MLMC for reflected diffusions

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Outline

- multilevel Monte Carlo
- 1D particles with mass
 - standard treatment
 - expanded domain
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 - results
- 1D massless particles
 - new treatment
 - results
- multi-dimensional generalisations

Multilevel Monte Carlo

MLMC is based on the telescoping sum

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

where \widehat{P}_{ℓ} represents an approximation using on level ℓ .

In SDE applications with uniform timestep $h_\ell=2^{-\ell}\,h_0$, if the weak convergence is

$$\mathbb{E}[\widehat{P}_{\ell}-P]=O(2^{-\alpha\,\ell}),$$

 \widehat{Y}_ℓ is an unbiased estimator for $\mathbb{E}[\widehat{P}_\ell - \widehat{P}_{\ell-1}]$, based on N_ℓ samples, with variance

$$\mathbb{V}[\widehat{Y}_{\ell}] = O(N_{\ell}^{-1} 2^{-\beta \ell}),$$

and expected cost

$$\mathbb{E}[C_{\ell}] = O(N_{\ell} 2^{\gamma \ell}), \dots$$

Multilevel Monte Carlo

... then the finest level L and the number of samples N_{ℓ} on each level can be chosen to achieve an RMS error of ε at an expected cost

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

Multilevel Monte Carlo

The standard estimator for SDE applications is

$$\widehat{Y}_{\ell} = N_{\ell}^{-1} \sum_{n=0}^{N_{\ell}} \left(\widehat{P}_{\ell}(W^{(n)}) - \widehat{P}_{\ell-1}(W^{(n)}) \right)$$

using the same Brownian motion $W^{(n)}$ for the n^{th} sample on the fine and coarse levels.

However, there is some freedom in how we construct the coupling provided \widehat{Y}_ℓ is an unbiased estimator for $\mathbb{E}[\widehat{P}_\ell - \widehat{P}_{\ell-1}]$.

Also, uniform timestepping is not required – it is fairly straightforward to implement MLMC using non-nested adaptive timestepping.

(G, Lester, Whittle: MCQMC14 proceedings)



Position x_t , and velocity u_t subject to steady and stochastic forcing:

$$du_t = a(x_t, u_t, t) dt + b(x_t, t) dw_t$$

$$dx_t = u_t dt$$

Domain $x \ge 0$, with reflection so that when it hits x = 0 at time τ then the velocity is reflected, so

$$u_{\tau^+} = -u_{\tau^-}.$$

Euler-Maruyama treatment with uniform timestep *h*:

$$\widehat{u}_{n+1} = s_n (\widehat{u}_n + a(\widehat{x}_n, \widehat{u}_n, t) h + b(\widehat{x}_n, t_n) \Delta w_n)
\widehat{x}_{n+1} = s_n (\widehat{x}_n + \widehat{u}_n h)$$

with $s_n = \pm 1$ chosen so that $\widehat{x}_{n+1} \geq 0$.

Problem: only $O(h^{1/2})$ strong convergence

Reason: doesn't account for reflection occurring part-way through a timestep.

Key idea: if A(X, U, t), B(X, t) are sufficiently smooth, get O(h) convergence using an extended domain:

$$dU_t = A(X_t, U_t, t) dt + B(X_t, t) dW_t$$

$$dX_t = U_t dt,$$

with

$$A(X, U, t) = \begin{cases} a(X, U, t), & X \ge 0 \\ -a(-X, -U, t), & X < 0 \end{cases}$$

$$B(X, t) = \begin{cases} b(X, t), & X \ge 0 \\ b(-X, t), & X < 0 \end{cases}$$

and then take x = |X| as output.



Why does that give O(h) strong convergence, but the original doesn't?

If we define

$$\left(\begin{array}{c} u_t \\ x_t \end{array}\right) = S(X_t) \ \left(\begin{array}{c} U_t \\ X_t \end{array}\right),$$

where $S(X) \equiv \operatorname{sign}(X)$, then u_t, x_t satisfy

$$du_t = a(x_t, u_t, t) dt + b(x_t, t) S(X_t) dW_t$$

$$dx_t = u_t dt,$$

By setting $dw_t = S(X_t) dW_t$, we see that this is equivalent in distribution to the original model problem.

Note: strong convergence is now at fixed W_t – not the same as fixed w_t .

New MLMC treatment:

$$\widehat{u}_{n+1}^{p} = \widehat{u}_{n} + a(\widehat{x}_{n}, \widehat{u}_{n}, t_{n}) h + b(\widehat{x}_{n}, t_{n}) \widehat{s}_{n} \Delta W_{n}
\widehat{x}_{n+1}^{p} = \widehat{x}_{n} + \widehat{u}_{n} h$$

followed by a correction/reflection step:

$$\begin{array}{rcl} \widehat{u}_{n+1} & = & \mathrm{sign}(\widehat{x}_{n+1}^{\rho}) \ \widehat{u}_{n+1}^{\rho} \\ \widehat{x}_{n+1} & = & \mathrm{sign}(\widehat{x}_{n+1}^{\rho}) \ \widehat{x}_{n+1}^{\rho} \\ \widehat{s}_{n+1} & = & \mathrm{sign}(\widehat{x}_{n+1}^{\rho}) \ \widehat{s}_{n} \end{array}$$

with same Brownian path for coarse and fine levels.

Can show that when a and b are both constant, the coarse and fine paths are identical at coarse timesteps.

Test case 1:

$$x_0 = 0.2$$
, $u_0 = -0.2$, $a(x, t) = 0$, $b(x, t) = 0.5$.

in domain $0 \le x \le 1$, with reflection at both boundaries.

Output of interest: $\int_0^1 x_t dt$ approximated by $\sum_{n=1}^{2^n} h_\ell \widehat{x}_n$.

Test case 2: changes drift, volatility to

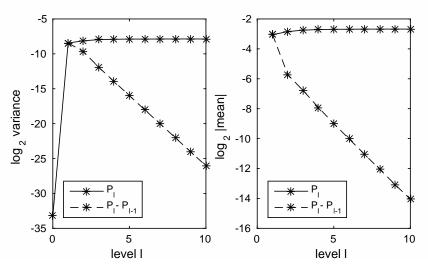
$$a(x, t) = -0.2, b(x, t) = 0.5 + 0.5 x.$$

- standard O(h) numerical analysis no longer applies



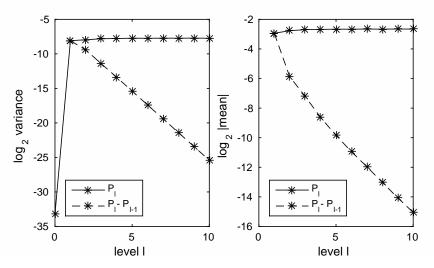
$$\mathbb{V}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] \sim h_{\ell}^2$$

$$\mathbb{E}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] \sim h_{\ell}$$



$$\mathbb{V}[\widehat{P}_{\ell}\!-\!\widehat{P}_{\ell-1}]\sim h_{\ell}^2$$

$$\mathbb{E}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] \sim h_{\ell}$$



Without mass, the SDE is

$$dx_t = a(x_t, t) dt + b(x_t, t) dw_t$$

and if the domain is $x \ge 0$, particles are prevented from crossing x = 0.

Euler-Maruyama treatment with uniform timestep h:

$$\widehat{x}_{n+1} = \left| \widehat{x}_n + a(\widehat{x}_n, t) h + b(\widehat{x}_n, t_n) \Delta w_n \right|$$

Again only $O(h^{1/2})$ strong convergence, even when b is uniform

Thinking about the extended domain leads to

$$dx_t = a(x_t, t) dt + b(x_t, t) S(X_t) dW_t$$

where $S(X) \equiv \operatorname{sign}(X)$, and hence the numerical approximation is

$$\widehat{x}_{n+1}^p = \widehat{x}_n + a(\widehat{x}_n, t_n) h + b(\widehat{x}_n, t_n) \widehat{s}_n \Delta W_n$$

followed by a correction/reflection step:

$$\widehat{x}_{n+1} = \operatorname{sign}(\widehat{x}_{n+1}^p) \widehat{x}_{n+1}^p$$

$$\widehat{s}_{n+1} = \operatorname{sign}(\widehat{x}_{n+1}^p) \widehat{s}_n$$

with same Brownian path for coarse and fine levels.

Note: if b is not uniform then we need to use first order Milstein approximation to get O(h) strong convergence.



Test case 1:

$$x_0 = 0.2$$
, $a(x, t) = 0$, $b(x, t) = 0.5$.

in domain $0 \le x \le 1$, with reflection at both boundaries.

Output of interest: $\int_0^1 x_t dt$ approximated by $\sum_{n=1}^{2^c} h_\ell \widehat{x}_n$.

Test case 2: changes drift, volatility to

$$a(x, t) = -0.2, b(x, t) = 0.5 + 0.5 x.$$

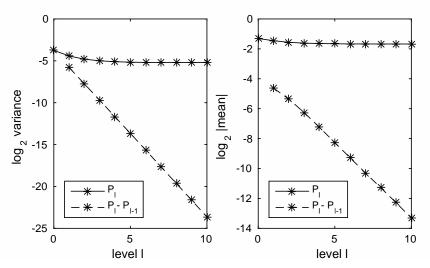
- standard O(h) numerical analysis no longer applies



Test case 1:

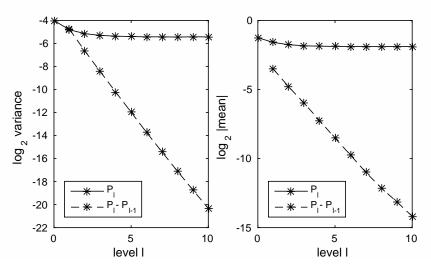
$$\mathbb{V}[\widehat{P}_{\ell}\!-\!\widehat{P}_{\ell-1}]\sim h_{\ell}^2$$

$$\mathbb{E}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] \sim h_{\ell}$$



$$\mathbb{V}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] \sim h_{\ell}^{3/2}$$

$$\mathbb{E}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] \sim h_{\ell}$$



Why is the variance $O(h^{3/2})$?

Ad-hoc explanation:

- O(1) path density near x=0
- $O(h^{1/2})$ movement in each timestep
- ullet $\Longrightarrow O(h^{1/2})$ probability of crossing boundary in each timestep
- ullet $\Longrightarrow O(h^{-1/2})$ total crossings per path
- ullet each crossing gives error which is O(h) but has near-zero mean
- if crossings are approximately independent, then

$$\mathbb{V}[\widehat{P}_{\ell} - \widehat{P}_{\ell-1}] = O(h^{-1/2} \times h^2) = O(h^{3/2})$$

Note: in the case with mass, the velocity is O(1), the movement in each timestep is O(h), so the number of crossings is $O(1) \implies V_{\ell} = O(h^2)$.

Start with particles with mass, and 2D domain with boundary at x=0.

lf

$$a = \begin{pmatrix} a_x(x,y) \\ a_y(x,y) \end{pmatrix}, \quad b = I,$$

then for x < 0 extended domain SDE naturally has

$$A = \begin{pmatrix} -a_x(-x,y) \\ a_y(-x,y) \end{pmatrix}, \quad B = I,$$

This leads to an SDE in the regular half-plane in which the x-component of $\mathrm{d}W$ is "flipped" each time the boundary is hit.

However, if

$$a = \left(\begin{array}{c} a_X(x,y) \\ a_Y(x,y) \end{array} \right), ~~ b = \left(\begin{array}{cc} 1 & 0 \\ \beta & 1 \end{array} \right),$$

then when x < 0

$$A = \begin{pmatrix} -a_x(-x,y) \\ a_y(-x,y) \end{pmatrix}, \quad B = \begin{pmatrix} 1 & 0 \\ -\beta & 1 \end{pmatrix},$$

because if X = -x then

$$\mathbb{E}[\,\mathrm{d} X\,\mathrm{d} y\,] = -\,\mathbb{E}[\,\mathrm{d} x\,\mathrm{d} y\,].$$

This gives a discontinuity in B, and there seems no way to get O(h) strong convergence.



Hence, multi-dimensional extension for particles with mass will only work in simple cases.

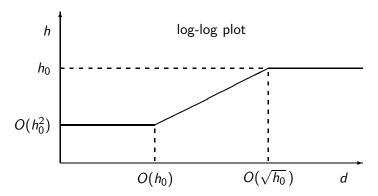
For massless particles, there is extra complication of oblique reflections.

If diffusion is isotropic (i.e. b is a multiple of the identity matrix) and reflections are normal, then can probably get O(h) strong convergence.

Otherwise, probably not.

What else can we do for massless case?

Go back to starting point with standard reflection treatment, and use adaptive timesteps based on distance d to boundary.



 $O(h_0^2)$ timestep near boundary $\Longrightarrow O(h_0)$ strong error

In MLMC, on level ℓ would use

$$h = \min \left(2^{-\ell} h_0, \; \max \left(2^{-2\ell} h_0, (\frac{1}{3} d / \|b\|_2)^2 \right) \right)$$

where d is the distance to the boundary

- $2^{-\ell}h_0$ in interior
- $2^{-2\ell}h_0$ next to boundary
- $(\frac{1}{3}d/\|b\|_2)^2$ in layer in-between

The factor $\frac{1}{3}$ implies $3\sqrt{h} \|b\|_2 < d$ so that boundary crossings from the intermediate zone (or interior zone) are unlikely.

Most will come from boundary zone, with resultant $O(2^{-\ell})$ strong error.

Key observation: the computational cost is proportional to $\int_D h^{-1} dx$ and this is $O(2^\ell)$.

Hence, in the usual MLMC theorem, we should get

$$\alpha = 1, \quad \beta = 2, \quad \gamma = 1,$$

and hence obtain $O(\varepsilon^{-2})$ complexity.

Conclusions

- simple reflection "trick" improves the MLMC variance for 1D reflected diffusions, for particles with or without mass
- the extension to multiple dimensions should work in simple cases, but not in more general cases
- more difficult cases can use adaptive timestepping

Webpages: http://people.maths.ox.ac.uk/gilesm/