

Pricing in affine forward variance models

Jim Gatheral



Seminario di finanza matematica, probabilità
Bologna, October 13, 2021

Outline of this talk

- Affine forward variance (AFV) models
 - The classical Heston model
 - The rough Heston model
- The CGF and its associated Volterra integral equation
 - Numerical solution using the Adams scheme
 - Rational approximation of the rough Heston solution
- Simulation of AFV models
 - The RSQE scheme
 - The HQE scheme
 - Convergence

Computing option prices from the characteristic function

- It is quite straightforward to get option prices by inverting the characteristic function of a given stochastic process (if it is known in closed-form).
- For example, formula (5.6) of [Gat06] is a special case of formula (2.10) of [Lew00]:

Lemma 1 (Formula (2.10) of Lewis)

$$C(S, K, t, T) = S - \sqrt{SK} \frac{1}{\pi} \int_0^{\infty} \frac{du}{u^2 + \frac{1}{4}} \operatorname{Re} \left[e^{-iuk} \varphi_t(T; u - i/2) \right]$$

with $k = \log \left(\frac{K}{S} \right)$.

- An analogous formula holds for puts.

Alternative methods

Alternative methods include:

- The COS method of [FO09]
- The SINC method of [BBRR21].
- Both of these methods, in particular the SINC method, are significantly faster and more accurate than the Lewis formula when calculating the prices of many options with the same expiration.
 - *i.e.* when computing smiles.

Summary so far

- Given the characteristic function, it is straightforward to price European options.
- In AFV models, the characteristic function is given by

$$\mathbb{E}_t \left[e^{iaX_T} \right] =: \varphi_t(T; a) = \exp \left\{ \int_t^T \xi_t(s) g(T - s; ia) ds \right\}$$

where $g(\cdot; u)$ is the unique global continuous solution of the convolution Riccati equation (2):

$$g(\tau; u) = R_V \left(u, (\kappa \star g)(\tau, u) \right),$$

and $R_V(u, w) = \frac{1}{2}(u^2 - u) + \rho u w + \frac{1}{2} w^2$.

- All we therefore need to do is to solve (2).

The fractional Adams scheme

The fractional Adams scheme of [DFF04] is for the numerical approximation of the solution of equations of the form

$$h(\tau) = \frac{1}{\Gamma(\alpha)} \int_0^\tau (\tau - s)^{\alpha-1} F(h(s)) ds. \quad (4)$$

- The rough Heston convolution Riccati equation is of this type.

This leads to the following scheme:

$$\hat{h}(t_{k+1}) = \sum_{0 \leq j \leq k} a_{j,k+1} F(\hat{h}(t_j)) + a_{k+1,k+1} F(\hat{h}(t_{k+1})), \quad (5)$$

with

$$\begin{aligned} a_{0,k+1} &= \frac{\Delta^\alpha}{\Gamma(\alpha+2)} [k^{\alpha+1} - (k-\alpha)(k+1)^\alpha] \\ a_{j,k+1} &= \frac{\Delta^\alpha}{\Gamma(\alpha+2)} [(k-j+2)^{\alpha+1} + (k-j)^{\alpha+1} - 2(k-j+1)^{\alpha+1}]; \\ &\quad 1 \leq j \leq k \\ a_{k+1,k+1} &= \frac{(\Delta t)^\alpha}{\Gamma(\alpha+2)}. \end{aligned} \quad (6)$$

- $\hat{h}(t_{k+1})$ is on both sides of (5) so this scheme is implicit.
- Thus, we first compute a predictor of $\hat{h}^P(t_{k+1})$ of $\hat{h}(t_{k+1})$ and plug that predictor back into (5).
- We construct the predictor $\hat{h}^P(t_{k+1})$ by freezing \hat{g} at the beginning of each interval:

$$\hat{h}^P(t_{k+1}) = \frac{1}{\Gamma(\alpha)} \int_0^{t_{k+1}} (t_{k+1} - s)^{\alpha-1} \tilde{g}(s) ds,$$

with

$$\tilde{g}(t) = \hat{g}(t_j); \quad t \in [t_j, t_{j+1}), \quad 0 \leq j \leq k.$$

Therefore,

$$\hat{h}^P(t_{k+1}) = \sum_{0 \leq j \leq k} b_{j,k+1} F(\hat{h}(t_j)),$$

where

$$b_{j,k+1} = \frac{\Delta^\alpha}{\Gamma(\alpha + 1)} [(k - j + 1)^\alpha - (k - j)^\alpha], \quad 0 \leq j \leq k.$$

Thus, the final explicit numerical scheme is given by

$$\hat{h}(t_{k+1}) = \sum_{0 \leq j \leq k} a_{j,k+1} F(\hat{h}(t_j)) + a_{k+1,k+1} F(\hat{h}^P(t_j)),$$

where the weights $a_{j,k+1}$ are defined in (6).

Rational approximation of the Heston solution

- The Adams scheme presented above for solving the rough Heston fractional differential equation is slow!
- In [GR19], we showed how to approximate the solution of the rough Heston fractional Riccati equation by a rational function.
 - This approximation solution is just as fast as the classical Heston solution and appears to be more accurate than the Adams scheme for any reasonable number of time steps!

Rational approximation to the rough Heston solution

Wlog, set $\nu = 1$ and $x = t$. Then the rough Heston fractional Riccati ODE (3) reads

$$\begin{aligned} D^\alpha h(x; a) &= -\frac{1}{2} a(a+i) + i\rho a h(x; a) + \frac{1}{2} h(x; a)^2 \\ &= \frac{1}{2} (h(x; a) - r_-) (h(x; a) - r_+) \end{aligned}$$

with

$$A = \sqrt{a(a+i) - \rho^2 a^2}; \quad r_\pm = \{-i\rho a \pm A\}.$$

The idea is to paste together short- and long-time expansions of the solution using a rational (Padé) approximation.

Short-time expansion

From (for example) the exponentiation theorem of [AGR2020], $h(x; a)$ can be written as

$$h(x; a) = \sum_{j=0}^{\infty} \frac{\Gamma(1 + j\alpha)}{\Gamma(1 + (j+1)\alpha)} \beta_j(a) x^{(j+1)\alpha}$$

with

$$\beta_0(a) = -\frac{1}{2} a(a + i)$$

$$\beta_k(a) = \frac{1}{2} \sum_{i,j=0}^{k-2} \mathbb{1}_{\{i+j=k-2\}} \beta_i(a) \beta_j(a) \frac{\Gamma(1 + i\alpha)}{\Gamma(1 + (i+1)\alpha)} \frac{\Gamma(1 + j\alpha)}{\Gamma(1 + (j+1)\alpha)} + i \rho a \frac{\Gamma(1 + (k-1)\alpha)}{\Gamma(1 + k\alpha)} \beta_{k-1}(a).$$

Solving the rough Heston Riccati equation for long times

- In analogy with the classical Heston solution, we expect that for a suitable range of a ,

$$\lim_{x \rightarrow \infty} h(x; a) = r_-.$$

- In that case, for large x , we could linearize the fractional Riccati equation as follows.

$$\begin{aligned} D^\alpha h(x; a) &= \frac{1}{2} (h(x; a) - r_-) (h(x; a) - r_+) \\ &\approx -\frac{1}{2} (r_+ - r_-) (h(x; a) - r_-) \\ &= -A (h(x; a) - r_-). \end{aligned}$$

continued...

- The above linear fractional differential equation has the exact solution

$$h_{\infty}(a, x) = r_- [1 - E_{\alpha}(-Ax^{\alpha})],$$

where $E_{\alpha}(\cdot)$ is the Mittag-Leffler function.

- As $x \rightarrow \infty$,

$$E_{\alpha}(-Ax^{\alpha}) = -\frac{1}{A} \frac{x^{-\alpha}}{\Gamma(1-\alpha)} + \mathcal{O}(|Ax^{\alpha}|^{-2}).$$

- Thus, as $x \rightarrow \infty$,

$$h_{\infty}(a, x) - r_- = \frac{r_-}{A} \frac{x^{-\alpha}}{\Gamma(1-\alpha)} + \mathcal{O}(|Ax^{\alpha}|^{-2}).$$

Large x expansion

- The form of the asymptotic solution motivates the following expansion of h for large x :

$$h(x; a) = r_- \sum_{k=0}^{\infty} \gamma_k \frac{x^{-k\alpha}}{A^k \Gamma(1 - k\alpha)}.$$

- The coefficients γ_k satisfy the recursion

$$\gamma_1 = -\gamma_0 = -1$$

$$\gamma_k = -\gamma_{k-1} + \frac{r_-}{2A} \sum_{i,j=1}^{\infty} \mathbb{1}_{\{i+j=k\}} \gamma_i \gamma_j \frac{\Gamma(1 - k\alpha)}{\Gamma(1 - i\alpha) \Gamma(1 - j\alpha)}.$$

Rational approximation

- Now we have small- and large- x expansions we can compute global rational approximations to $h(x; a)$ of the form

$$h^{(m,n)}(x; a) = \frac{\sum_{i=1}^m p_i y^i}{\sum_{j=0}^n q_j y^j}$$

with $y = x^\alpha$ that match these expansions up to order m and n respectively.

- Only the diagonal approximants $h^{(n,n)}$ are admissible approximations of h .

$h^{(3,3)}$ is the best

- From various numerical experiments, the particular approximation $h^{(3,3)}$ seems to be amazingly close to the true solution for reasonable choices of model parameters.
- Though the excellent quality of the global approximation $h^{(3,3)}$ might at first seem very surprising, it is consistent with many Padé approximation stories from the literature.
- In our case, $h^{(3,3)}$ is clearly better than either $h^{(2,2)}$ or $h^{(4,4)}$.
 - $h^{(5,5)}$ is another very good approximation, but still not as good as $h^{(3,3)}$. $h^{(5,5)}$ is obviously also slower to compute.
 - Higher order approximations may turn out to beat $h^{(3,3)}$. However, $h^{(3,3)}$ may still be best in practice if speed of computation is taken into account.

Computing $h^{(3,3)}$

- We have the series expansion of h for small y :

$$h_s(y) = b_1 y + b_2 y^2 + b_3 y^3 + \mathcal{O}(y^4).$$

- We have the series expansion of h for large y :

$$h_\ell(y) = g_0 + \frac{g_1}{y} + \frac{g_2}{y^2} + \mathcal{O}\left(\frac{1}{y^3}\right).$$

- Matching the coefficients of the rational approximation

$$h^{(3,3)}(y) = \frac{p_1 y + p_2 y^2 + p_3 y^3}{1 + q_1 y + q_2 y^2 + q_3 y^3}$$

to $h_s(y)$ and $h_\ell(y)$ respectively gives a linear system of six equations for the six unknowns $\{p_i, q_i\}$.

Model parameters

- Recall that in our formulation of the rough Heston model,

$$\kappa(\tau) = \frac{\nu}{\Gamma(\alpha)} \tau^{\alpha-1},$$

with $\alpha = H + 1/2$.

- We choose model parameters roughly consistent with those found from calibration to SPX options on May 19, 2017 in [EGR19]:

$$\xi(u) = 0.025; H = 0.05; \nu = 0.4; \rho = -0.65. \quad (7)$$

Padé vs Adams smile errors

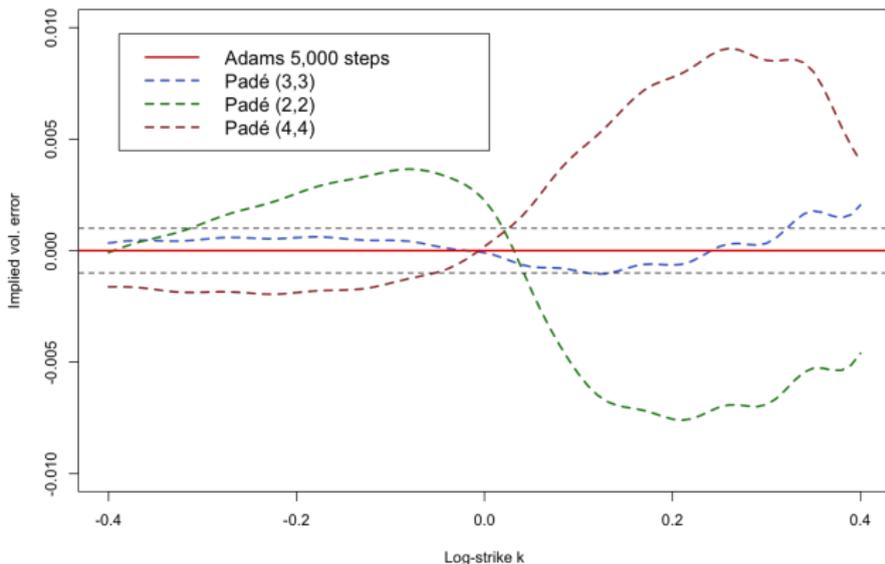


Figure 2: The red curve is the Adams smile with 5,000 steps. Again, $h^{(3,3)}$ errors are uncannily small.

Simulation

- We now turn our attention to simulation of AFV models as prescribed in [Gat21].
- Inspired by [BLP17] and [And08] we propose two algorithms to simulate AFV models.
 - Now that we have two alternative methods of computing the volatility smile under rough Heston, we can easily check convergence of these algorithms.

Discretization of the spot and variance processes

From the AFV dynamics,

$$d\xi_t(u) = \kappa(u - t) \sqrt{v_t} dW_t,$$

it follows that

$$\begin{aligned} v_T = \xi_T(T) &= \xi_0(T) + \int_0^T d\xi_s(T) \\ &= \xi_0(T) + \int_0^T \kappa(T - s) \sqrt{v_s} dW_s. \end{aligned} \quad (8)$$

- Wlog, let $t = 0$ and $\xi(u) = \xi_0(u)$. Let the time step $\Delta = T/N$ where N is the number of steps.
- As in [BLP17], we have the following exact decomposition of (8):

$$v_{n\Delta} = \xi(n\Delta) + \sum_{k=1}^n \int_{(k-1)\Delta}^{k\Delta} \kappa(n\Delta - s) \sqrt{v_s} dW_s.$$

Discretization of the v -process

- With simpler notation,

$$v_n = \xi_n + \sum_{k=1}^n \int_{(k-1)\Delta}^{k\Delta} \kappa(n\Delta - s) \sqrt{v_s} dW_s =: \hat{\xi}_n + u_n, \quad (9)$$

where the \mathcal{F}_{n-1} -adapted variable $\hat{\xi}_n$ is given by

$$\hat{\xi}_n = \mathbb{E}[v_n | \mathcal{F}_{n-1}] = \xi_n + \sum_{k=1}^{n-1} \int_{(k-1)\Delta}^{k\Delta} \kappa(n\Delta - s) \sqrt{v_s} dW_s, \quad (10)$$

and the martingale increment u_n by

$$u_n = \int_{(n-1)\Delta}^{n\Delta} \kappa(n\Delta - s) \sqrt{v_s} dW_s. \quad (11)$$

The X -process

- We also need to simulate the n th increment of the component of the log-stock price process $X = \log S$ parallel to the volatility process¹,

$$\chi_n = \int_{(n-1)\Delta}^{n\Delta} \sqrt{v_s} dW_s. \quad (12)$$

- We then have the following discretization of the X process:

$$X_n = X_{n-1} - \frac{1}{4} (v_n + v_{n-1}) \Delta + \sqrt{1 - \rho^2} \sqrt{\bar{v}_n \Delta} Z_n^\perp + \rho \chi_n,$$

where Z_n^\perp is standard normal, independent of χ_n and u_n .

¹We write the increments as χ_n to emphasize that they should be approximately χ^2 distributed random variables.

Choice of kernel

- In our computations we will focus on the *power-law* kernel:

$$\kappa(\tau) = \sqrt{2H\eta} \tau^{\alpha-1} =: \tilde{\eta} \tau^{\alpha-1}. \quad (13)$$

- We also need

Definition

For $i, j \geq 0$,

$$\mathcal{K}_i(\Delta) = \int_0^\Delta \kappa(s + i\Delta) ds;$$

$$\mathcal{K}_{i,j}(\Delta) = \int_0^\Delta \kappa(s + i\Delta) \kappa(s + j\Delta) ds.$$

- The $\mathcal{K}_{i,j}(\Delta)$ with $i \neq j$ are not in general computable in closed-form but are easy to compute numerically.

Covariances and correlations

- It can be shown that

$$\text{var}[u_n | \mathcal{F}_{n-1}] = \bar{v}_n \mathcal{K}_{0,0}(\Delta) + \mathcal{O}(\Delta^{1+2H}), \quad (14)$$

where

$$\bar{v}_n := \frac{1}{2H+1} \left[\hat{\xi}_n + 2H v_{n-1} \right].$$

- Similarly

$$\begin{aligned} \text{var}[\tilde{\xi}_{n+1} | \mathcal{F}_{n-1}] &\approx \bar{v}_n \mathcal{K}_{1,1}(\Delta) \\ \text{var}[\chi_n | \mathcal{F}_{n-1}] &\approx \bar{v}_n \Delta \\ \text{cov}[u_n, \tilde{\xi}_{n+1} | \mathcal{F}_{n-1}] &\approx \bar{v}_n \mathcal{K}_{0,1}(\Delta) \\ \text{cov}[u_n, \chi_n | \mathcal{F}_{n-1}] &\approx \bar{v}_n \mathcal{K}_0(\Delta) \\ \text{cov}[\chi_n, \tilde{\xi}_{n+1} | \mathcal{F}_{n-1}] &\approx \bar{v}_n \mathcal{K}_1(\Delta). \end{aligned} \quad (15)$$

Given a suitable kernel, all of these may be easily computed.

The correlation matrix

- Because variances and covariances in an AFV model are linear in ξ , the correlation matrix takes the simple form.

$$R = \begin{pmatrix} 1 & \rho_{u\chi} & \rho_{u\xi} \\ \rho_{u\chi} & 1 & \rho_{\xi\chi} \\ \rho_{u\xi} & \rho_{\xi\chi} & 1 \end{pmatrix}. \quad (16)$$

where

$$\rho_{u\chi} = \frac{\mathcal{K}_0(\Delta)}{\sqrt{\Delta}\sqrt{\mathcal{K}_{0,0}(\Delta)}}$$

$$\rho_{u\xi} = \frac{\mathcal{K}_{0,1}(\Delta)}{\sqrt{\mathcal{K}_{0,0}(\Delta)}\sqrt{\mathcal{K}_{1,1}(\Delta)}}$$

$$\rho_{\xi\chi} = \frac{\mathcal{K}_1(\Delta)}{\sqrt{\Delta}\sqrt{\mathcal{K}_{1,1}(\Delta)}}$$

are all independent of n .

The power-law kernel

- In the case of the power-law kernel $\kappa(\tau) = \tilde{\eta} \tau^{\alpha-1}$, these correlations are functions of H only.
- Specifically

$$\rho_{u\chi} = \frac{\sqrt{2H}}{H + 1/2},$$

and the other correlations may be easily computed numerically.

- In Figure 3, we plot these correlations as a function of H .

Plot of the correlation matrix in the power-law kernel case

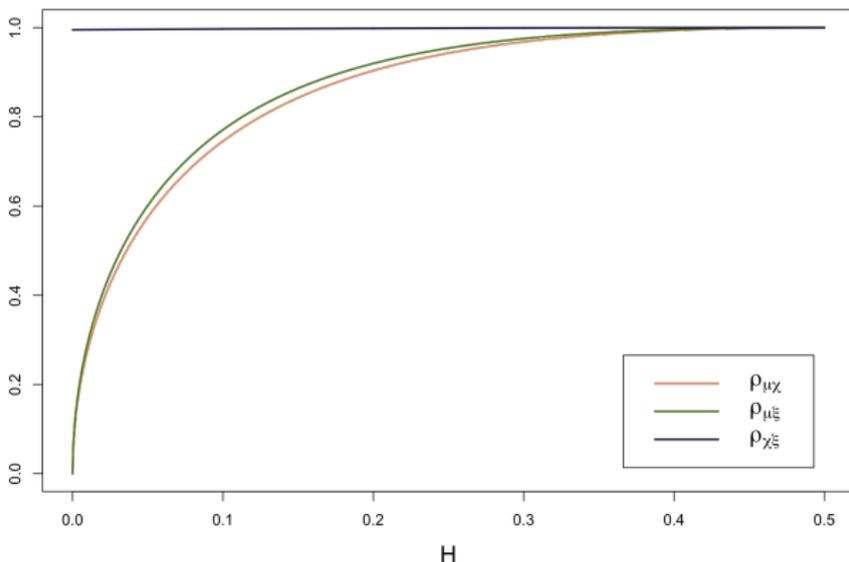


Figure 3: The correlations $\rho_{u\chi}$, $\rho_{u\xi}$, and $\rho_{\chi\xi}$ vs. H in the power-law kernel case.

A further approximation

- By assumption, the kernel behaves as a power-law kernel for Δ sufficiently small.
- Figure 3 thus suggests the following approximation whose motivation is easy to see by thinking of $\mathcal{K}_i(\Delta)$ as Δ times the average value of $\kappa(s + i\Delta)$ over the interval $(0, \Delta]$.

Approximate covariance

For $i \geq 0$ and $j \geq 1$,

$$\mathcal{K}_{i,j}(\Delta) \Delta \approx \mathcal{K}_i(\Delta) \mathcal{K}_j(\Delta). \quad (17)$$

An approximate correlation matrix

With the approximation (17),

$$\mathcal{K}_{0,1}(\Delta) \approx \frac{1}{\Delta} \mathcal{K}_1(\Delta) \mathcal{K}_0(\Delta); \quad \mathcal{K}_{1,1}(\Delta) \approx \frac{1}{\Delta} \mathcal{K}_1(\Delta)^2.$$

Substituting these expressions into (15) gives the correlation matrix

$$\bar{R} = \begin{pmatrix} 1 & \bar{\rho} & \bar{\rho} \\ \bar{\rho} & 1 & 1 \\ \bar{\rho} & 1 & 1 \end{pmatrix}, \quad (18)$$

where

$$\bar{\rho} \approx \frac{\mathcal{K}_0(\Delta)}{\sqrt{\mathcal{K}_{0,0}(\Delta)} \Delta} = \rho_{u\chi}.$$

Consequences for simulation

- At each step, we need to generate (at least) three random variables: u_n , χ_n , and $\hat{\xi}_{n+1}$.

$$u_n = \int_{(n-1)\Delta}^{n\Delta} \kappa(n\Delta - s) \sqrt{v_s} dW_s$$

$$\chi_n = \int_{(n-1)\Delta}^{n\Delta} \sqrt{v_s} dW_s$$

$$\hat{\xi}_{n+1} = \xi_{n+1} + \sum_{k=1}^n \int_{(k-1)\Delta}^{k\Delta} \kappa((n+1)\Delta - s) \sqrt{v_s} dW_s.$$

- When the model is Markovian ($H = 1/2$), we need only generate u_n at the n th time step; χ_n and $\hat{\xi}_{n+1}$ are perfectly correlated with u_n .
 - In practice, in the non-Markovian case ($H < 1/2$), we need only generate one other random variable consistent with the correlation matrix \bar{R} .

Average values of the kernel

- Echoing the notation of [BLP17], let

$$b_j^{*2} = \frac{1}{\Delta} \mathcal{K}_{j-1,j-1}(\Delta). \quad (19)$$

- b_j^{*2} thus gives the RMS average of the kernel at the j th lag.

The evolution of the forward variance curve

- The approximation

$$\int_{(k-1)\Delta}^{k\Delta} \kappa((n+1)\Delta - s) \sqrt{v_s} dW_s \approx b_{n+1-k}^* \chi_k$$

gives

$$\hat{\xi}_{n+1} \approx \xi_{n+1} + \sum_{k=1}^n b_{n+1-k}^* \chi_k.$$

- Similarly (though not needed for the algorithm), for $m > n$,

$$\mathbb{E}[v_m | \mathcal{F}_n] \approx \xi_m + \sum_{k=1}^n b_{m-k}^* \chi_k.$$

- We see that the entire forward variance curve evolves according to the weighted historical path of the $X = \log S$ process.

The Andersen Quadratic Exponential (QE) scheme

- Naïve simulation of the v process leads to negative values
- Andersen's Quadratic Exponential (QE) scheme [And08] guarantees v positive
 - Conditional means and variances are matched at each step
- We simulate according to the value of the ratio

$$\psi_n := \frac{\text{var}[v_n | \mathcal{F}_{n-1}]}{\hat{\xi}_n^2} = \frac{\text{var}[u_n | \mathcal{F}_{n-1}]}{\hat{\xi}_n^2}. \quad (20)$$

The Andersen Quadratic Exponential (QE) scheme

- If $\psi_n \leq 2$, simulate v_n as

$$v_n = \alpha_n (\beta_n + Z_n)^2$$

with $Z_n \sim N(0, 1)$ and

$$\beta_n^2 = \frac{2}{\psi_n} - 1 + \sqrt{\frac{2}{\psi_n} - 1} \sqrt{\frac{2}{\psi_n} - 1}; \quad \alpha_n = \frac{\hat{\xi}_n}{1 + \beta_n^2}.$$

- On the other hand, if $\psi_n \geq 1$, simulate v_n as

$$v_n = -\mathbb{1}_{\{U_n < p_n\}} \gamma_n \log \frac{U_n}{p_n}$$

with $U_n \sim \mathcal{U}(0, 1)$ and

$$p_n = \frac{2}{1 + \psi_n}; \quad \gamma_n = \frac{1}{2} \hat{\xi}_n (1 + \psi_n).$$

Conditional means and variances

- It is straightforward to check that means and variances are correctly matched in both cases.
- Since the two regions of applicability overlap, Andersen suggests to use algorithm ψ^- if $\psi_n < 3/2$ and algorithm ψ^+ if $\psi_n \geq 3/2$.

A Riemann-sum QE scheme

- Inspired by the Riemann-sum scheme of [BLP17] and the rough-Donsker scheme of [HJM17], we simulate the u_n , $\hat{\xi}_{n+1}$ and χ_n as if all three were perfectly correlated, equivalent to setting $\bar{\rho} = 1$ in (18).
 - From Figure 3 such an approximation may be justified if H is not too much less than $\frac{1}{2}$.

The RSQE scheme

The RSQE scheme

- 1 Given χ_k , for $k < n$, with ϵ very small, compute

$$\hat{\xi}_n = \max \left[\epsilon, \xi_n + \sum_{k=1}^{n-1} b_{n-k+1}^* \chi_k \right].$$

- 2 With $\text{var}[v_n | \mathcal{F}_{n-1}] = b_1^{*2} \bar{v}_n \Delta$, simulate v_n using the QE scheme.

3 $u_n = v_n - \hat{\xi}_n.$

4 $\chi_n = \frac{u_n}{b_1^*}.$

- 5 Finally,

$$X_n = X_{n-1} - \frac{1}{4} (v_n + v_{n-1}) \Delta + \sqrt{1 - \rho^2} \sqrt{\bar{v}_n \Delta} Z_n^\perp + \rho \chi_n.$$

A hybrid QE scheme

- The RSQE scheme matches unconditional means and variances at each step but it does not match the covariance structure of the process.
- For example, consider the conditional covariance between u_n and χ_n which is given by

$$\text{cov}[u_n, \chi_n | \mathcal{F}_{n-1}] = \int_{(n-1)\Delta}^{n\Delta} \kappa(n\Delta - s) \mathbb{E}[v_s | \mathcal{F}_{n-1}] ds \approx \bar{v}_n \mathcal{K}_0(\Delta).$$

- The RSQE scheme sets $u_n = b_1^* \chi_n$ so that

$$\text{cov}[u_n, \chi_n | \mathcal{F}_{n-1}] \approx b_1^* \text{var}[\chi_n | \mathcal{F}_{n-1}] = \bar{v}_n \sqrt{\mathcal{K}_{0,0}(\Delta) \Delta},$$

which is equivalent to the approximation

$$\mathcal{K}_0(\Delta) \approx \sqrt{\mathcal{K}_{0,0}(\Delta) \Delta}. \quad (21)$$

- Approximation (21), though accurate for small Δ when the kernel κ has no singularity at zero, is obviously very inaccurate when H is small.
- The essence of the hybrid scheme with $\kappa = 1$ of [BLP17] is to correct the error in the approximation (21) by simulating another random variable, uncorrelated with u_n , so as to match the covariance of u_n and χ_n .
 - For this, we need a bivariate version of Andersen's QE scheme.

A bivariate version of Andersen's QE scheme

- As before, let

$$u_n = \int_{(n-1)\Delta}^{n\Delta} \kappa(n\Delta - s) \sqrt{v_s} dW_s; \quad \chi_n = \int_{(n-1)\Delta}^{n\Delta} \sqrt{v_s} dW_s.$$

- Linear regression gives

$$u_n \approx \beta_{u\chi} \chi_n + \varepsilon_n,$$

where $\beta_{u\chi} = \mathcal{K}_0(\Delta)/\Delta$, and ε_n and χ_n are uncorrelated.

- Since $v_n = \hat{\xi}_n + u_n \geq 0$, we must ensure that $\beta_{u\chi} \chi_n + \varepsilon_n + \hat{\xi}_n \geq 0$.
- A scheme to achieve this is given in the following lemma.

The hybrid QE (HQE) scheme

The HQE scheme

- 1 Given χ_k , for $k < n$, with ϵ very small, compute

$$\hat{\xi}_n = \max \left[\epsilon, \xi_n + \sum_{k=1}^{n-1} b_{n-k+1}^* \chi_k \right].$$

- 2 Simulate χ_n and ε_n using the bivariate QE scheme

- 3 $v_n = \hat{\xi}_n + \frac{1}{\Delta} \mathcal{K}_0(\Delta) \chi_n + \varepsilon_n.$

- 4 Finally,

$$X_n = X_{n-1} - \frac{1}{4} (v_n + v_{n-1}) \Delta + \sqrt{1 - \rho^2} \sqrt{\bar{v}_n} \Delta Z_n^\perp + \rho \chi_n.$$



Jim Gatheral.

The volatility surface: A practitioner's guide.

John Wiley & Sons, 2006.



Jim Gatheral and Martin Keller-Ressel.

Affine forward variance models.

Finance and Stochastics, 23(3):501–533, 2019.



Jim Gatheral and Radoš Radoičić.

Rational approximation of the rough Heston solution.

International Journal of Theoretical and Applied Finance, 22(3):1950010–19, 2019.



Jim Gatheral

Efficient simulation of affine forward variance models.

SSRN, <https://www.ssrn.com/abstract=3876680>, 2021.



Blanka Horvath, Antoine Jacquier, and Aitor Muguruza.

Functional central limit theorems for rough volatility.

SSRN, <https://www.ssrn.com/abstract=3078743>, 2017.



Alan Lewis,

Option Valuation under Stochastic Volatility with Mathematica Code.

Finance Press: Newport Beach, CA, 2000.