1 Problem setting

1.1 The rBergomi model

We consider the rBergomi model for the price process S_t as defined in [1], normalized to $r = 0^1$, which is defined by

(1.1)
$$dS_t = \sqrt{v_t} S_t dZ_t,$$
$$v_t = \xi_0(t) \exp\left(\eta \widetilde{W}_t^H - \frac{1}{2} \eta^2 t^{2H}\right),$$

where the Hurst parameter 0 < H < 1 and $\eta > 0$. We refer to v_t as the variance process, and $\xi_0(t) = \mathrm{E}\left[v_t\right]$ is the forward variance curve. Here, \widetilde{W}^H is a certain Riemann-Liouville fBm process², defined by

$$\widetilde{W}_t^H = \int_0^t K^H(t, s) dW_s^1, \quad t \ge 0,$$

where the kernel $K^H: \mathbb{R}_+ \to \mathbb{R}_+$ is

(1.3)
$$K^{H}(t-s) = \sqrt{2H}(t-s)^{H-1/2}, \quad \forall \ 0 \le s \le t.$$

By construction, \widetilde{W}^H is a centered, locally $(H - \epsilon)$ - Hölder continuous, Gaussian process with $\operatorname{Var}\left[\widetilde{W}_t^H\right] = t^{2H}$, and a dependence structure defined by

$$\mathrm{E}\left[\widetilde{W}_{u}^{H}\widetilde{W}_{v}^{H}\right] = u^{2H}G\left(\frac{v}{u}\right), \quad v > u,$$

where for $x \ge 1$ and $\gamma = \frac{1}{2} - H$

(1.4)
$$G(x) = 2H \int_0^1 \frac{ds}{(1-s)^{\gamma}(x-s)^{\gamma}}.$$

In (1.1) and (1.2), W^1, Z denote two *correlated* standard Brownian motions with correlation $\rho \in]-1,0]$, so that we can represent Z in terms of W^1 as

$$Z = \rho W^1 + \overline{\rho} W^{\perp} = \rho W^1 + \sqrt{1 - \rho^2} W^{\perp},$$

where (W^1, W^{\perp}) are two independent standard Brownian motions. Therefore, the solution to (1.1), with $S(0) = S_0$, can be written as

$$S_t = S_0 \exp\left(\int_0^t \sqrt{v(s)} dZ(s) - \frac{1}{2} \int_0^t v(s) ds\right), \quad S_0 > 0$$

$$v_u = \xi_0(u) \exp\left(\eta \widetilde{W}_u^H - \frac{\eta^2}{2} u^{2H}\right), \quad \xi_0 > 0.$$
(1.5)

 $^{^{1}}r$ is the interest rate.

²The so-called Riemann-Liouville processes are deduced from the standard Brownian motion by applying Riemann-Liouville fractional operators, whereas the standard fBm requires a weighted fractional operator [?, ?].

The filtration $(\mathcal{F}_t)_{t\geq 0}$ can here be taken as the one generated by the two-dimensional Brownian motion (W^1,W^{\perp}) under the risk neutral measure \mathbb{Q} , resulting in a filtered probability space $(\Omega,\mathcal{F},\mathcal{F}_t,\mathbb{Q})$. The stock price process S is clearly then a local $(\mathcal{F}_t)_{t\geq 0}$ -martingale and a supermartingale. We shall henceforth use the notation $\mathrm{E}\left[.\right] = E^{\mathbb{Q}}\left[.\mid \mathcal{F}_0\right]$ unless we state otherwise.

Remark 1.1. The rBergomi model is non-Markovian in the instantaneous variance v_t , that is $E^{\mathbb{Q}}[v_u \mid \mathcal{F}_t] \neq = E^{\mathbb{Q}}[v_u \mid v_t]$. However, it is Markovian in the state vector by definition, that is $E^{\mathbb{Q}}[v_u \mid \mathcal{F}_t] = \xi_t(u)$.

1.2 Option pricing under the rBergomi model

We are interested in pricing European call options under the rBergomi model. Assuming $S_0 = 1$, and using the conditioning argument on the σ -algebra generated by W^1 (an argument first used by [2] in the context of Markovian stochastic volatility models), we can show that the call price is given by

$$C_{RB}(T, K) = E \left[(S_T - K)^+ \right]$$

$$= E \left[E \left[(S_T - K)^+ \mid \sigma(W^1(t), t \le T) \right] \right]$$

$$= E \left[C_{BS} \left(S_0 = \exp \left(\rho \int_0^T \sqrt{v_t} dW_t^1 - \frac{1}{2} \rho^2 \int_0^T v_t dt \right), \ k = K, \ \sigma^2 = (1 - \rho^2) \int_0^T v_t dt \right) \right],$$

where $C_{BS}(S_0, k, \sigma^2)$ denotes the Black-Scholes call price, for initial spot price S_0 , strike price k and volatility σ^2 .

We point out that the analytical smoothing, based on conditioning, performed in (1.6) enables us to uncover the available regularity, and hence get a smooth, analytic integrand inside the expectation. Therefore, applying a deterministic quadrature technique such as ASGQ or QMC becomes an adequate option for computing the call price as we will investigate later.

2 Quantity of interest and error bounds

We recall that our goal is to compute the expectation in (1.6). We need 2N-dimensional Gaussian inputs for the used hybrid scheme (N) is the number of time steps in the time grid), namely

- $\mathbf{W}^{(1)} = \{W_i^{(1)}\}_{i=1}^N$: The N Gaussian random variables that are defined in Section 1.1.
- $\mathbf{W}^{(2)} = \{W_j^{(2)}\}_{j=1}^N$: An artificially introduced N Gaussian random variables that are used for left-rule points in the hybrid scheme.

We can rewrite (1.6) as

$$C_{RB}(T, K) = E\left[C_{BS}\left(S_{0} = \exp\left(\rho \int_{0}^{T} \sqrt{v_{t}} dW_{t}^{1} - \frac{1}{2}\rho^{2} \int_{0}^{T} v_{t} dt\right), \ k = K, \ \sigma^{2} = (1 - \rho^{2}) \int_{0}^{T} v_{t} dt\right)\right]$$

$$\approx \int_{\mathbb{R}^{2N}} C_{BS}\left(G(\mathbf{w}^{(1)}, \mathbf{w}^{(2)})\right) \rho_{N}(\mathbf{w}^{(1)}) \rho_{N}(\mathbf{w}^{(2)}) d\mathbf{w}^{(1)} d\mathbf{w}^{(2)}$$

$$:= C_{RB}^{N},$$
(2.1)

where G maps 2N independent standard Gaussian random inputs to the parameters fed to Black-Scholes formula, and ρ_N is the multivariate Gaussian density, given by

$$\rho_N(\mathbf{z}) = \frac{1}{(2\pi)^{N/2}} e^{-\frac{1}{2}\mathbf{z}^T\mathbf{z}}.$$

Therefore, the initial integration problem that we are solving lives in 2N-dimensional space, which becomes very large as the number of time steps N, used in the hybrid scheme, increases.

If we denote by \mathcal{E}_{tot} the total error of approximating the expectation in (1.6) using the ASGQ estimator, Q_N , then we have a natural error decomposition

(2.2)
$$\mathcal{E}_{\text{tot}} \le |C_{\text{RB}} - C_{\text{RB}}^N| + |C_{\text{RB}}^N - Q_N| \le \mathcal{E}_B(N) + \mathcal{E}_Q(\text{TOL}_{\text{ASGQ}}, N),$$

where \mathcal{E}_Q is the quadrature error, \mathcal{E}_B is the bias, TOL_{ASGQ} is a user selected tolerance for ASGQ method, and C_{RB}^N is the biased price computed with N time steps as given by (2.1).

On the other hand, the total error of approximating the expectation in (1.6) using the randomized QMC or MC estimator, $Q_N^{\text{MC(QMC)}}$ can be bounded by

(2.3)
$$\mathcal{E}_{\text{tot}} \le \left| C_{\text{RB}} - C_{\text{RB}}^{N} \right| + \left| C_{\text{RB}}^{N} - Q_{N}^{\text{MC (QMC)}} \right| \le \mathcal{E}_{B}(N) + \mathcal{E}_{S}(M, N),$$

where \mathcal{E}_S is the statistical error³, M is the number of samples used for MC or randomized QMC method.

3 ASGQ error estimate

As discussed, potential ways of estimating the quadrature error are presented below.

3.1 First way: Similar to the one implemented by Joakim

I think this way is almost similar to the one implemented by Joakim. He may correct me if I missed some things.

In our case, once we fix N, we define from (2.1)

$$F^N = C_{BS}(G(\mathbf{W}^{(1)}, \mathbf{W}^{(2)})).$$

We introduce the set $C^0(\mathbb{R})$ of real-valued continuous functions over \mathbb{R} , and the subspace of polynomials of degree at most q over \mathbb{R} , $\mathbb{P}^q(\mathbb{R}) \subset C^0(\mathbb{R})$. Next, we consider a sequence of univariate Lagrangian interpolant operators in each dimension Y_n $(1 \leq n \leq 2N)$, that is, $\{U_n^{m(\beta_n)}\}_{\beta_n \in \mathbb{N}_+}$ (we refer to the value β_n as the interpolation level). Each interpolant is built over a set of $m(\beta_n)$ collocation points, $\mathcal{H}^{m(\beta_n)} = \{y_n^1, y_n^2, \dots, y_n^{m(\beta_n)}\} \subset \mathbb{R}$, thus, the interpolant yields a polynomial approximation,

$$U^{m(\beta_n)}: C^0(\mathbb{R}) \to \mathbb{P}^{m(\beta_n)-1}(\mathbb{R}), \quad U^{m(\beta_n)}[F^N](y_n) = \sum_{j=1}^{m(\beta_n)} \left(f(y_n^j) \prod_{k=1; k \neq j}^{m(\beta_n)} \frac{y_n - y_n^k}{y_n^j - y_n^k} \right).$$

The statistical error estimate of MC or randomized QMC is $C_{\alpha} \frac{\sigma_M}{\sqrt{M}}$, where M is the number of samples and $C_{\alpha} = 1.96$ for 95% confidence interval.

The 2N-variate Lagrangian interpolant can then be built by a tensorization of univariate interpolants: denote by $C^0(\mathbb{R}^{2N})$ the space of real-valued 2N-variate continuous functions over \mathbb{R}^{2N} and by $\mathbb{P}^{\mathbf{q}}(\mathbb{R}^{2N}) = \bigotimes_{n=1}^{2N} \mathbb{P}^{\mathbf{q}_n}(\mathbb{R})$ the subspace of polynomials of degree at most q_n over \mathbb{R} , with $\mathbf{q} = (q_1, \ldots, q_{2N}) \in \mathbb{N}^{2N}$, and consider a multi-index $\boldsymbol{\beta} \in \mathbb{N}_+^{2N}$ assigning the interpolation level in each direction, y_n , then the multivariate interpolant can then be written as

$$U^{m(\boldsymbol{\beta})}: C^0(\mathbb{R}^{2N}) \to \mathbb{P}^{m(\boldsymbol{\beta})-1}(\mathbb{R}^{2N}), \quad U^{m(\boldsymbol{\beta})}[F^N](\mathbf{y}) = \bigotimes_{n=1}^{2N} U^{m(\beta_n)}[F^N](\mathbf{y}).$$

Given this construction, we can define the ASGQ interpolant for approximating F^N , using a set of multi indices $\mathcal{I} \in \mathbb{N}^{2N}$ as

(3.1)
$$I^{\mathcal{I}}[F^N] = \sum_{\beta \in \mathcal{I}} \Delta U_N^{\beta},$$

where

$$\Delta_i U_N^{\beta} = \begin{cases} U_N^{\beta} - U_N^{\beta'}, \text{ with } \beta' = \beta - e_i, \text{ if } \beta_i > 0 \\ U_N^{\beta}, & \text{otherwise,} \end{cases}$$

where e_i denotes the *i*th 2N-dimensional unit vector. Then, ΔU_N^{β} is defined as

$$\Delta U_N^{\beta} = \left(\prod_{i=1}^{2N} \Delta_i\right) U_N^{\beta}.$$

We define the interpolation error induced by ASGQ as

$$(3.2) e_N = F^N - I^{\mathcal{I}}[F^N].$$

One can have a bound on the interpolation error of ASGQ, e_N , by tensorizing one dimensional error estimates, and then simply integrate that bound to get the ASGQ error, $\mathcal{E}_Q(\text{TOL}_{\text{ASGQ}}, N)$, defined in (2.2). However, we think that this will not lead to a sharp error estimate for ASGQ. Another strategy for estimating the ASGQ error, is to estimate $\text{E}\left[e_N\right]$ using MC by sampling directly e_N (what I think Jaokim code is doing actually).

If we define $Y = F^N + (Q_N^{\mathcal{I}} - I^{\mathcal{I}}[F^N])$ (where $Q_N^{\mathcal{I}}$ is the ASGQ estimator), then we have

(3.3)
$$E[Y] = E[F^N]$$

$$Var[Y] = Var[e_N] < Var[\mathcal{A}_{MC}],$$

where \mathcal{A}_{MC} is the MC estimator for $E[F^N]$.

(3.3) shows that ASGQ can be seen as a control variate for MC estimator and consequently as a powerful variance reduction tool.

This way of estimating the quadrature error comes with the disadvantage of exciting the strong error which has a poor behavior in our context resulting maybe to having a non-sharp error estimate. In fact, by the central limit theorem, we expect that

$$|\mathbf{E}[e_N]| = \left| \int \underbrace{F^N - I^{\mathcal{I}}[F^N](y)}_{Y(y)} dy \right|$$

$$\approx \frac{C_{\alpha}}{\sqrt{M}} \sqrt{\mathrm{Var}(Y)}.$$
(3.4)

In our context of the rBergomi model we know that the strong error is of order H, that is we expect to have $Var(Y) = \mathcal{O}(h^H)$ (h is the mesh size and H is the Hurst parameter which is of order ≈ 0.1 . As a consequence, it may be that using this way will not provide a sharp enough error estimate for the quadrature error!

3.2 Second way

To avoid exciting the strong error when estimating the quadrature error and just act on the weak error, we can use a second way that is inspired of randomized QMC. In fact, we suggest to use a randomized version of ASGQ where the randomization involves randomized rotation and scaling for quadrature rules since we deal with unbounded domains and Hermite quadrature rule. Although this comes with the advantage of just acting on the weak error, it has the issue of reducing anisotropy which is a main feature for a good performance of ASGQ.

We can formulate this more in case the first way fail!

3.3 Third way

One can learn the error curve as a way to reduce the extra burden that comes from estimating the ASGQ error but this not yet formulated yet. I will try to formulate it if the two previous options fail!

4 Numerical experiments of estimating quadrature error

In this section, I show the obtained quadrature estimates for one of my examples for N = 2, 4, 8 where N is the number of time steps. Looking at Figure 4.1, I have some questions that are not clear to me:

- \mathbf{Q}_1 : I did not get why the quadrature error is kind of constant and then it decays so fast for very large number of collocation points. Is it normal to observe such behavior? Also, this behavior is less weird when taking N=8.
 - A_1 : This behavior as per Raul is not weird so It maybe that with few points we already rich an accuracy that remains constant ans as we have larger number of quadrature points we see this sharp decay, meaning the accuracy changes a lot for high number of quadrature points!
- \mathbf{Q}_2 : Why the interpolation error seems to be constant?
 - A_2 : This maybe due to the fact that the statistical error is already dominating the bias as explained by (3.4)! so we do not see what really happens for the bias term
- Does it seem that using the interpolation error as a bound for the quadrature error is not that good?

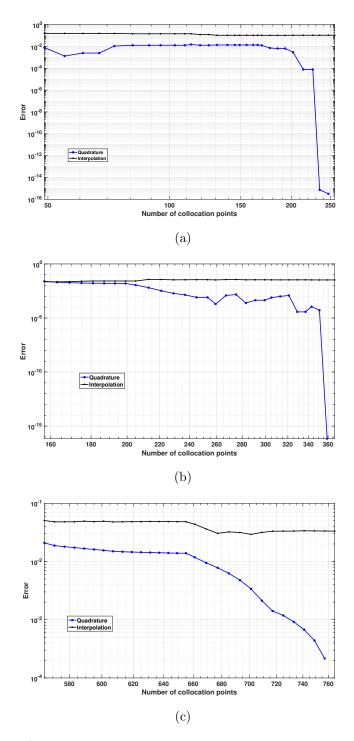


Figure 4.1: Quadrature (computed as the difference between the quadrature solution and a reference solution computed with very high number of quadrature points), and interpolation error (as explained in the first way of Section 3) for 3 different cases of number of time steps: a) N=2, b) N=4, c) N=8

References Cited

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- [2] Marc Romano and Nizar Touzi. Contingent claims and market completeness in a stochastic volatility model. *Mathematical Finance*, 7(4):399–412, 1997.