Randomized Markovian approximation for rough volatility models

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Overview

- Topic: Markovian Approximation for Rough Volatility Models
- Contribution: Introduce stochastic nodes and weights.
 - Nodes and weights are chosen for each sample path.
 - L²-error minimization reduces to minimizing the time-integrated variance of a certain random variable.
 - This framework naturally incorporates variance reduction techniques.
- **Results:** When the number of nodes is small, this approach outperforms all other deterministic methods (except BL2).

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Classical Stochastic Volatility Models

Volatility follows a stochastic differential equation (SDE):

$$V_t = V_0 + \int_0^t b(V_s) \,\mathrm{d}s + \int_0^t \sigma(V_s) \,\mathrm{d}W_s. \tag{1}$$

- Markovian: Future volatility depends only on its current state.
- Example: Heston model, SABR model

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Rough Volatility Models (RVM)

Volatility follows a stochastic Volterra equation:

$$V_t = V_0 + \int_0^t \frac{K(t-s)b(V_s) \,\mathrm{d}s}{+} \int_0^t \frac{K(t-s)\sigma(V_s) \,\mathrm{d}W_s}{,} \quad (2)$$

where the kernel K(t) is typically given by the **fractional kernel**:

$$K(t) = rac{t^{H-1/2}}{\Gamma(H+1/2)}, \quad 0 < H < 1/2.$$
 (3)

- Reduces to a **fractional Brownian motion** with Hurst parameter *H* when b = 0 and $\sigma = 1$.
- Non-Markovian: Exhibits memory effects.
- Highly irregular and rougher than classical models.

• If
$$H = \frac{1}{2}$$
, the kernel reduces to $K(t) \equiv 1$.

Example: Rough Heston Model

The Rough Heston Model extends the classical Heston model:

$$dS_t = S_t \sqrt{V_t} dB_t, \quad S_0 > 0,$$

$$V_t = V_0 + \int_0^t \frac{K(t-s)(\theta - \lambda V_s)}{(\theta - \lambda V_s)} ds + \int_0^t \frac{K(t-s)\eta}{\sqrt{V_s}} dW_s,$$
(5)

where W_t and B_t are ρ -correlated Brownian motions.

 Admits a semi-analytical characteristic function [El Euch and Rosenbaum, 2019], allowing Fourier pricing techniques.

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Volatility Path V_t in the Rough Heston Model



Why RVM?

- Empirical Motivation: Empirical studies [Gatheral et al., 2018] show that volatility exhibits rough behavior, resembling fractional Brownian motion.
- Better Fit for IV Smiles: RVM explains the steep implied volatility (IV) smiles observed in short-maturity options [Fukasawa, 2021].

Challenges in RVM

Non-Markovian Nature:

The fractional kernel K(t) introduces memory effects, making the process V_t depend on its entire past.

Significantly increases the computational cost

Addressing the Non-Markovian Challenge

Key Idea:

Approximate the fractional kernel K(t) using a sum of exponentials:

$$K(t) = \int_0^\infty c_H y^{-H - \frac{1}{2}} e^{-yt} \, \mathrm{d}y \approx \sum_{i=1}^n w_i e^{-y_i t} =: K^n(t).$$
(6)

- This approximation, known as a **Markovian approximation** [Abi Jaber and El Euch, 2019], transforms the original model into an *n*-dimensional Markov process.
- Enable the use of **standard numerical methods** for efficient simulation.

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Markovian Approximation

The Markovian approximation is defined as:

$$V_t^n = V_0 + \int_0^t \frac{\mathcal{K}^n(t-s)b(V_t^n) \,\mathrm{d}t}{t} + \int_0^t \frac{\mathcal{K}^n(t-s)\sigma(V_t^n) \,\mathrm{d}W_t}{t} \quad (7)$$

 V_t^n is a *n*-dimensional Markovian process since:

• V_t^n can be expressed as $V_t^n = V_0 + \sum_{i=1}^n w_i V_t^{n,i}$, where $(V_t^{n,1}, \dots, V_t^{n,n})$ is a solution to the *n*-dimensional SDE [Alfonsi and Kebaier, 2024]:

$$V_{t}^{n,i} = -\int_{0}^{t} y_{i} V_{s}^{n,i} \, \mathrm{d}s + \int_{0}^{t} b(V_{0} + \sum_{i=1}^{n} w_{i} V_{s}^{n,i}) \, \mathrm{d}s + \int_{0}^{t} \sigma(V_{0} + \sum_{i=1}^{n} w_{i} V_{s}^{n,i}) \, \mathrm{d}W_{s}.$$
(8)

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\mathbb{L}^2 -error of Markovian Approximation

Proposition ([Alfonsi and Kebaier, 2024])

For every T > 0, there exists some C > 0 (depending on T, V_0 , b, σ), such that

$$\mathbb{E}[|V_{T} - V_{T}^{n}|^{2}] \leq C \int_{0}^{T} |K(t) - K^{n}(t)|^{2} dt.$$
(9)

 The choice of weights (w_i)ⁿ_{i=1} and nodes (y_i)ⁿ_{i=1} plays a crucial role in reducing the kernel approximation error ∫₀^T |K(t) − Kⁿ(t)|²dt.

Trade-off Accuracy and Efficiency

- Increasing *n* improves approximation accuracy by providing more weights and nodes for a better integral approximation.
- However, a larger *n* increases the dimensionality of the SDE, leading to higher computational cost and greater discretization error.

Achieving good accuracy with small n is crucial

Selection of Weights and Nodes

Deterministic Approaches:

- Use quadrature rules to approximate the kernel.
- Use same nodes and weights for all sample paths.
- Consequently, the bias introduced by the approximation is the same for every path.
- This effect is especially pronounced when *n* is small.

Introducing stochastic selection of nodes

Construction of Randomized Approximation Kernel

- Let Y be a random variable with density f_Y supported on $(0, \infty)$.
- The kernel K(t) can be expressed as an **expectation**:

$$\begin{split} \mathcal{K}(t) &= \int_{0}^{\infty} c_{H} y^{-H - \frac{1}{2}} e^{-yt} \, \mathrm{d}y = \int_{0}^{\infty} c_{H} y^{-H - \frac{1}{2}} e^{-yt} \frac{1}{f_{Y}(y)} f_{Y}(y) \, \mathrm{d}y \\ &= \mathbb{E}[\frac{c_{H} Y^{-H - \frac{1}{2}}}{f_{Y}(Y)} e^{-Yt}]. \end{split}$$

(10)

• Let $Z_Y(t) := \frac{c_H Y^{-H-\frac{1}{2}}}{f_Y(Y)} e^{-Yt}$. By the law of large numbers, approximate this expectation by $K^{n,Y}(t)$ as a finite sum:

$$\mathcal{K}^{n,Y}(t) := \frac{1}{n} \sum_{i=1}^{n} Z_{Y_i}(t), \qquad (11)$$

where $Y_i \sim Y$ are i.i.d.

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Randomized Markovian Approximation

A randomized Markovian approximation is defined as:

$$V_{t}^{n,Y} = V_{0} + \int_{0}^{t} \frac{K^{n,Y}(t-s)b(V_{t}^{n,Y}) dt}{t} + \int_{0}^{t} \frac{K^{n,Y}(t-s)\sigma(V_{t}^{n,Y}) dW_{t}}{(12)}$$

• As in the non-randomized case:

- V_t^{n,Y} can be expressed using the solution of an *n*-dimensional SDE.
- The \mathbb{L}^2 -error is controlled by the kernel approximation error:

$$\mathbb{E}[|V_{T} - V_{T}^{n,Y}|^{2}] \leq C\mathbb{E}\left[\int_{0}^{T} |K(t) - K^{n,Y}(t)|^{2} dt\right]$$

= $\frac{C}{n} \int_{0}^{T} \operatorname{Var}[Z_{Y}(t)] dt.$ (13)

Selection of density f_Y

• To minimize $\int_0^T \operatorname{Var}[Z_Y(t)] dt$, the **optimal density** is given by:

$$f_Y(y) = \frac{y^{-H-1}\sqrt{1 - e^{-2yT}}}{\int_0^\infty y^{-H-1}\sqrt{1 - e^{-2yT}} \,\mathrm{d}y}.$$
 (14)

- However, this optimal density is impractical for random number generation.
- Instead, use a **mimic density** f_Y as an alternative:

$$f_{Y}(y) = \begin{cases} C_{1}y^{-\alpha}, & 0 < y \le G, \\ C_{2}y^{-\beta}, & G < y. \end{cases}$$
(15)

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Variance Reduction Techniques

Objective: Reduce $\int_0^T \operatorname{Var}[Z_Y(t)] dt$ to minimize the kernel approximation error.

- Control Variates (CV):
 - Use an auxiliary variable with a known expectation to adjust the kernel approximation.
 - Minimize variance by introducing an optimal correction term.
- Antithetic Variates (AV):
 - In the case where n = 2, construct antithetic pairs (Y_1, Y_2) .
 - Leverage negative correlation to reduce variance.

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Effect of Large Node in Markovian Approximation

• In the Markovian approximation, the SDE (8) includes the mean-reverting term:

$$-\int_0^t y_i V_s^{n,i} \,\mathrm{d}s. \tag{16}$$

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- Large y_i leads to **rapid decay**, increasing discretization error.
- This issue arises in both **quadrature-based** and **randomized** approximation:
 - Quadrature: Large nodes may be inevitable due to the integration rule.
 - ► Randomized: Since Y has support on (0,∞), large nodes can occasionally appear.

Strategy to Mitigate Large Node Truncation of PDF:

For L > 0, let $c_L := \mathbb{P}(Y \le L) = \int_0^L f_Y(y) \, \mathrm{d}y$.

• Define a density f_{Y^L} as:

$$f_{Y^L}(y) := \begin{cases} \frac{f_Y(y)}{c_L}, & y < L\\ 0, & L \le y. \end{cases}$$
(17)

• The random variable $Y^L \sim f_{Y^L}$ takes values in (0,L). Kernel Approximation Error:

$$\mathbb{E}\Big[\int_0^T |K(t) - K^{n,Y^L}(t)|^2 \,\mathrm{d}t\Big] = \frac{1}{n} \int_0^T \operatorname{Var}[Z_{Y^L}(t)] \,\mathrm{d}t + \int_0^T (\int_L^\infty c_H y^{-H - \frac{1}{2}} e^{-yt} \,\mathrm{d}y)^2 \,\mathrm{d}t \text{ (truncation error).}$$
(18)

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Numerical Result

- Model: Rough Heston model
- Option Type: European Call Option
- Parameters:

$$H = 0.1, \quad \lambda = 0.3, \quad \nu = 0.3, \quad \theta = 0.02, \\ \rho = -0.7, \quad V_0 = 0.02, \quad S_0 = 1, \quad T = 1.$$
(19)

- Monte Carlo Paths: 10⁷
- **Discretization:** Multi-Factor Euler Scheme [Alfonsi and Kebaier, 2024]
- Time Steps: 512
- Number of Nodes: *n* = 2
- Benchmark: Fourier-based Option Price

Maximum Relative Error (MRE)

$$\mathsf{MRE} = \max_{k \in [-0.2, 0.2]} \frac{|\mathsf{IV}_{\mathsf{analytical}}(k) - \mathsf{IV}_{\mathsf{approximation}}(k)|}{|\mathsf{V}_{\mathsf{analytical}}(k)} \times 100 \quad (20)$$

- IV_{analytical}(k): IV obtained using the Fourier-based method
- IV_{approximation}(k): IV obtained using the Markovian approximation and Monte Carlo simulation
- k: Log-moneyness
- The error is computed over the log-moneyness range [-0.2, 0.2], using 17 equally spaced points.

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MRE of IV for different values of L



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IV: Deterministic vs Randomized



AE ([Abi Jaber and El Euch, 2019]), AK ([Alfonsi and Kebaier, 2024]), GG ([Bayer and Breneis, 2023a]), BL2 ([Bayer and Breneis, 2023b, Appendix F]), and RMA (Randomized Markovian and the second seco

MRE: Deterministic vs Randomized

- RMA outperforms AE, AK, and GG.
- BL2 has the lowest error but relies on numerical optimization with weak theoretical justification.

Improvement

- The current approximation kernel only approximates the integral representation of K(t) over [0, L].
- As an improvement, the approximate kernel is scaled by a constant *C* to reduce the approximation error.
- Specifically, we seek to minimize the error

$$E\int_0^T \left| K(t) - C K^{n,Y}(t) \right|^2 \mathrm{d}t.$$
 (21)

• The optimal constant C is given by

$$C = \frac{E \int_0^T \mathcal{K}(t) \, \mathcal{K}^{n,Y}(t) \, dt}{E \int_0^T \left(\mathcal{K}^{n,Y}(t) \right)^2 \, \mathrm{d}t}.$$
 (22)

Improvement Result



MRE for logmoneyness in [-0.2, 0.2].



MRE for logmoneyness in [-0.1, 0.1].

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Summary

- Introduced the probabilistic approach to determining weights and nodes in the Markovian approximation.
- Provided the optimal density and the mimic density to minimize the kernel approximation error.
- Demonstrated the effectiveness of variance reduction techniques.
- Showed that the randomized Markovian approximation outperforms all quadrature methods except BL2.

Future Work

• Analyze the impact of specific nodes and weights on approximation accuracy and develop alternative strategies to \mathbb{L}^2 minimization.

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