbb{E}[C_\ell] \leq c_3 2^{\gamma \ell}$$

MLMC Theorem

then there exists a positive constant c_4 such that for any $\varepsilon < 1$ there exist L and N_ℓ for which the multilevel estimator

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

has a mean-square-error with bound $\mathbb{E} \left[\left(\hat{Y} - \mathbb{E}[P] \right)^2 \right] < \varepsilon^2$

with an expected computational cost C with bound

$$C \leq \begin{cases} c_4 \varepsilon^{-2}, & \beta > \gamma, \\ c_4 \varepsilon^{-2} (\log \varepsilon)^2, & \beta = \gamma, \\ c_4 \varepsilon^{-2 - (\gamma - \beta)/\alpha}, & 0 < \beta < \gamma. \end{cases}$$

MLMC Theorem

Two observations of optimality:

- ▶ MC simulation needs $O(\varepsilon^{-2})$ samples to achieve RMS accuracy ε , so when $\beta > \gamma$, the cost is optimal — $O(1)$ cost per sample on average.
(Would need multilevel QMC to further reduce costs)
- ▶ When $\beta < \gamma$, another interesting case is when $\beta = 2\alpha$, which corresponds to $\mathbb{E}[\widehat{Y}_\ell]$ and $\sqrt{\mathbb{E}[\widehat{Y}_\ell^2]}$ being of the same order as $\ell \rightarrow \infty$.
In this case, the total cost is $O(\varepsilon^{-\gamma/\alpha})$, which is the cost of a single sample on the finest level — again optimal.

MLMC

Numerical algorithm:

1. start with $L=0$
2. if $L < 2$, get an initial estimate for V_L using $N_L = 1000$ samples, otherwise extrapolate from earlier levels
3. for $\ell \leq L$, determine optimal N_ℓ to achieve $\sum_{\ell=0}^L V_\ell / N_\ell \leq \varepsilon^2 / 2$
4. perform extra calculations as needed, updating estimates of V_ℓ
5. if $L < 2$ or the bias estimate is greater than $\varepsilon / \sqrt{2}$, set $L := L+1$ and go back to step 2

Application: SDEs

With SDEs, level ℓ corresponds to approximation using M^ℓ timesteps, giving approximate payoff \widehat{P}_ℓ at cost $C_\ell = O(M^\ell)$.

Usually choose M in the range 2 – 4; often 4 for Euler-Maruyama and 2 for Milstein discretisation.

Simplest estimator for $\mathbb{E}[\widehat{P}_\ell - \widehat{P}_{\ell-1}]$ for $\ell > 0$ is

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

using same driving Brownian path for both levels.

The implementation is easy – for a coarse timestep of size Mh we simply sum the M fine path Brownian increments ΔW to get the coarse path Brownian increment.

Application: SDEs

Euler-Maruyama discretisation has $O(h^{1/2})$ strong convergence so

$$\mathbb{E}[(\widehat{S}_{\ell,T} - S_T)^2] = O(h_\ell) \implies \mathbb{E}[(\widehat{S}_{\ell,T} - \widehat{S}_{\ell-1,T})^2] = O(h_\ell)$$

Hence for Lipschitz European payoff functions $P \equiv f(S_T)$,

$$\mathbb{V}[\widehat{P}_\ell - \widehat{P}_{\ell-1}] \leq \mathbb{E}[(\widehat{P}_\ell - \widehat{P}_{\ell-1})^2] \leq K^2 \mathbb{E}[(\widehat{S}_{T,\ell} - \widehat{S}_{T,\ell-1})^2] = O(h_\ell)$$

In terms of the MLMC theorem, this means we have

$$\begin{aligned} C_\ell = O(M^\ell) &\implies \gamma = \log_2 M, \\ V_\ell = O(h_\ell) = O(M^{-\ell}) &\implies \beta = \log_2 M, \end{aligned}$$

so the overall cost to achieve ε RMS accuracy is $O(\varepsilon^{-2} |\log \varepsilon|^2)$.

Things are not so good for digital options – complexity is $O(\varepsilon^{-5/2})$.

MLMC SDE algorithm

Input: fine and coarse timesteps h^f, h^c , final time $T = N h^c$, refinement factor $M = h^c/h^f$, initial states $\widehat{S}^f = \widehat{S}^c = S_0$

for $n = 1, N$ **do**

$\Delta W^c := 0$

for $m = 1, M$ **do**

generate r.v. $\Delta W^f \sim N(0, h^f)$

$\Delta W^c := \Delta W^c + \Delta W^f$

$\widehat{S}^f := \widehat{S}^f + a(\widehat{S}^f) h^f + b(\widehat{S}^f) \Delta W^f$

end for

$\widehat{S}^c := \widehat{S}^c + a(\widehat{S}^c) h^c + b(\widehat{S}^c) \Delta W^c$

end for

$\widehat{P}_\ell - \widehat{P}_{\ell-1} := f(\widehat{S}^f) - f(\widehat{S}^c)$

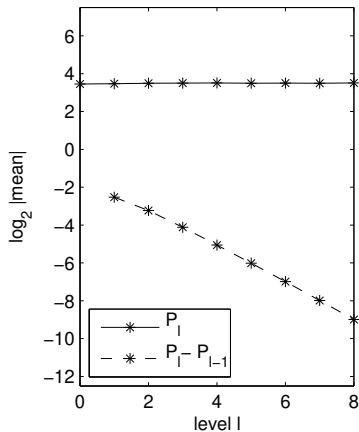
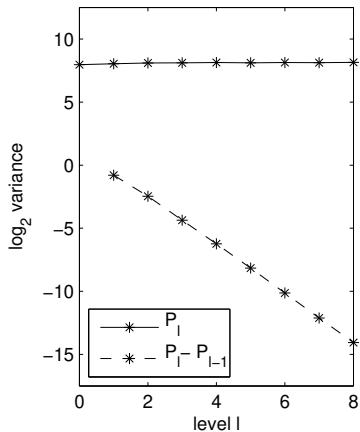
Application: SDEs

- ▶ Milstein discretisation – MBG (2008, 2012)
- ▶ MLQMC for SDEs – MBG, Waterhouse (2009)
- ▶ Greeks – Burgos (2011)
- ▶ American options – Belomestny & Schoenmakers (2011)
- ▶ jump-diffusion models – MBG, Xia (2012)
- ▶ Lévy-driven processes – Dereich (2010), Marxen (2010), Dereich & Heidenreich (2011), Kyprianou (2014)
- ▶ multi-dim Milstein without Lévy areas – MBG, Szpruch (2014)
- ▶ adaptive timesteps – Hoel, von Schwerin, Szepessy, Tempone (2012), MBG, Lester, Whittle (2014), Fang, MBG (2020)
- ▶ exponential Lévy processes – Xia (2017)

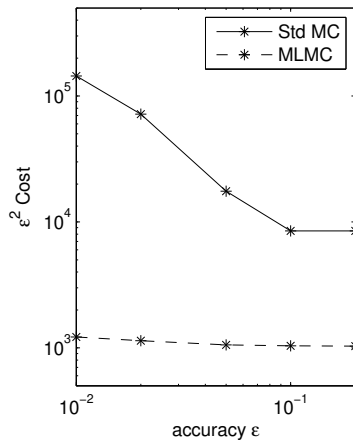
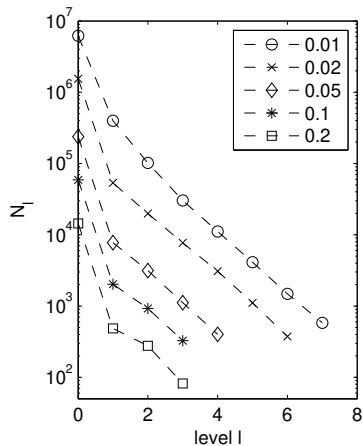
Basket call option

- ▶ 5 underlying assets, modelled by Geometric Brownian Motion with correlation between 5 driving Brownian motions
- ▶ Milstein numerical approximation
- ▶ standard call option based on average at final time T

Basket call option



Basket call option



Application: reduced precision

One simple use of two-level MLMC is with reduced precision floating point arithmetic:

- ▶ double precision on “fine” level
- ▶ single precision (or even half-precision?) on “coarse” level

This can be combined with SDE treatment by using this two-level treatment for each of the expectations $\mathbb{E}[\hat{P}_\ell - \hat{P}_{\ell-1}]$,

$$\mathbb{E}[\hat{P}_\ell - \hat{P}_{\ell-1}] = \mathbb{E}[\hat{P}_\ell^{\text{float}} - \hat{P}_{\ell-1}^{\text{float}}] + \mathbb{E}[(\hat{P}_\ell - \hat{P}_{\ell-1}) - (\hat{P}_\ell^{\text{float}} - \hat{P}_{\ell-1}^{\text{float}})]$$

For the low-accuracy computations can also use imprecise conversion of uniform r.v.'s to Normal r.v.'s

Application: approximate r.v.'s

Continuing with that idea, one approach for the CIR model

$$dr_t = a(b-r) dt + \sigma\sqrt{r} dW_t$$

is to use exact simulation which involves sampling from the non-central χ^2 distribution.

One method of sampling is to generate a $(0, 1)$ uniform r.v. U and then apply the inverse of the non-central χ^2 CDF.

Doing this accurately is expensive, but one can construct a cheap fairly accurate piecewise bilinear approximation – linear in both U and the non-centrality parameter (for fixed degrees of freedom $d = 4ab/\sigma^2$).

Looks a good approach for interest rate models and Heston stochastic volatility.

Application: nested expectation

The general form of a nested expectation is $\mathbb{E} \left[f(\mathbb{E}[Y | X]) \right]$.

The standard approach uses N outer samples $X^{(n)}$, and for each M inner samples $Y^{(m,n)}$. To achieve ε RMS accuracy usually requires $N = O(\varepsilon^{-2})$, $M = O(\varepsilon^{-1})$, so the total cost is $O(\varepsilon^{-3})$.

The MLMC approach uses $M_\ell = 2^\ell M_0$ inner samples on level ℓ

- ▶ cost is $O(2^\ell)$ so $\gamma = 1$
- ▶ an “antithetic” construction gives $\beta = 2$ if f is smooth, and $\beta = 3/2$ if f is continuous and piecewise smooth; both good enough for $O(\varepsilon^{-2})$ total cost
- ▶ if f is discontinuous $\beta = 1/2$ and the cost is $O(\varepsilon^{-5/2})$; can be improved to $O(\varepsilon^{-2} |\log \varepsilon|^2)$ using adaptive sampling

Application: VaR and CVaR

VaR loss L_η is defined implicitly by $\mathbb{P}[L > L_\eta] = \eta$.

Given an estimate \tilde{L}_η , Rockafellar & Uryasev (2000) show that CVaR is

$$\begin{aligned}\mathbb{E}[L \mid L > L_\eta] &= L_\eta + \eta^{-1} \mathbb{E}[\max(0, L - L_\eta)] \\ &= \min_x \{x + \eta^{-1} \mathbb{E}[\max(0, L - x)]\} \\ &= \tilde{L}_\eta + \eta^{-1} \mathbb{E}[\max(0, L - \tilde{L}_\eta)] + O(\tilde{L}_\eta - L_\eta)^2\end{aligned}$$

For ε RMS error,

- ▶ first estimate \tilde{L}_η to accuracy $O(\varepsilon^{1/2})$ at cost $o(\varepsilon^{-2})$
- ▶ then estimate $\eta^{-1} \mathbb{E}[\max(0, L - \tilde{L}_\eta)]$ to accuracy ε using MLMC; $\beta = 3/2$ so total cost is $O(\varepsilon^{-2})$

Can also use random sampling to reduce cost for portfolios with lots of products.

Final comments

- ▶ MLMC has become widely used in academia over the past 10 years, and also MLQMC in some areas (mainly PDEs)
- ▶ very large savings in some application areas (especially PDEs and stochastic modelling of chemical reactions)
- ▶ very limited uptake in the finance sector so far, but I think there are very good opportunities here
- ▶ research worldwide (inc. papers) is listed on a webpage:
`people.maths.ox.ac.uk/gilesm/mlmc_community.html`
- ▶ MLMC software and examples available on another webpage:
`people.maths.ox.ac.uk/gilesm/mlmc/`
- ▶ my papers are on:
`people.maths.ox.ac.uk/gilesm/mlmc.html`
best to start with review: 'Multilevel Monte Carlo methods'.
Acta Numerica, 24:259-328, 2015.