

Risk analysis and high-dimensional integrals

Alexander Podkopaev, Nikita Puchkin, Igor Silin

SKOLKOVO INSTITUTE OF SCIENCE AND TECHNOLOGY

December 10, 2016

1 Description of algorithm and PDE solver

We implement a solver for numerical solution of PDE

$$\begin{aligned}\frac{\partial}{\partial t}u(x, t) &= \sigma \frac{\partial^2}{\partial x^2}u(x, t) - V(x, t)u(x, t), \quad t \in [0, T], x \in \mathbb{R} \\ u(x, 0) &= f(x)\end{aligned}$$

The exact solution is given by the Feynman-Kac formula

$$u(x, T) = \int_{C\{x, 0; T\}} f(\xi(T)) \exp \left\{ - \int_0^T V(\xi(\tau), T - \tau) d\tau \right\} \mathcal{D}_\xi,$$

where the integration is done over the set $C\{x, 0; T\}$ of all continuous paths $\xi(T) : [0, T] \rightarrow \mathbb{R}$ from the Banach space $\Xi([0, T], \mathbb{R})$ starting at $\xi(0) = x$ and stopping at arbitrary endpoints at time T . \mathcal{D}_ξ is the Wiener measure and $\xi(t)$ is the Wiener process.

For numerical computation one can break the time range $[0, T]$ into n intervals by points

$$\tau_k = k\delta t, \quad 0 \leq k < n, \quad n : \tau_n = T$$

The average path of a Brownian particle $\xi(\tau_k)$ after k steps is defined as

$$\xi^{(k)} = \xi(\tau_k) = x + \xi_1 + \dots + \xi_k,$$

where every random step ξ_i , $1 \leq i \leq k$, is independently taken from a normal distribution $\mathcal{N}(0, 2\sigma\delta t)$. By definition $\xi^{(0)} = x$.

On the introduced uniform grid one approximates

$$\Lambda(T) = \int_0^T V(\xi(\tau), T - \tau) d\tau \simeq \sum_{i=0}^n w_i V_i^{(n)} \delta t, \quad V_i^{(n)} \equiv V(\xi(\tau_i), \tau_{n-i}),$$

where the set of weights $\{w_i\}_{i=0}^n$ is taken according to trapezoid rule or Simpson rule. Then

$$\exp\{-\Lambda(T)\} \simeq \prod_{i=0}^n \exp\{-w_i V_i^{(n)} \delta t\}$$

The Wiener measure transforms to n -dimensional measure

$$\mathcal{D}_\xi^{(n)} = \left(\frac{\lambda}{\pi}\right)^{\frac{n}{2}} \prod_{k=1}^n \exp\{-\lambda \xi_k^2\} d\xi_k, \quad \lambda = \frac{1}{4\sigma\delta t},$$

and a numerical approximation of the exact solution can be written in the form

$$u^{(n)}(x, T) = \int_{-\infty}^{\infty} \mathcal{D}_{xi} f(\xi^{(n)}) \prod_{i=0}^n e^{-w_i v_i^{(n)} \delta t}$$

The multidimensional integral can be represented in terms of n one-dimensional convolutions. Define

$$F_k^{(n)}(x) = \sqrt{\frac{\lambda}{\pi}} \int_{-\infty}^{\infty} \Phi_{k+1}^{(n)}(x + \xi) e^{-\lambda \xi^2} d\xi, \quad x \in \mathbb{R}, \quad k = n, n-1, \dots, 1,$$

where

$$\Phi_{k+1}^{(n)}(x) = F_{k+1}^{(n)}(x) \exp\{-w_k V(x, \tau_{n-k}) \delta t,$$

and

$$F^{(n)}(x)_{n+1} = f(x)$$

Then the numerical solution is given by formula

$$u^{(n)}(x, T) = F_1^{(n)}(x) e^{-w_0 V(x, T) \delta t}$$

Since $F_k^{(n)}(x)$ is represented as integral over all real values, it is replaced by an integral over a segment

$$F_k^{(n)}(x) \simeq \tilde{F}_k^{(n)}(x) = \sqrt{\frac{\lambda}{\pi}} \int_{-a_x}^{a_x - h_x} \Phi_{k+1}^{(n)}(x + \xi) e^{-\lambda \xi^2} d\xi$$

The function $F_k^{(n)}(x)$ is computed on the uniform mesh

$$x_i^{(k)} = -ka_x + ih_x, \quad 0 \leq i \leq kM, \quad h_x = \frac{a_x}{N_x}, \quad M = 2N_x$$

and the integration mesh is taken with the same step h_x

$$\xi_j = -a_x + jh_x, \quad 0 \leq j < M$$

Then

$$\begin{aligned} \tilde{F}_k^{(n)}(x_i^{(k)}) &\simeq \sum_{j=0}^{M-1} \mu_j \Phi_{k+1}^{(n)}(x_{i+j}^{(k+1)}) p(\lambda, \xi_j), \quad p(\lambda, \xi) = \sqrt{\frac{\lambda}{\pi}} e^{-\lambda \xi^2} \\ \Phi_{k+1}^{(n)}(x_i^{(k+1)}) &= \tilde{F}_{k+1}^{(n)}(x_i^{(k+1)}) \exp\{-w_k V(x_i^{(k+1)}, \tau_{n-k}) \delta t \end{aligned}$$

or in the matrix form

$$\begin{aligned} \tilde{F}_k^{(n)} &= \Phi_{k+1}^{(n)} \circ \tilde{\mu}, \quad \tilde{\mu}_j = \mu_j p(\lambda, \xi_j) \\ \Phi_{k+1}^{(n)} &= \tilde{F}_{k+1}^{(n)} \exp\{-w_k V(x^{(k+1)}, \tau_{n-k}) \delta t, \end{aligned}$$

where $a \circ b$ denotes a convolution of vectors $a \in \mathbb{R}^m$ and $b \in \mathbb{R}^k$, i. e. a vector $c \in \mathbb{R}^{m+k-1}$, such that

$$c_i = \sum_{j=0}^{k-1} a_{i+j} b_j, \quad a_i = 0, \quad \forall i : (i < 0) \vee (i \geq m)$$

The algorithm for computation of $u^{(n)}(x, T)$ is following.

1. Given T , choose a time step δt and the number of steps $n \frac{T}{\delta t}$.
2. Create 1D array τ of size n , $\tau_i = i\delta t$.
3. Create 1D array w of size $(n + 1)$, corresponding to the set of weights in trapezoid or Simpson rule.
4. Choose a_x , size of coordinate grid $M = 2N_x$ and a coordinate step $h_x = \frac{a_x}{N_x}$.
5. Initialize 1D array $\tilde{F}_{n+1}^{(n)}$ of size $(n + 1)M$, where $F_{n+1}^{(n)} = f(x)$ and $x_i = -(n + 1)a_x + ih_x$, $0 \leq i < (n + 1)M$.
6. For $k = n, n - 1, \dots, 1$ do
 - (a) Create 1D array $x^{(k+1)}$ of size $(k + 1)M$, $x_i^{(k+1)} = -(k + 1)a_x + ih_x$, $0 \leq i < (k + 1)M$
 - (b) Create 1D array of size $(k + 1)M$ $e^{-w_k V(x^{(k)}, \tau_{n-k})\delta t}$.
 - (c) Create 1D array $\Phi_{k+1}^{(n)} = F_{k+1}^{(n)} \odot e^{-w_k V(x^{(k+1)}, \tau_{n-k})\delta t}$.
 - (d) Create 1D array μ of size $M + 1$, corresponding to the set of weights.
 - (e) Create 1D array ξ of size $M + 1$, where $\xi_j = -a_x + jh_x$, $0 \leq j < M + 1$.
 - (f) Create 1D array $\tilde{\mu} = \mu \odot p(\lambda, \xi)$.
 - (g) Compute convolution $\tilde{F}_k^{(n)} = \Phi_{k+1}^{(n)} \circ \tilde{\mu}$. $\tilde{F}_k^{(n)}$ is 1D array of size kM .
7. Create 1D array $x^{(1)}$ of size M , $x_i^{(1)} = -a_x + ih_x$, $0 \leq i < M$
8. Create 1D array of size M $e^{-w_0 V(x^{(1)}, T)\delta t}$.
9. Compute the numerical solution $u^{(n)} = F_1^{(n)} \odot e^{-w_0 V(x^{(1)}, T)\delta t}$.

The proposed algorithm is implemented in class 'PDE_Solver'. An object of the class has following attributes:

- sigma
- V
- f
- a_x
- M
- h_x
- T
- n
- delta_t
- u – numerical solution

Names of all attributes correspond to notations used in this paper. Methods of an object of the class 'PDE_solver' are following:

- `Set_Limits(float a)` – sets $a_x = a$.
- `Convolve(1D array a, 1D array b)` – returns a 1D array $c = a \circ b$.
- `Convolve_Low-rank(1D array a, 1D array b)` – returns a 1D array $c = a \circ b$ using low-rank cross approximation.
- `Solve()` – computes a numerical solution u according to the proposed algorithm.