

Multilevel Monte Carlo for elliptic SPDEs

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Outline

- standard Monte Carlo simulation
- multilevel Monte Carlo simulation
- elliptic SPDE application
- conclusions

Monte Carlo simulation

In many applications want to estimate $\mathbb{E}[P(\omega)]$ where $\omega \in \Omega$ is an infinite-dimensional random variable.

- computational finance:
 - ω represents W_t the driving Brownian motion in an SDE (stochastic differential equation)
 - P is the financial payoff function
- simulation of oil reservoirs & nuclear waste repositories:
 - ω represents $k(x)$, the diffusivity in an elliptic SPDE
 - $$-\nabla \cdot (k(x) \nabla p) = 0$$
 - P might be the flux of oil or contaminants across some boundary

Monte Carlo simulation

In MC simulation we estimate the expectation using

$$\hat{Y} = N^{-1} \sum_{n=1}^N \hat{P}(\omega^{(n)})$$

where $\omega^{(n)}$ are N independent samples

Note there are two sources of error here:

- *sampling error* due to the finite number of samples
- *bias* because $\hat{P}(\omega)$ is an approximation to $P(\omega)$ due to
 - discretisation error (finite timesteps, finite grid size)
 - finite dimensional approximation to ω

Monte Carlo simulation

The mean square error is

$$\begin{aligned}\mathbb{E} \left[\left(\hat{Y} - \mathbb{E}[P] \right)^2 \right] &= \mathbb{E} \left[\left(\hat{Y} - \mathbb{E}[\hat{Y}] + \mathbb{E}[\hat{Y}] - \mathbb{E}[P] \right)^2 \right] \\ &= \mathbb{E} \left[\left(\hat{Y} - \mathbb{E}[\hat{Y}] \right)^2 \right] + \left(\mathbb{E}[\hat{Y}] - \mathbb{E}[P] \right)^2 \\ &= \mathbb{V}[\hat{Y}] + \left(\mathbb{E}[\hat{Y}] - \mathbb{E}[P] \right)^2 \\ &= N^{-1} \mathbb{V}[\hat{P}] + \left(\mathbb{E}[\hat{P}] - \mathbb{E}[P] \right)^2\end{aligned}$$

- first term is due to sampling error
- second term is due to bias

Monte Carlo simulation

To achieve RMS accuracy of ε requires:

- $N = O(\varepsilon^{-2})$

- bias = $O(\varepsilon)$

In a d -dimensional SPDE application with grid spacing h , if the bias is $O(h^\alpha)$ then need $h = O(\varepsilon^{1/\alpha})$, and total cost is $O(\varepsilon^{-(2+d/\alpha)})$, assuming efficient multigrid solution

(very challenging because of very rough coefficients
– Graham & Scheichl)

To get acceptable accuracy in 3D applications may need 10,000 simulations on a 128^3 grid \implies very expensive

Multilevel Monte Carlo

The multilevel objective is to greatly reduce this cost:

dim	$\alpha = 1$		$\alpha = 2$	
	MC	MLMC	MC	MLMC
1	ε^{-3}	ε^{-2}	$\varepsilon^{-2.5}$	ε^{-2}
2	ε^{-4}	$\varepsilon^{-2}(\log \varepsilon)^2$	ε^{-3}	$\varepsilon^{-2}(\log \varepsilon)^2$
3	ε^{-5}	ε^{-3}	$\varepsilon^{-3.5}$	$\varepsilon^{-2.5}$

How? Use multigrid philosophy:

- fine grid accuracy at coarse grid cost
- geometric sequence of grids

but no iteration in Monte Carlo simulation?

Multilevel Monte Carlo

Consider Monte Carlo simulations with different levels of refinement, $l = 0, 1, \dots, L$, with level L being the finest.

If \hat{P}_l is the approximation of P on level l , then

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

Idea is to independently estimate each of the terms on the r.h.s., in a way which minimises the overall variance for a fixed computational cost.

Finest level is still the same, but will use very few samples at that level.

Multilevel Monte Carlo

Simplest estimator for $\mathbb{E}[\hat{P}_l - \hat{P}_{l-1}]$ for $l > 0$ is

$$\hat{Y}_l = N_l^{-1} \sum_{n=1}^{N_l} \left(\hat{P}_l^{(n)} - \hat{P}_{l-1}^{(n)} \right)$$

using same stochastic sample $\omega^{(n)}$ for both levels

Variance is $N_l^{-1} V_l$ where $V_l = \mathbb{V}[\hat{P}_l - \hat{P}_{l-1}]$

Key point: V_l gets progressively smaller as l increases because \hat{P}_l, \hat{P}_{l-1} both accurately approximate P for same ω

Multilevel Monte Carlo

If C_l is cost of one sample on level l , the variance of the

combined estimator is $\sum_{l=0}^L N_l^{-1} V_l$ and its computational

cost is $\sum_{l=0}^L N_l C_l$ so the variance is minimised for fixed cost

by choosing $N_l \propto \sqrt{V_l/C_l}$, and then the cost on level l is

proportional to $N_l C_l \propto \sqrt{V_l C_l}$

To make RMS error ε

- choose constant of proportionality so variance is $\frac{1}{2} \varepsilon^2$
- choose L so that $\left(\mathbb{E}[\hat{P}_l] - \mathbb{E}[P] \right)^2 < \frac{1}{2} \varepsilon^2$

MLMC Theorem

If there exist independent estimators \widehat{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

$$\text{i) } \left| \mathbb{E}[\widehat{P}_l - P] \right| \leq c_1 2^{-\alpha l}$$

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

$$\text{iii) } \mathbb{V}[\widehat{Y}_l] \leq c_2 N_l^{-1} 2^{-\beta l}$$

$$\text{iv) } C_l \leq c_3 2^{\gamma l}$$

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

$$\hat{Y} = \sum_{l=0}^L \hat{Y}_l,$$

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

with a computational cost C with bound

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

Papers

- My first paper (*Operations Research, 2006 – 2008*) applied idea to SDE path simulation, and proved slightly less general form of the theorem
- Second paper (*MCQMC 2006*) improved multilevel variance convergence using better discretisation
- Third paper with D. Higham & X. Mao (*Finance and Stochastics, 2009*) performed numerical analysis of discretisation in first paper
- New paper with K. Debrabant and A. Rößler analyses discretisation in second paper

Multilevel method is a generalisation of two-level control variate method of Kebaier (2005), and related to multilevel parametric integration by Heinrich (2001).

Elliptic SPDE

We consider the elliptic PDE

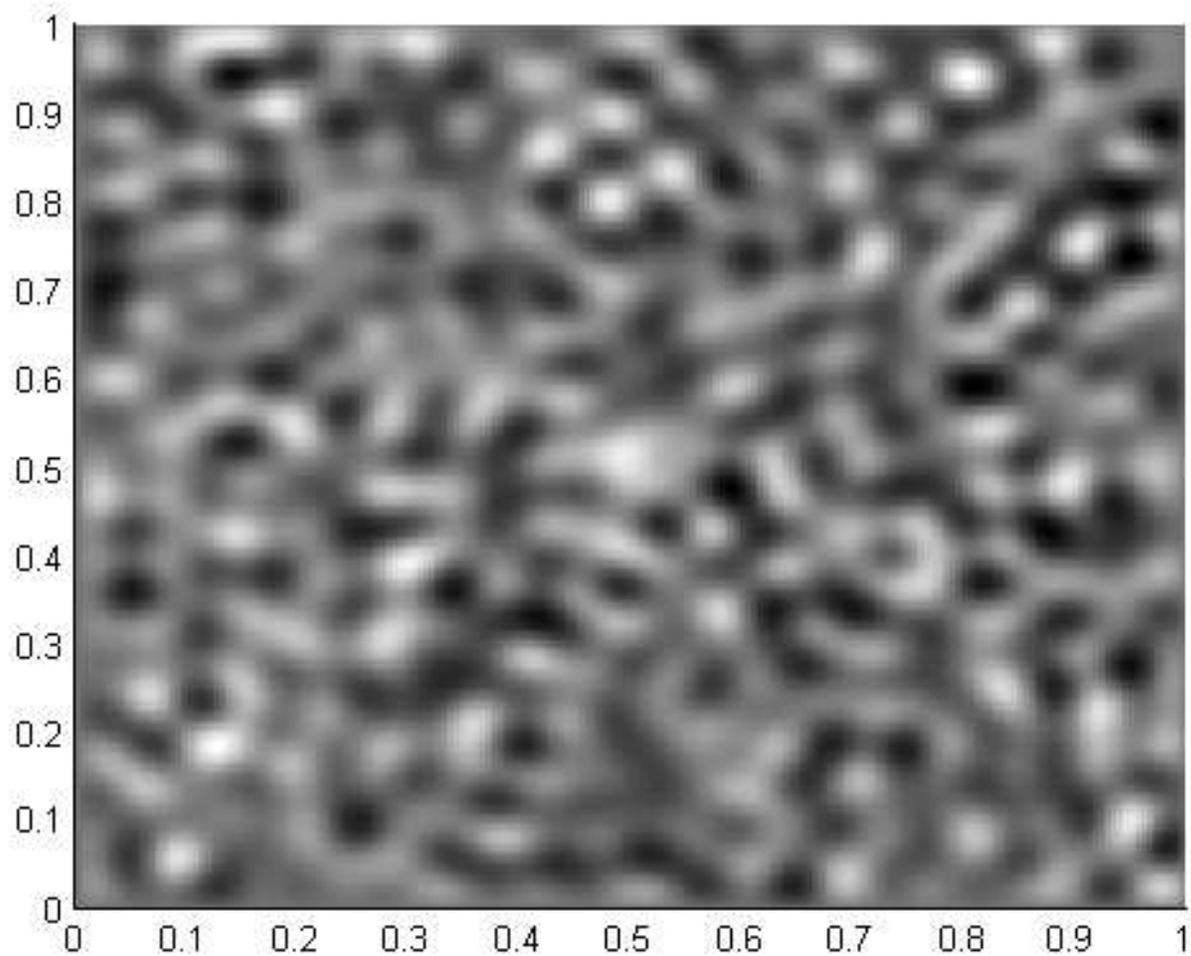
$$-\nabla \cdot (k(x, \omega) \nabla p(x, \omega)) = f(x, \omega), \quad x \in D,$$

with **random coefficient** $k(x, \omega)$ and random data $f(x, \omega)$.

We model k as a **lognormal random field**, i.e. $\log k$ is a Gaussian field with mean 0 and covariance function

$$R(x, y) = \sigma^2 \exp(-\|x - y\|/\lambda)$$

Elliptic SPDE



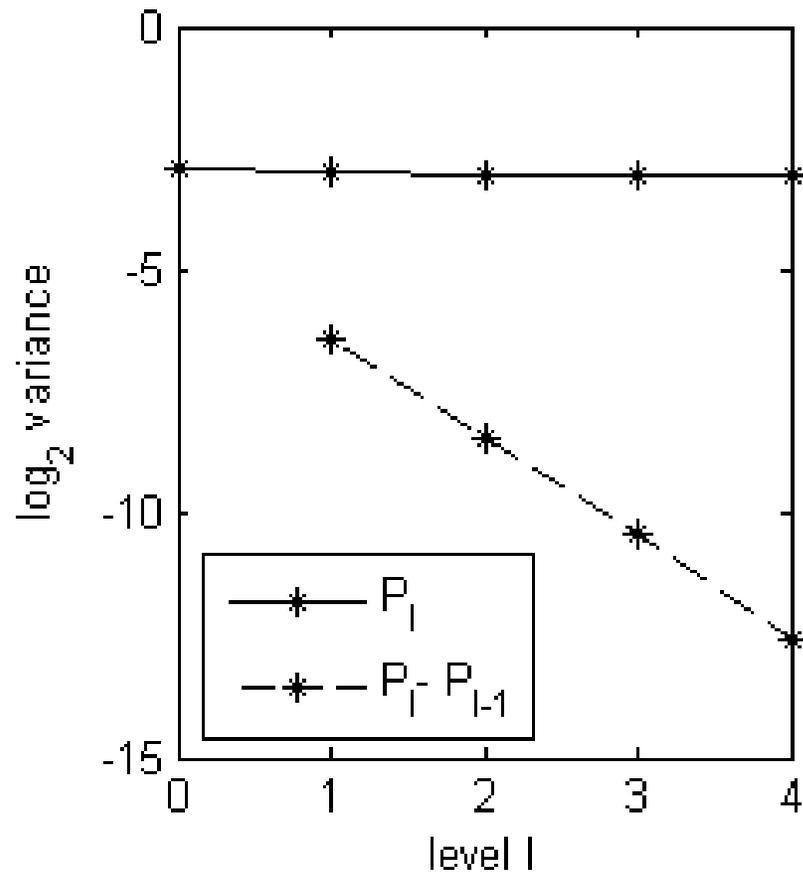
A typical realisation for $D = [0, 1]^2$, $\lambda = 0.001$, $\sigma^2 = 1$.

Elliptic SPDE

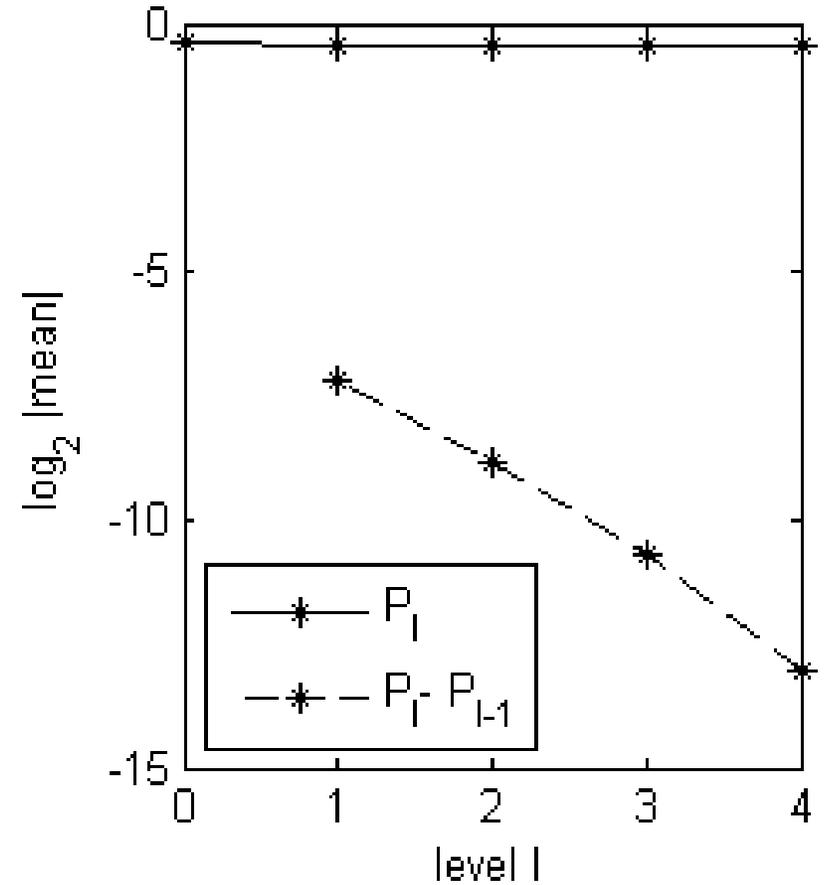
Discretisation:

- cell-centred Finite Volume discretisation on a uniform grid \mathcal{T}_h – for rough coefficients we need to make h very small
- sampling of the random coefficient currently based on truncated Karhunen-Lòeve expansion, evaluated at cell centres – but the method of sampling is not essential to the algorithm
- each level of refinement has twice as many grid points in each direction

1D Results

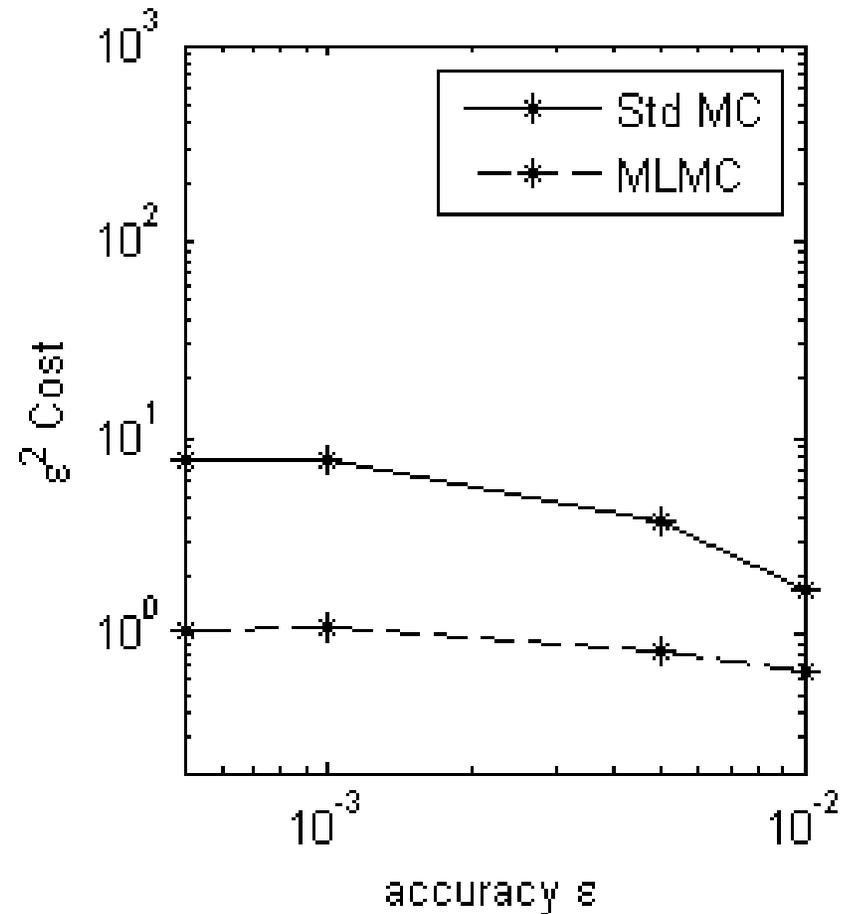
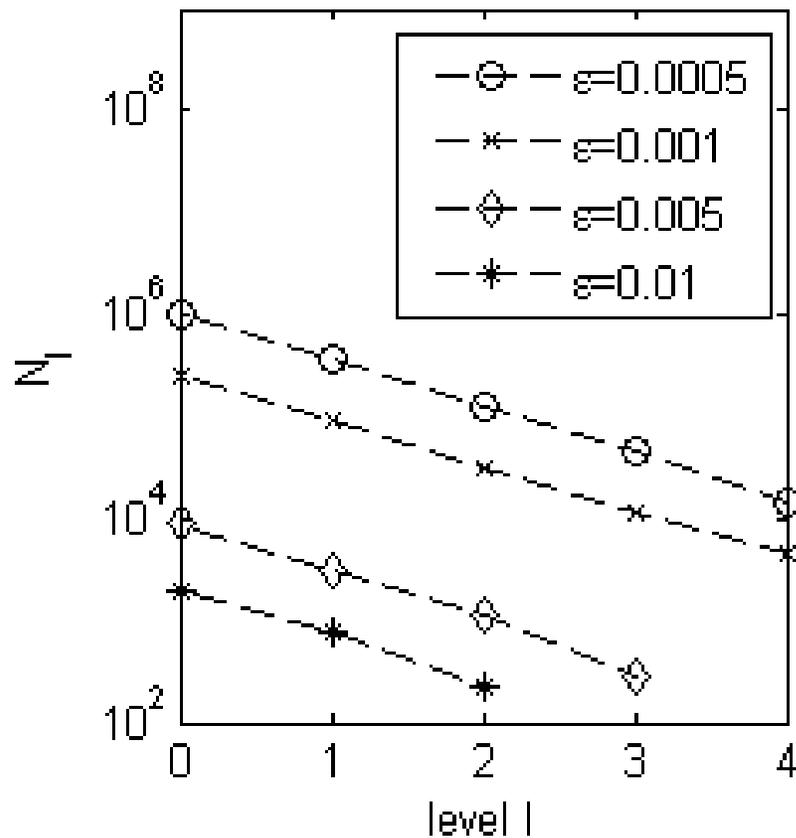


$$\mathbb{V}[\hat{P}_l - \hat{P}_{l-1}] \sim h_l^2,$$

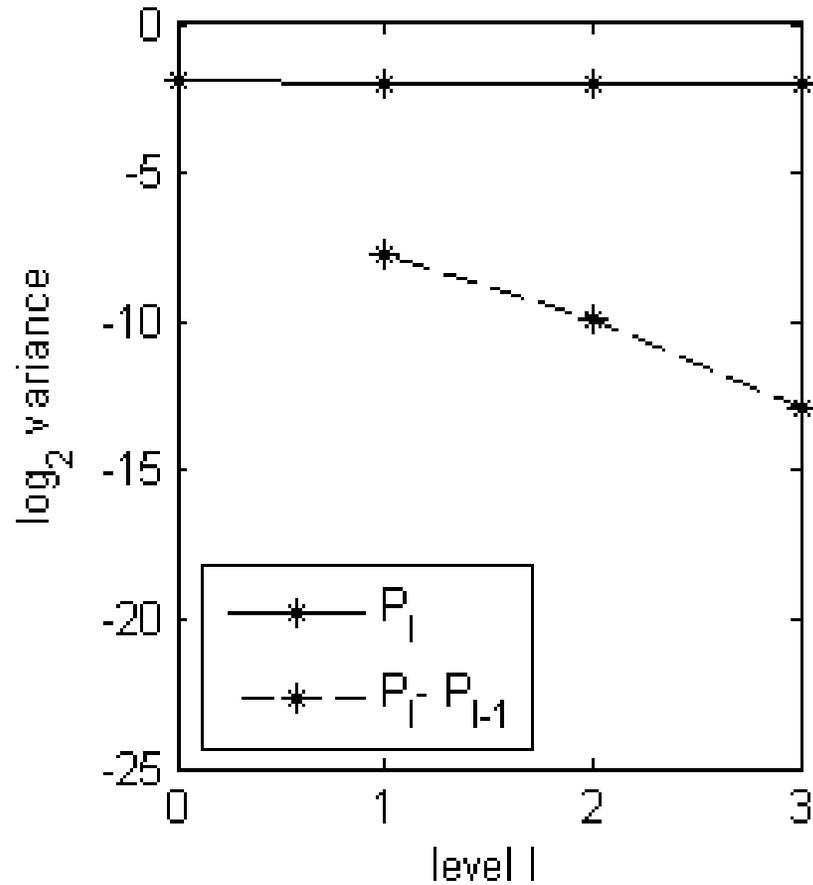


$$\mathbb{E}[\hat{P}_l - \hat{P}_{l-1}] \sim h_l^2$$

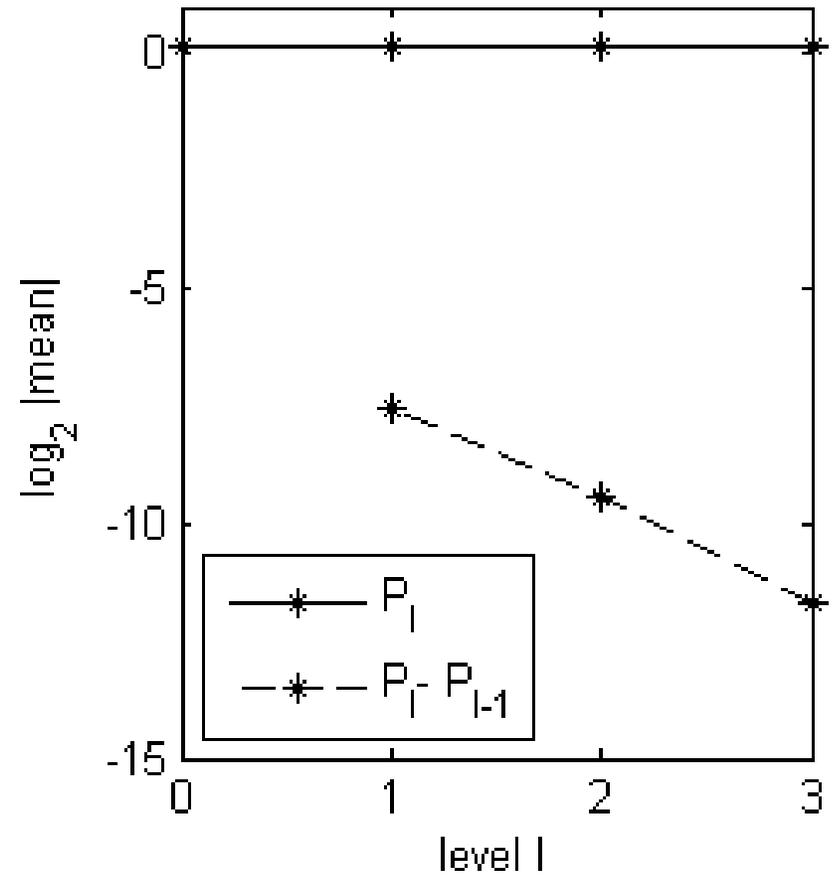
1D Results



2D Results

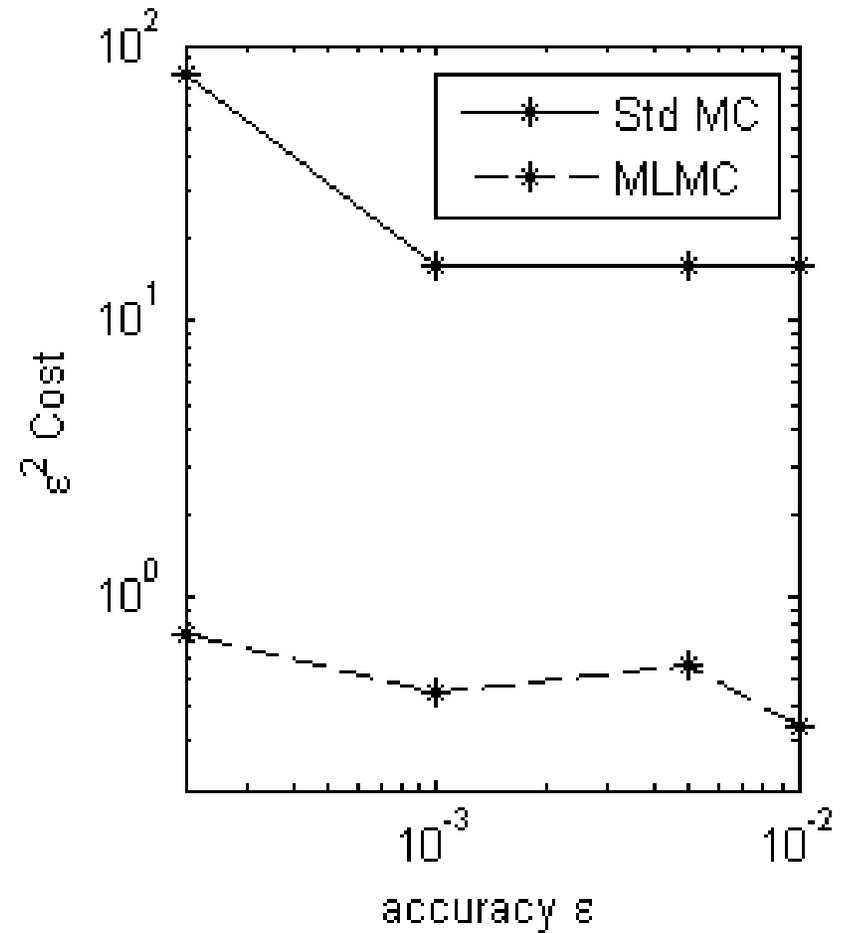
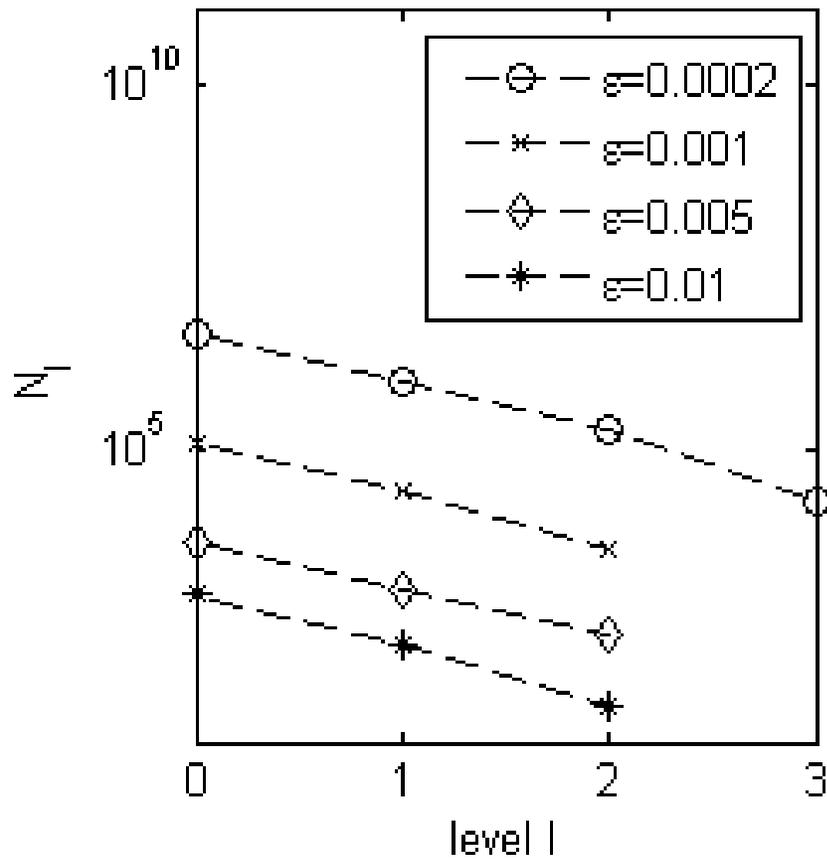


$$\mathbb{V}[\hat{P}_l - \hat{P}_{l-1}] \sim h_l^2,$$



$$\mathbb{E}[\hat{P}_l - \hat{P}_{l-1}] \sim h_l^2$$

2D Results



Conclusions

- standard Monte Carlo is prohibitively expensive for 2D and 3D elliptic SPDE applications
- multilevel Monte Carlo greatly reduces the cost, making this feasible for engineering applications
- we believe it is a viable competitor to polynomial chaos approach, particularly for applications with minimal spatial correlation
- numerical analysis is very hard, but we're making some headway with finite element analysis, at least to gain insight into its effectiveness
- future work will look at combining the multilevel approach with quasi-Monte Carlo sampling – has been very effective for SDE applications in finance