Interest Rate Models 8. LIBOR Market Model, II

Andrew Lesniewski

Baruch College New York

Spring 2017

Outline

- Simulating Brownian motion
- 2 Discretizing SDEs
- 3 Generating Monte Carlo paths for LMM

Monte Carlo methods for LMM

- LMM does not allow for a natural implementation based on recombining trees, and thus all valuations have to be performed via Monte Carlo simulations.
- We shall describe two numerical schemes for generating Monte Carlo paths for LMM: Euler's scheme and Milstein's scheme.
- They both rely on replacing continuous time stochastic differential equations by suitable finite difference schemes.

- There exist many more of less refined methods for simulating a Wiener process; here we describe two of them.
- The random walk method is easy to implement at the expense of being rather noisy. It represents a Wiener process as a random walk sampled at a finite set of event dates t₀ < t₁ < ... < t_m:

$$Z(t_{-1}) = 0,$$

$$Z(t_n) = Z(t_{n-1}) + \sqrt{t_n - t_{n-1}} \, \xi_n, \quad n = 0, \dots, m,$$
(1)

where $t_{-1} = 0$, and where ξ_n are i.i.d. random variables with $\xi_n \sim N(0, 1)$.

• A good method of generating the ξ_n 's is to first generate a sequence of uniform pseudorandom numbers u_n (using, say, the Mersenne twister algorithm), and then set

$$\xi_n = N^{-1}(u_n), \tag{2}$$

where $N^{-1}(x)$ is the inverse cumulative normal function. $N^{-1}(x)$ can be efficiently and accurately computed using e.g. the Beasley-Springer-Moro algorithm, see [1].



- The spectral decomposition method generally leads to much better performance than the random walk method. It assures that the simulated process has the same covariance matrix C as the Wiener process Z(t) sampled at t_0, t_1, \ldots, t_m .
- The covariance matrix is explicitly given by:

$$C_{ij} = E[Z(t_i)Z(t_j)]$$

$$= \min(t_i, t_j).$$
(3)

Consider the eigenvalue problem for C:

$$CE_j = \lambda_j E_j, \qquad j = 0, \ldots, m,$$
 (4)

with orthonormal E_i 's.

• Since the covariance matrix $\mathcal C$ is positive definite, all of its eigenvalues λ_j are nonnegative, and we will assume that

$$\lambda_0 \geq \ldots \geq \lambda_m \geq 0.$$
 (5)



• We will denote the *n*-th component of the vector E_j by $E_j(t_n)$, and consider the random variable

$$Z(t_n) = \sum_{0 \le j \le m} \sqrt{\lambda_j} \, E_j(t_n) \xi_j, \tag{6}$$

where ξ_i are, again, i.i.d. random variables with $\xi_i \sim N(0, 1)$.

- These numbers are best calculated by applying the inverse cumulative normal function to a sequence of Sobol numbers. Alternatively, one could use a sequence of uniform pseudorandom numbers; this, however, leads to a higher sampling variance.
- Then, for each $n = 0, ..., m, Z(t_n) \sim N(0, t_n)$, and

$$E[Z(t_i)Z(t_j)] = \sum_{0 \le k \le m} \lambda_k E_k(t_i)E_k(t_j)$$

$$= C_{ij}.$$
(7)

- We can thus regard $Z(t_n)$ a realization of the discretized Wiener process¹.
- For computational efficiency, we may want to truncate (6) at some p < m. This eliminates the *high frequencies* from $Z(t_n)$, and lowers the variance. The price for this may be systematically lower accuracy.

¹ This realization of the discretized Wiener process is related to the well known Karhounen-Loeve expansion of the (continuous time) Wiener process.

Multi factor Brownian motion

• We now consider the case of a multi-factor Brownian motion $Z_a(t)$, with

$$\mathsf{E}[dZ_a(t)\,dZ_b(t)] = \rho_{ab}dt.$$

• The Cholesky decomposition of ρ yields

$$\rho = LL^{\mathrm{T}},\tag{8}$$

where *L* is a $d \times d$ dimensional, lower triangular matrix.

For example, if

$$\rho = \begin{bmatrix} 1 & \rho_{12} \\ \rho_{12} & 1 \end{bmatrix},\tag{9}$$

then

$$L = \begin{bmatrix} 1 & 0 \\ \rho & \sqrt{1 - \rho_{12}^2} \end{bmatrix},\tag{10}$$

Multi factor Brownian motion

- Now, if $X \in \mathbb{R}^d$ is a vector of independent standard normal variables, then LX is a multivariate normal variable with correlation matrix ρ .
- Indeed,

$$\begin{aligned} \mathsf{E}[(LX)_a(LX)_b] &= \sum_{0 \le k, l \le d} L_{ak} L_{bl} \mathsf{E}[X_k X_l] \\ &= \sum_{0 \le k, l \le d} L_{ak} L_{bl} \delta_{kl} \\ &= \sum_{0 \le k \le d} L_{ak} L_{bk} \\ &= \sum_{0 \le k \le d} L_{ak} (L^{\mathsf{T}})_{kb} \\ &= (LL^{\mathsf{T}})_{kl} \\ &= \rho_{kl}. \end{aligned}$$

- Numerical solution of a stochastic differential equation amounts to generating paths of the state variables given a path of the stochastic drivers of the system, namely the underlying Brownian motion.
- This requires approximating the continuous time system by a discrete time stochastic system.
- Consider first a one factor SDE,

$$dX(t) = A(t, X(t))dt + B(t, X(t))dZ(t), X(0) = X_0.$$
(11)

This is equivalent to

$$X(s) = X(t) + \int_{t}^{s} A(u, X(u)) du + \int_{t}^{s} B(u, X(u)) dZ(u).$$
 (12)



lacktriangle Now, if f(t,x) is twice continuously differentiable, then Ito's lemma states

$$df(t,X(t)) = \mathcal{L}^{0}f(t,X(t))dt + \mathcal{L}^{1}f(t,X(t))dZ(t),$$
(13)

where the operators \mathcal{L}^i are defined by

$$\mathcal{L}^{0} = \frac{\partial}{\partial t} + A \frac{\partial}{\partial x} + \frac{1}{2} B^{2} \frac{\partial^{2}}{\partial x^{2}}, \qquad (14)$$

and

$$\mathcal{L}^1 = B \, \frac{\partial}{\partial x} \,. \tag{15}$$

Applying Ito's lemma (13) to A yields

$$A(s,X(s)) = A(t,X(t)) + \int_{t}^{s} \mathcal{L}^{0} A(u,X(u)) du + \int_{t}^{s} \mathcal{L}^{1} B(u,X(u)) dZ(u)$$

$$\approx A(t,X(t)) + \mathcal{L}^{0} A(t,X(t)) \int_{t}^{s} du + \mathcal{L}^{1} B(t,X(t)) \int_{t}^{s} dZ(u).$$

We can thus approximate

$$\int_{t}^{t+\delta} A(s,X(s)) ds \approx A(t,X(t)) \delta + \mathcal{L}^{0} A(t,X(t)) I_{(0,0)} + \mathcal{L}^{1} B(t,X(t)) I_{(1,0)}.$$

Here

$$I_{(0,0)} = \int_{t}^{t+\delta} \int_{t}^{s} du \, ds,$$

$$I_{(1,0)} = \int_{t}^{t+\delta} \int_{t}^{s} dZ(u) \, ds,$$
(16)

are iterated integrals.



Similarly, we make the following approximation:

$$\begin{split} B(s,X(s)) &= B(t,X(t)) + \int_t^s \mathcal{L}^0 B(u,X(u)) du + \int_t^s \mathcal{L}^1 B(u,X(u)) dZ(u) \\ &\approx B(t,X(t)) + \mathcal{L}^0 B(t,X(t)) \int_t^s du + \mathcal{L}^1 B(t,X(t)) \int_t^s dZ(u) \,. \end{split}$$

Therefore,

$$\int_{t}^{t+\delta}B(s,X(s))dZ(s)\approx B(t,X(t))\Delta Z(t)+\mathcal{L}^{0}B(t,X(t))I_{(0,1)}+\mathcal{L}^{1}B(t,X(t))I_{(1,1)},$$

where $\Delta Z(t) = Z(t + \delta) - Z(t)$, and

$$I_{(0,1)} = \int_{t}^{t+\delta} \int_{t}^{s} du \, dZ(s),$$

$$I_{(1,1)} = \int_{t}^{t+\delta} \int_{t}^{s} dZ(u) \, dZ(s).$$
(17)



As a result, we obtain the following approximation:

$$X(t+\delta) = X(t) + A(t,X(t))\delta + B(t,X(t))\Delta Z(t) + \mathcal{L}^{0}A(t,X(t))I_{(0,0)} + \mathcal{L}^{1}A(t,X(t))I_{(1,0)} + \mathcal{L}^{0}B(t,X(t))I_{(0,1)} + \mathcal{L}^{1}B(t,X(t))I_{(1,1)}.$$
(18)

Note:

$$I_{(0,0)} = \int_{t}^{t+\delta} (s-t)ds = \frac{1}{2} \delta^{2},$$

$$I_{(1,1)} = \int_{t}^{t+\delta} (Z(s) - Z(t))dZ(s) = \frac{1}{2} ((\Delta Z)^{2} - \delta),$$

$$I_{(0,1)} = \int_{t}^{t+\delta} (s-t)dZ(s) = \delta \Delta Z - I_{(1,0)},$$

$$I_{(1,0)} = \int_{t}^{t+\delta} (Z(s) - Z(t))ds.$$
(19)

• In particular, it is very fortuitous that $I_{(1,1)}$ can be computed in a closed, easy to simulate form.

Euler's scheme

- This approximation leads to practical discretization schemes of (11). We consider a sequence of times $0 = t_0 < t_1 < \ldots < t_m = T$.
- The first such scheme, Euler's scheme, consists in retaining the first three terms on the right hand side of (18):

$$X_{n+1} = X_n + A(t_n, X_n)\delta_n + B(t_n, X_n)\Delta Z_n,$$
(20)

where $\delta_n = t_{n+1} - t_n$, and $\xi_n \sim N(0,1)$. The random variables ξ_n are assumed independent.

Milstein's scheme

- In the second scheme, Milstein's scheme, in addition to the terms present in Euler's scheme, we also retain the last term on the right hand side of (18).
- Note that this term is of order of magnitude δ , while the three discarded terms are of order of magnitude $\delta^{3/2}$ and δ^2 .
- Explicitly, Milstein's scheme is given by

$$X_{n+1} = X_n + A(t_n, X_n)\delta_n + B(t_n, X_n)\Delta Z_n + \frac{1}{2}B(t_n, X_n)B'(t_n, X_n)(\Delta Z_n^2 - \delta_n),$$

where ' denotes the derivative with respect to x.



• We now consider an n-dimensional state variable $X \in \mathbb{R}^n$ driven by a d-dimensional Brownian motion $Z(t) \in \mathbb{R}^d$,

$$dX_{i}(t) = A_{i}(t, X(t))dt + \sum_{1 < a < d} B_{ia}(t, X(t))dZ_{a}(t),$$
(21)

where, for simplicity, we assume that the components of Z are independent.

This implies that

$$X_{i}(t+\delta) = X_{i}(t) + \int_{t}^{t+\delta} A_{i}(s,X(s))ds + \sum_{1 \leq a \leq d} \int_{t}^{t+\delta} B_{ia}(s,X(s))dZ_{a}(s). \tag{22}$$

 The following calculations generalize the calculations we carried out above for the case of a single factor SDE.



• If f(t, x) is twice continuously differentiable, then Ito's lemma states

$$df(t, X(t)) = \mathcal{L}^{0} f(t, X(t)) dt + \sum_{1 \le a \le d} \mathcal{L}^{a} f(t, X(t)) dZ_{a}(t),$$
 (23)

where the operators \mathcal{L}^i are defined by

$$\mathcal{L}^{0} = \frac{\partial}{\partial t} + \sum_{1 \le i \le n} A_{i} \frac{\partial}{\partial x_{i}} + \frac{1}{2} \sum_{1 \le i, j \le n} \sum_{1 \le a \le d} B_{ia} B_{ja} \frac{\partial^{2}}{\partial x_{i} \partial x_{j}}, \qquad (24)$$

and

$$\mathcal{L}^{a} = \sum_{1 \le i \le n} B_{ia} \frac{\partial}{\partial x_{i}}, \text{ for } a = 1, \dots, d.$$
 (25)

Applying Ito's lemma (23) to A_i yields

$$\begin{split} A_{i}(s,X(s)) &= A_{i}(t,X(t)) + \int_{t}^{s} \mathcal{L}^{0}A_{i}(u,X(u))du \\ &+ \sum_{1 \leq a \leq d} \int_{t}^{s} \mathcal{L}^{a}B_{ia}(u,X(u))dZ_{a}(u) \\ &\approx A_{i}(t,X(t)) + \mathcal{L}^{0}A_{i}(t,X(t)) \int_{t}^{s} du \\ &+ \sum_{1 \leq a \leq d} \mathcal{L}^{a}B_{ia}(t,X(t)) \int_{t}^{s} dZ_{a}(u) \,. \end{split}$$

We can thus approximate

$$\int_{t}^{t+\delta} A_{i}(s,X(s))ds \approx A_{i}(t,X(t))\delta + \mathcal{L}^{0}A_{i}(t,X(t))I_{(0,0)} + \sum_{1\leq a\leq d} \mathcal{L}^{a}B_{ia}(t,X(t))I_{(a,0)}.$$

Here

$$I_{(0,0)} = \int_{t}^{t+\delta} \int_{t}^{s} du \, ds,$$

$$I_{(a,0)} = \int_{t}^{t+\delta} \int_{t}^{s} dZ_{a}(u) \, ds,$$
(26)

are iterated integrals.



Similarly, we make the following approximation:

$$\begin{split} B_{ia}(s,X(s)) &= B_{ia}(t,X(t)) + \int_t^s \mathcal{L}^0 B_{ia}(u,X(u)) du \\ &+ \sum_{1 \leq b \leq d} \int_t^s \mathcal{L}^b B_{ia}(u,X(u)) dZ_b(u) \\ &\approx B_{ia}(t,X(t)) + \mathcal{L}^0 B_{ia}(t,X(t)) \int_t^s du \\ &+ \sum_{1 \leq b \leq d} \mathcal{L}^b B_{ia}(t,X(t)) \int_t^s dZ_b(u) \,. \end{split}$$

Therefore,

$$\int_{t}^{t+\delta} B_{ia}(s,X(s)) dZ_{a}(s) \approx B_{ia}(t,X(t)) \Delta Z_{a}(t) + \mathcal{L}^{0} B_{ia}(t,X(t)) I_{(0,a)}$$
$$+ \sum_{1 \leq b \leq d} \mathcal{L}^{b} B_{ia}(t,X(t)) I_{(a,b)},$$

where

$$I_{(0,a)} = \int_{t}^{t+\delta} \int_{t}^{s} du \, dZ_{a}(s) ,$$

$$I_{(a,b)} = \int_{t}^{t+\delta} \int_{t}^{s} dZ_{a}(u) \, dZ_{b}(s) .$$
(27)

• The integral I_(a,b), for a ≠ b, is known as the Levy area. There is no close form expression for the Levy area, and it is computationally expensive to simulate.

As a result, we obtain the following approximation:

$$X_{i}(t+\delta) = X_{i}(t) + A_{i}(t,X(t))\delta + \sum_{1 \leq a \leq d} B_{ia}(t,X(t))\Delta Z_{a}(t)$$

$$+ \mathcal{L}^{0}A_{i}(t,X(t))I_{(0,0)} + \sum_{1 \leq a \leq d} \mathcal{L}^{a}A_{i}(t,X(t))I_{(a,0)}$$

$$+ \sum_{1 \leq b \leq d} \left(\mathcal{L}^{0}B_{ib}(t,X(t))I_{(0,b)} + \sum_{1 \leq a \leq d} \mathcal{L}^{a}B_{ib}(t,X(t))I_{(a,b)} \right),$$
(28)

where

$$\Delta Z_a(t) = Z_a(t+\delta) - Z_a(t).$$

Note that:

$$I_{(0,0)} = \int_{t}^{t+\delta} (s-t)ds$$

$$= \frac{1}{2} \delta^{2},$$

$$I_{(a,a)} = \int_{t}^{t+\delta} (Z_{a}(s) - Z_{a}(t))dZ_{a}(s)$$

$$= \frac{1}{2} ((\Delta Z_{a})^{2} - \delta),$$

$$I_{(0,a)} = \int_{t}^{t+\delta} (s-t)dZ_{a}(s)$$

$$= \delta \Delta Z_{a} - I_{(a,0)},$$

$$I_{(a,0)} = \int_{t}^{t+\delta} (Z_{a}(s) - Z_{a}(t))ds$$
(29)

Note, in particular, that $I_{(a,a)}$ admits a simple, closed form expression.

Integrability condition

 In order to deal with the Levy areas I_(a,b), we impose the following integrability condition:

$$\mathcal{L}^a B_{ib} = \mathcal{L}^b B_{ia}, \tag{30}$$

or explicitly

$$\sum_{1 \le k \le n} B_{ka} \frac{\partial B_{ib}}{\partial x_k} = \sum_{1 \le k \le n} B_{kb} \frac{\partial B_{ia}}{\partial x_k} . \tag{31}$$

Note that then

$$I_{(a,b)} + I_{(b,a)} = \int_{t}^{t+\delta} \int_{t}^{s} (dZ_{a}(u) \ dZ_{b}(s) + dZ_{b}(u) \ dZ_{a}(s))$$

= $\Delta Z_{a} \Delta Z_{b}$. (32)

• In other words, the Levy areas $I_{(a,b)}$ and $I_{(b,a)}$ conspire to add up to a simple, easy to simulate expression!



Integrability condition

Therefore, when the integrability condition holds, (28) can be written as

$$X_{i}(t+\delta) = X_{i}(t) + A_{i}(t,X(t))\delta + \sum_{1 \leq a \leq d} B_{ia}(t,X(t))\Delta Z_{a}(t) + \frac{1}{2} \mathcal{L}^{0}A_{i}(t,X(t))\delta^{2} + \sum_{1 \leq a \leq d} \left(\left(\mathcal{L}^{a}A_{i}(t,X(t))\mathcal{L}^{0}B_{ia}(t,X(t))\right) I_{(a,0)} + \mathcal{L}^{0}B_{ia}(t,X(t))\Delta Z_{a}\delta + \sum_{1 \leq a \leq d} \left(\frac{1}{2} \mathcal{L}^{a}B_{ia}(t,X(t))(\Delta Z_{a}^{2} - \delta) + \sum_{a+1 \leq b \leq d} \mathcal{L}^{a}B_{ib}(t,X(t))\Delta Z_{a}\Delta Z_{b} \right).$$
(33)

This approximation leads to the following two discretization schemes.

Euler's scheme

 Euler's scheme is obtained by discarding all but the first three terms on the right hand side of (28):

$$X_{i,n+1} = X_{i,n} + A_{i,n}\delta_n + \sum_{1 \le a \le d} B_{ia,n} \Delta Z_{a,n}.$$
 (34)

• Euler's scheme is of order of convergence 1/2 meaning that the approximate solution converges in a suitable norm to the actual solution at the rate of $\delta t^{1/2}$, as $\delta t \equiv \max \delta t_0 \to 0$.

Milstein's scheme

Milstein's scheme includes the last term in (33):

$$X_{i,n+1} = X_{i,n} + A_{i,n}\delta_n + \sum_{1 \leq a \leq d} B_{ia,n} \Delta Z_{a,n}$$

$$+ \sum_{1 \leq a \leq d} \left(\frac{1}{2} \mathcal{L}^a B_{ia,n} (\Delta Z_{a,n}^2 - \delta_n) + \sum_{a+1 \leq b \leq d} \mathcal{L}^a B_{ib,n} \Delta Z_{a,n} \Delta Z_{b,n} \right).$$

This can be rewritten in a more symmetric form as:

$$X_{i,n+1} = X_{i,n} + \left(A_{i,n} - \frac{1}{2} \sum_{1 \le a \le d} \mathcal{L}^{a} B_{ia,n}\right) \delta_{n} + \sum_{1 \le a \le d} B_{ia,n} \Delta Z_{a,n}$$

$$+ \frac{1}{2} \sum_{1 \le a,b \le d} \mathcal{L}^{a} B_{ib,n} \Delta Z_{a,n} \Delta Z_{b,n}.$$
(35)

• Milstein's scheme is of order of convergence 1 meaning that the approximate solution converges in a suitable norm to the actual solution at the rate of δt , as $\delta t \to 0$.

Discretizing LMM: Euler's scheme

• We choose a sequence of event dates t_0, t_1, \ldots, t_m , and denote by $L_{jn} \simeq L_j(t_n)$ the approximate solution. We also set

$$\Delta_{j,n} = \Delta_j(t_n, L_n),$$

$$B_{ja,n} = B_{ja}(t_n, L_{j,n}),$$
(36)

and $\delta_n = t_{n+1} - t_n$.

Applied to LMM, Euler's scheme (34) reads:

$$L_{j,n+1} = L_{j,n} + \Delta_{j,n} \delta_n + \sum_{1 \le a \le d} B_{ja,n} \Delta Z_{a,n} , \qquad (37)$$

where, as before, $\Delta Z_{a,n} = Z_a(t_{n+1}) - Z_a(t_n)$ is the discretized Brownian motion.



Discretizing LMM: Milstein's scheme

- Fortunately, LMM is in the category of models which satisfy the integrability condition required for Milstein's scheme to work.
- In order to lighten up the notation, let us define:

$$\Upsilon_{jab,n} \equiv B_{ja}(t_n, L_{j,n}) \frac{\partial B_{jb}(t_n, L_{j,n})}{\partial L_j}.$$
 (38)

Then Milstein's scheme (35) applied to the LMM model reads:

$$L_{j,n+1} = L_{j,n} + \left(\Delta_{j,n} - \frac{1}{2} \sum_{1 \le a \le d} \Upsilon_{jaa,n}\right) \delta_{n} + \sum_{1 \le a \le d} B_{ja,n} \Delta Z_{a,n} + \frac{1}{2} \sum_{1 \le a,b \le d} \Upsilon_{jab,n} \Delta Z_{a,n} \Delta Z_{n,b}.$$
(39)

- A bit of a challenge lies in handling the drift terms. Because of their complexity, their calculation (at each time step) takes up to 50% of the total computation time.
- On the other hand, they are relatively small as compared to the initial values of the LIBOR forwards, and it would be desirable to develop an efficient methodology for accurate approximate evaluation of the drift terms.
- The first and simplest approach consist in "freezing" the values of $F_j(t)$ at the initial value $F_{j,0} \equiv F_j(0)$. We precompute the values

$$\Delta_{j,0} \equiv \Delta_j(t, F_0), \tag{40}$$

and use them for the drift terms throughout the simulation. This approximation, the frozen curve approximation, is rather crude, and does not perform very well when applied to pricing longer dated instruments.

 Going one step in the low noise expansion beyond the frozen curve approximation produces satisfying results.



 The second approach is a refinement of the frozen curve approximation, and consists in the following. From Ito's lemma,

$$\Delta_{j}(t, F(t)) = \Delta_{j,0} + \int_{0}^{t} \mathcal{L}^{a} \Delta_{j}(s, F(s)) dZ_{a}(s)$$

$$\simeq \Delta_{j,0} + \mathcal{L}^{a} \Delta_{j}(0, F_{0}) Z_{a}(t),$$
(41)

where we have suppressed all terms of order higher than 1/2.

We thus arrive at the following approximation, the order 1/2 approximation:

$$\Delta_{j,1/2}(t) \equiv \Delta_{j,0} + \mathcal{L}^a \Delta_j(0, F_0) Z_a(t). \tag{42}$$



- The coefficients $\mathcal{L}^a\Delta_j$ in the formula above are explicitly given by the following expressions.
- Under the forward measure Q_k:

$$\mathcal{L}^{a}\Delta_{j} = U_{ja}C_{j} \times \begin{cases} -\sum_{j+1 \leq i \leq k} \frac{\rho_{ji}\delta_{i}C_{i}}{1+\delta_{i}F_{i}} \left[U_{ja} \frac{\partial C_{j}}{\partial F_{j}} + U_{ia} \left(\frac{\partial C_{i}}{\partial F_{i}} - \frac{\delta_{i}C_{i}}{1+\delta_{i}F_{i}} \right) \right], & \text{if } j < k \\ 0, & \text{if } j = k, \\ \sum_{k+1 \leq i \leq j} \frac{\rho_{ji}\delta_{i}C_{i}}{1+\delta_{i}F_{i}} \left[U_{ja} \frac{\partial C_{j}}{\partial F_{j}} + U_{ia} \left(\frac{\partial C_{i}}{\partial F_{i}} - \frac{\delta_{i}C_{i}}{1+\delta_{i}F_{i}} \right) \right], & \text{if } j > k. \end{cases}$$

$$(43)$$

Under the spot measure:

$$\mathcal{L}^{a}\Delta_{j} = U_{ja}C_{j}\sum_{\gamma(t)< j < j} \frac{\rho_{ji}\delta_{i}C_{i}}{1 + \delta_{i}F_{i}} \left[U_{ja}\frac{\partial C_{j}}{\partial F_{j}} + U_{ia}\left(\frac{\partial C_{i}}{\partial F_{i}} - \frac{\delta_{i}C_{i}}{1 + \delta_{i}F_{i}}\right) \right]. \tag{44}$$



• The order 3/4 approximation, uses the next order term in the low noise expansion:

$$\Delta_{j,3/4}\left(t
ight)\equiv\Delta_{j,0}(t,L_{0})+\Gamma_{ja}Z_{a}\left(t
ight)+\Omega_{j}t.$$

Under the forward measure Q_k, the coefficients Γ_{ia} are given by:

$$\Gamma_{ja} = \begin{cases} -C_{j} \sum\limits_{j+1 \leq i \leq k} \frac{\rho_{ji} \delta_{i} C_{i}}{1 + \delta_{i} I_{i}} \left[U_{ja} \frac{\partial C_{j}}{\partial I_{j}} + U_{ia} \left(\frac{\partial C_{i}}{\partial I_{i}} - \frac{\delta_{i} C_{i}}{1 + \delta_{i} I_{i}} \right) \right], & \text{if } j < k, \\ 0, & \text{if } j = k, \\ C_{j} \sum\limits_{k+1 \leq i \leq j} \frac{\rho_{ji} \delta_{i} C_{i}}{1 + \delta_{i} I_{i}} \left[U_{ja} \frac{\partial C_{j}}{\partial I_{j}} + U_{ia} \left(\frac{\partial C_{i}}{\partial I_{i}} - \frac{\delta_{i} C_{i}}{1 + \delta_{i} I_{i}} \right) \right], & \text{if } j > k, \end{cases}$$

The coefficients Ω_i are given by:

$$\Omega_{j} = \begin{cases} -\sum\limits_{j+1 \leq i \leq k} \frac{\rho_{ji} \delta_{i}}{1 + \delta_{i} L_{i}} \left[\Delta_{j,0} C_{i} \frac{\partial C_{j}}{\partial L_{j}} + \Delta_{i,0} C_{j} \left(\frac{\partial C_{i}}{\partial L_{i}} - \frac{\delta_{i} C_{i}}{1 + \delta_{i} L_{i}} \right) \right], & \text{if } j < k, \\ 0, & \text{if } j = k, \\ \sum\limits_{k+1 \leq i \leq j} \frac{\rho_{ji} \delta_{i}}{1 + \delta_{i} L_{i}} \left[\Delta_{j,0} C_{i} \frac{\partial C_{j}}{\partial L_{j}} + \Delta_{i,0} C_{j} \left(\frac{\partial C_{i}}{\partial L_{i}} - \frac{\delta_{i} C_{i}}{1 + \delta_{i} L_{i}} \right) \right], & \text{if } j > k. \end{cases}$$

Under the spot measure,

$$\Gamma_{ja} = C_j \sum_{\gamma(t) \leq i \leq j} \frac{\rho_{ji} \delta_i C_i}{1 + \delta_i L_i} \Big[U_{ja} \frac{\partial C_j}{\partial L_j} + U_{ia} \Big(\frac{\partial C_i}{\partial L_i} - \frac{\delta_i C_i}{1 + \delta_i L_i} \Big) \Big],$$

and

$$\Omega_{j} = \sum_{\gamma(t) \leq i \leq j} \frac{\rho_{ji}\delta_{i}}{1 + \delta_{i}L_{i}} \left[\Delta_{j,0}C_{i}\frac{\partial C_{j}}{\partial L_{j}} + \Delta_{i,0}C_{j}\left(\frac{\partial C_{i}}{\partial L_{i}} - \frac{\delta_{i}C_{i}}{1 + \delta_{i}L_{i}}\right) \right].$$

The order 3/4 approximation leads to excellent accuracy.

On might easily refine this approach by computing terms of higher order in stochastic Taylor's expansion. This leads, however, to more complex and computationally expensive formulas, and the benefit of using an asymptotic expansion disappears. The order 1/2 approximation appears to offer the best performance versus accuracy profile.



References



Glasserman, P.: *Monte Carlo Methods in Financial Engineering*, Springer Verlag (2003).



Kloeden, P. E., and Platen, E.: *Numerical Solution of Stochastic Differential Equations*, Springer Verlag (1992).