Design of high-order short-time approximations as a problem of matching the covariance of a Brownian motion

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One of the outstanding problems in the numerical discretization of the Feynman–Kac formula calls for the design of arbitrary-order short-time approximations that are constructed in a stable way, yet only require knowledge of the potential function. In essence, the problem asks for the development of a functional analogue to the Gauss quadrature technique for one-dimensional functions. In PRE 69 (2004) 056701, it has been argued that the problem of designing an approximation of order $\nu$ is equivalent to the problem of constructing discrete-time Gaussian processes that are supported on finite-dimensional probability spaces and match certain generalized moments of the Brownian motion. Since Gaussian processes are uniquely determined by their covariance matrix, it is tempting to reformulate the moment-matching problem in terms of the covariance matrix alone. Here, we show how this can be accomplished.

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1 Introduction

Since its introduction, the Feynman–Kac formula [1–3] has played a fundamental role in the development of numerical algorithms capable of accounting for the physical properties of quantum systems made up of distinguishable particles. When utilized in tandem with the Monte Carlo integration technique, a powerful method is obtained: the path-integral Monte Carlo [4], which is capable of probing the quantum effects without any untestable approximations. The Feynman–Kac formula expresses the density matrix of a thermodynamic system as the expected value of a functional of the Brownian bridge

$$\rho(x, x'; \beta) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-(x'-x)^2/2\sigma^2} \mathbb{E} \left\{ -\beta \int_0^1 V[x_r(u) + \sigma B_u^0] du \right\}. \quad (1)$$

A second formulation of the Feynman–Kac formula is in terms of the full Brownian motion and reads

$$\langle x | e^{-\beta H} | \psi \rangle = \int_\mathbb{R} \rho(x, x'; \beta) \psi(x') dx' = \mathbb{E} \left\{ \exp\left( -\beta \int_0^1 V(x + \sigma B_u^0) du \right) \psi(x + \sigma B_1) \right\}, \quad (2)$$

where $\psi(x)$ is any square integrable function. In the above, $\rho(x, x'; \beta)$ is the density matrix for a one-dimensional canonical system characterized by the inverse temperature $\beta = 1/(k_BT)$ and made up of identical particles of mass $m_0$ moving in the
potential \( V(x) \). The stochastic element that appears in Eq. (1), \( \{ B_{u}, 0 \leq u \leq 1 \} \), is a so-called standard Brownian bridge, defined as follows: if \( \{ B_{u}, u \geq 0 \} \) is a standard Brownian motion starting at zero, then the Brownian bridge is the stochastic process \( \{ B_{u}, 0 \leq u \leq 1 \mid B_{1} = 0 \} \), i.e., a Brownian motion conditioned on the event \( B_{1} = 0 \). As is well known, a Brownian bridge can be realized as the process \( \{ B_{u} - uB_{1}, 0 \leq u \leq 1 \} \) [5]. To complete the description of Eqs. (1) and (2), we have \( x_{r}(u) = x + (x' - x)u \) and \( \sigma = (\hbar^{2}/m\beta_{0})^{1/2} \).

A problem of direct interest to the chemical physicist is the development of approximations supported on finite-dimensional probability spaces that have fast convergence for smooth enough potentials, as measured against the number of evaluations of the potential function. Most desirably, such approximations should utilize only the potential function in their construction. For reasons of stability, they should converge, perhaps at a slower rate, for all continuous potentials that are bounded from below. Until recently, the fastest method available (as order of convergence) has been the trapezoidal Trotter discrete path integral method. The technique is usually derived by means of the Lie–Trotter product formula and an appropriate short-time high-temperature approximation. The formal asymptotic convergence of the trapezoidal Trotter method and of related techniques was extensively studied by Suzuki [6, 7] for bounded operators and by Ichinose and Tamura [8, 9], among others. In particular, results of the last two authors [9] show that the symmetric Trotter–Suzuki approximation has optimal convergence \( O(1/n^{2}) \) for sufficiently smooth potentials as far as pointwise convergence of their integral kernels is concerned. This type of convergence is also implied in the present paper. On the other hand, the non-existence theorem of Suzuki [6] makes it implausible that faster convergence can be achieved by utilizing short-time approximations constructed as functions of the kinetic and potential operators.

Recently, a more general approach has been put forward by the present author [10], who argued that, for sufficiently smooth potentials, there might exist direct short-time high-temperature approximations of arbitrary polynomial convergence order. Of course, these short-time approximations are generally not functions of the kinetic and potential operators. The construction of such approximations is based upon an “experimental” theorem on the pointwise convergence of their integral kernels of the Lie–Trotter product formulas. Although not rigorously proved, this theorem seems quite plausible. The short-time approximations considered are based on carefully designed finite-dimensional approximations to the Brownian motion entering the Feynman–Kac formula. Basically, the Brownian motion is replaced by some discrete-time Gaussian process that is supported on a finite-dimensional probability space. A set of functional equations involving some generalized moments of the Gaussian process have been shown to control the order of convergence \( \nu \). Because the number of equations increases in an exponential fashion with \( \nu \), explicit solutions have been obtained only for \( \nu = 3, 4 \).

In the present work, we exploit the fact that both the Brownian motion and its replacement are Gaussian processes and are, therefore, uniquely determined by their covariance matrices. We thus show how to express the functional equations in terms of the covariance matrices alone. Hopefully, the new equations will prove
A problem of matching the covariance of a Brownian motion
easier to utilize in a complete mathematical proof of the existence of short-time
approximations of arbitrary order. In the appendix, we give a general convergence
theorem regarding the construction of finite-dimensional discrete approximations
to the Feynman–Kac formula.

2 Statement of the moment-matching problem

Perhaps one of the oldest techniques for simulating a Brownian bridge (or Browni-
nan motion) is via random series. As such, let \( \{ \lambda_k(\tau) \}_{k \geq 0} \) be any orthonormal basis
in \( L^2[0, 1] \) such that \( \lambda_0(\tau) \equiv 1 \), let

\[
\Lambda_k(u) = \int_0^u \lambda_k(\tau) d\tau \quad \text{for} \quad k \geq 0
\]

and let \( \bar{a} := \{ a_0, a_1, \ldots \} \) be a sequence of independent identically distributed stan-

dard normal random variables. By the Ito–Nisio theorem [11], the random series
\[
\sum_{k \geq 0} a_k \Lambda_k(u)
\]

is uniformly convergent almost surely and equal in distribution with a
standard Brownian motion \( B_u \) starting at zero. By the construction of a Brow-
nian bridge as the process \( \{ B_u - uB_1, 0 \leq u \leq 1 \} \) and the fact that \( \Lambda_0(u) = u \), it

follows that \( \sum_{k \geq 1} a_k \Lambda_k(u) = \sum_{k \geq 0} a_k \Lambda_k(u) - a_0 u \) is equal in distribution with a
standard Brownian bridge. If \( \Omega \) is the set of all sequences \( \bar{a} := \{ a_0, a_1, \ldots \} \) and if

\[
dP[\bar{a}] = \prod_{k=0}^{\infty} d\mu(a_k) \quad \text{with} \quad d\mu(z) = (2\pi)^{1/2} \exp(-z^2/2)dz
\]

is the probability measure on \( \Omega \) associated with the sequence of independent random
variables \( \bar{a} := \{ a_0, a_1, \ldots \} \), then the Feynman–Kac formula given by Eq. (1) reads

\[
\rho(x, x', \beta) = \rho_{fp}(x, x', \beta) \int_\Omega dP[\bar{a}] \exp\left\{ -\beta \int_0^1 \left[ x_r(u) + \sigma \sum_{k=1}^\infty a_k \Lambda_k(u) \right] du \right\}.
\]

Here, \( \rho_{fp}(x, x', \beta) = \exp[-(x' - x)^2/2\sigma^2]/(2\pi\sigma^2)^1/2 \) is recognized as the density
matrix of a free particle. The alternative formulation given by Eq. (2) becomes

\[
\rho(x, x', \beta) = \int_\Omega dP[\bar{a}] \exp\left\{ -\beta \int_0^1 \left[ x + \sigma \sum_{k=0}^\infty a_k \Lambda_k(u) \right] du \right\} \psi(x + \sigma a_0).
\]

Eqs. (3) and (4) are appropriately called random series representations of the
Feynman–Kac formula in the chemical-physics literature [12, 13].

Eq. (3) is suggestive of some sort of numerical approximation to the Feynman–Kac formula, namely

\[
\rho_n(x, x', \beta) = \rho_{fp}(x, x', \beta) \int_\Omega dP[\bar{a}] \exp\left\{ -\beta \sum_{i=1}^{n_a} w_i V \left[ x_r(\theta_i) + \sigma \sum_{k=1}^{n_a} a_k \Lambda_k(\theta_i) \right] \right\}.
\]

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\]
Here, the non-negative weights \( w_i \) (we assume \( \sum_i w_i = 1 \)) and the knots \( \theta_i \) define a quadrature technique on the interval \([0, 1]\). It goes without saying that such approximations are convergent as \( n_q \to \infty \) and \( n_\nu \to \infty \) under mild assumptions for the potential \( V(x) \): boundness from below and continuity. In order not to disrupt from the flow of the presentation, we give the simple proof in the appendix (see Corollary 1). Thus, we are in no shortage of quadrature formulas. We just want the faster ones.

Another utilization of Eq. (5) is as a short-time approximation in a Lie–Trotter product of the form

\[
\rho_n(x, x'; \beta) = \int_\mathbb{R} dx_1 \ldots \int_\mathbb{R} dx_n \rho_0^{(\nu)} \left( x, x_1; \frac{\beta}{n + 1} \right) \ldots \rho_0^{(\nu)} \left( x_n, x'; \frac{\beta}{n + 1} \right).
\]

(6)

The gain in interpreting Eq. (6) as a short-time approximation is that the requirement that the functions \( \{ \Lambda_k(u); 1 \leq k \leq n_\nu \} \) are constructed according to the Ito–Nisio prescription can be relaxed. We denote this by utilizing a tilde overscript, so that the short-time approximation reads

\[
\rho_0^{(\nu)}(x, x'; \beta) = \rho_{fp}(x, x'; \beta) \int_\mathbb{R} d\mu(a_1) \ldots \int_\mathbb{R} d\mu(a_{n_\nu}) \times \\
\times \exp \left\{ -\beta \int_0^1 V \left[ x_r(u) + \sigma \sum_{k=1}^{n_\nu} a_k \tilde{\Lambda}_k(u) \right] d\omega(u) \right\}.
\]

(7)

The functions \( \{ \tilde{\Lambda}_k(u); 1 \leq k \leq n_\nu \} \) are required to be continuous (that is, bounded on the quadrature knots) and vanish at both end points. At this moment, we should emphasize that this special choice of short-time approximation is made while bearing in mind its usefulness in Monte Carlo simulations. Due to some special properties, especially the availability of the fast sampling algorithm [14], the subsequence of Lie–Trotter products with \( n = 2^k - 1 \) is of utmost practical interest. A simple proof of the convergence of this Lie–Trotter subsequence is given in the appendix (see Corollary 2), again for continuous and bounded from below potentials.

A second requirement that we ask of the construction given by Eq. (7) has to do with the symmetry of the density matrix, which should reflect itself in the symmetry of the short-time approximation. As such, we require that the discrete probability measure

\[
d\omega(u) = \sum_{i=1}^{n_\nu} w_i \delta(u - \theta_i) du
\]

(8)

defining the quadrature technique on \([0, 1]\) must be symmetric about \( \frac{1}{2} \). Also, the finite dimensional process \( \sum_{k=1}^{n_\nu} a_k \tilde{\Lambda}_k(u) \) must be invariant under the transformation \( u' = 1 - u \). That is, we require the equality in distribution

\[
\sum_{k=1}^{n_\nu} a_k \tilde{\Lambda}_k(u) \overset{d}{=} \sum_{k=1}^{n_\nu} a_k \tilde{\Lambda}_k(1 - u).
\]

(9)
The time symmetry of the process \( \sum_{k=1}^{\nu} a_k \tilde{\Lambda}_k(u) \) can be enforced, for example, by restricting the functions \( \{\tilde{\Lambda}_k(u); 1 \leq k \leq \nu\} \) to the class of symmetric and antisymmetric functions. To understand these requirements, notice that the Hermiticity of the density matrix stems from the symmetry of the Lebesgue measure on \([0, 1]\) as well as from the time symmetry of the standard Brownian bridge \( B^0_0 \), i.e., the fact that \( \{B^0_1 \cdot u; 0 \leq u \leq 1\} \) is a Brownian bridge equal in distribution to \( \{B^0_0; 0 \leq u \leq 1\} \). Because the random sum \( \sum_{k=1}^{\nu} a_k \tilde{\Lambda}_k(u) \) is intended as a replacement for the Brownian bridge, it is convenient to introduce the notations

\[
\tilde{B}^0_u \equiv \sum_{k=1}^{\nu} a_k \tilde{\Lambda}_k(u) \quad \text{and} \quad \tilde{B}_u \equiv a_0 u + \tilde{B}^0_u \quad \text{for } 0 \leq u \leq 1.
\]  

(10)

Thus, the Brownian bridge (motion) is approximated by a simple Gaussian process that is supported on a finite-dimensional probability space. Moreover, only the values of the process for the discrete times represented by the quadrature knots \( \theta_1, \theta_2, \ldots, \theta_{\nu} \) are relevant for the construction of the short-time approximation.

The symbol \( (\nu) \) appearing as a superscript in Eq. (7) denotes the order of convergence of the short-time approximation. This is defined as the largest \( \nu \) for which the so-called convergence operator \( T_\nu \psi \) expressed by

\[
(T_\nu \psi)(x) = \lim_{\beta \to 0^+} \frac{\int_{\mathbb{R}} \left[ \rho^{(\nu)}_0(x, x'; \beta) - \rho(x, x'; \beta) \right] \psi(x') \, dx'}{\beta^{\nu+1}}.
\]  

(11)

is well defined at least on the class of infinitely differentiable and compactly supported functions \( \psi(x) \). In Ref. [10], it is claimed but not rigorously demonstrated that

\[
\lim_{n \to \infty} (n + 1)^\nu \left[ \rho_n(x, x'; \beta) - \rho(x, x'; \beta) \right] = \beta^{\nu+1} \int_0^1 \left\langle x \middle| e^{-\theta \beta H} T_\nu e^{-(1-\theta)\beta H} \middle| x' \right\rangle \, d\theta,
\]  

(12)

where \( \rho_n(x, x'; \beta) \) is defined by Eq. (6). A rigorous proof of Eq. (12) is beyond the mathematical abilities of the present author. Based on the more or less formal arguments presented in the aforementioned reference, it is very likely that the statement is true. The author would be very grateful to the mathematically more inclined reader who may want to investigate the problem and prove or disprove the assertion (again, the case \( n = 2^k - 1 \) suffices for all practical purposes).

Nonetheless, Eq. (12) states that the convergence of the Lie–Trotter product is as fast as \( 1/n^\nu \), which explains the nomenclature regarding the order of convergence. We should emphasize that merely the existence of the convergence operator expressed by Eq. (11) sets some constraints on the smoothness of the potential function. The natural class of potentials to study the problem of constructing short-time approximations of arbitrary order is the class of continuously and infinitely differentiable functions for which

\[
\frac{1}{\sqrt{2\pi\alpha}} \int_{\mathbb{R}} e^{-z^2/(2\alpha)} \left| V^{(k)}(x+z) \right|^2 \, dz < \infty,
\]  

(13)
for all $x \in \mathbb{R}$ and $\alpha > 0$ and for all integers $k \geq 0$ and $j \geq 1$. This condition is necessary in order to ensure that we recover the original potential $V(x)$, derivatives $V^{(k)}(x)$, or products of such functions from their Gaussian transforms, in the limit that $\alpha \to 0$ (see Theorem 3 of Ref. [15]). We mention that this class of potentials does not include some pathological infinitely differentiable and bounded from below potentials such as $\exp(x^4)$ or even $\cos(\exp(x^4))$.

In these conditions, according to Theorem 4 of Ref. [10], the convergence operator $T_\nu$ exists if and only if

$$
E \left[ (B_1)^{j_1} (M_0)^{j_2} (M_1)^{j_3} \cdots (M_{2\mu - 2})^{j_{2\mu}} \right] = E \left[ (\tilde{B}_1)^{j_1} (\tilde{M}_0)^{j_2} (\tilde{M}_1)^{j_3} \cdots (\tilde{M}_{2\mu - 2})^{j_{2\mu}} \right]
$$

(14)

for all $2\mu$-tuples of non-negative integers $(j_1, j_2, \ldots, j_{2\mu})$ such that $\sum_{k=1}^{2\mu} kj_k = 2\mu$ and $1 \leq \mu \leq \nu$. The random variables $M_k$ and $\tilde{M}_k$ are defined by

$$
M_k \equiv \int_0^1 (B_u)^k du \quad \text{and} \quad \tilde{M}_k \equiv \int_0^1 (\tilde{B}_u)^k d\omega(u) = \sum_{i=1}^{n_q} w_i (\tilde{B}_{\theta_i})^k,
$$

(15)

respectively.

3 Reformulation of the generalized moment conditions in terms of covariance matrices

Since they are Gaussian processes with continuous paths, the Brownian motion and its approximation are uniquely determined by the covariance matrices

$$
\gamma(u, \tau) = E(B_u B_\tau) = u\tau + \sum_{k=1}^{\infty} \Lambda_k(u) \Lambda_k(\tau) = \min\{u, \tau\}
$$

(16)

and

$$
\tilde{\gamma}(u, \tau) = E(\tilde{B}_u \tilde{B}_\tau) = u\tau + \sum_{k=1}^{n_\mu} \tilde{\Lambda}_k(u) \tilde{\Lambda}_k(\tau),
$$

(17)

respectively. As such, at least in principle, the relations given by Eq. (14) can be formulated in terms of these covariance matrices alone. In this section, we show how this can be done.

Let $J_{2\mu}$ denote the set of solutions of the Diophantine equation $\sum_{k=1}^{2\mu} kj_k = 2\mu$. For each $\zeta \in J_{2\mu}$, we define the integer

$$
d(\zeta) = j_3 + j_4 + \cdots + j_{2\mu}
$$

(18)

and

$$
n(\zeta) = [j_1 + j_3 + 2j_4 + \cdots + (2\mu - 2)j_{2\mu}] = \mu - j_2 - d(\zeta),
$$

(19)

respectively. We also define the differentiation functional $D_\zeta$ acting on the space of infinitely differentiable functions $f(\lambda_0, \lambda_1, \ldots, \lambda_{d(\zeta)})$ that associates to each $f$ the
A problem of matching the covariance of a Brownian motion following partial derivative evaluated at the origin

\[ D \zeta f \equiv \frac{\partial^{j_1}}{\partial \lambda_0^{j_1}} \frac{\partial^{j_2}}{\partial \lambda_1 \partial \lambda_2} \cdots \frac{\partial^{j_{\mu}}}{\partial \lambda_{(2\mu-2)j_{2\mu}}} \cdots \frac{\partial^{2j_4}}{\partial \lambda_{j_3+1} \partial \lambda_{j_3+2} \cdots \partial \lambda_{j_3+j_4}} \cdots \frac{\partial^{j_3}}{\partial \lambda_{(2\mu-2)j_{2\mu}+1} \partial \lambda_{(2\mu-2)j_{2\mu}+2} \cdots \partial \lambda_{(2\mu-2)d(\zeta)}} f(0,0,\ldots,0). \]  

(20)

The differential order of \( D \zeta \) is \( j_1 + j_3 + 2j_4 + \cdots + (2\mu-2)j_{2\mu} = 2n(\zeta) \).

The reader will understand the need for this rather cumbersome notation shortly. Going back to Eq. (14), let us notice that the equality

\[ M_0 = \int_0^1 1du = 1 = \int_0^1 1d\omega(u) = \tilde{M}_0 \]

means that the factors containing \( M_0 \) and \( \tilde{M}_0 \) cancel out. Next, we utilize the identities

\[ (B_1)^{j_1} = j_1! \frac{d^{j_1}}{d\lambda^{j_1}} e^{\lambda B_1} \bigg|_{\lambda=0} \quad \text{and} \quad M_k = k! \int_0^1 du \frac{d^k}{d\lambda^k} e^{\lambda B_u} \bigg|_{\lambda=0} \]

to express the first factor in Eq. (14) as

\[ j_1!(1!)^{j_3} \cdots ((2\mu-2)!)^{j_{2\mu}} E \left[ \frac{D^{j_1}}{d\lambda^{j_1}} e^{\lambda B_1} \bigg|_{\lambda=0} \right] \left( \int_0^1 du \frac{d^{j_3}}{d\lambda^{j_3}} e^{\lambda B_u} \bigg|_{\lambda=0} \right) \cdots \left( \int_0^1 du \frac{d^{j_{2\mu}}}{d\lambda^{j_{2\mu}}} e^{\lambda B_u} \bigg|_{\lambda=0} \right)^{j_{2\mu}}. \]  

(21)

Expanding the parenthesis and interchanging the order of integration and differentiation, we obtain the result

\[ j_1!(1!)^{j_3} \cdots ((2\mu-2)!)^{j_{2\mu}} D \zeta f_\zeta, \]

where

\[ f_\zeta(\lambda_0, \lambda_1, \ldots, \lambda_{d(\zeta)}) = \int_0^1 du_1 \cdots \int_0^1 du_{d(\zeta)} E \exp \left( \sum_{i=0}^{d(\zeta)} \lambda_i B_{u_i} \right) \]

and \( u_0 = 1 \).

To evaluate the function \( f_\zeta \), one may utilize any random series and compute

\[ E \exp \left( \sum_{j=0}^{d(\zeta)} \lambda_i B_{u_i} \right) = E \exp \left( \sum_{i=0}^{d(\zeta)} \lambda_i \sum_{k=0}^{\infty} a_k \Lambda_k(u_i) \right) = \prod_{k=0}^{\infty} \left\{ \int d\mu(z) \exp \left( z \sum_{i=0}^{d(\zeta)} \lambda_i \Lambda_k(u_i) \right) \right\} = \prod_{k=0}^{\infty} \exp \left\{ \frac{1}{2} \sum_{i=0}^{d(\zeta)} \lambda_i \Lambda_k(u_i)^2 \right\}. \]  

7
The last term equals
\[
\exp \left\{ \frac{1}{2} \sum_{i=0}^{d(\zeta)} \sum_{j=0}^{d(\zeta)} \lambda_i \lambda_j \left( \sum_{k=0}^{\infty} \Lambda_k(u_i) \Lambda_k(u_j) \right) \right\} = \exp \left\{ \frac{1}{2} \sum_{i=0}^{d(\zeta)} \sum_{j=0}^{d(\zeta)} \lambda_i \lambda_j \gamma(u_i, u_j) \right\}.
\]

Therefore, the explicit expression of the function \(f_\zeta\) is
\[
f_\zeta(\lambda_0, \lambda_1, \ldots, \lambda_{d(\zeta)}) = \int_0^1 du_1 \cdots \int_0^1 du_{d(\zeta)} \exp \left\{ \frac{1}{2} \sum_{i=0}^{d(\zeta)} \sum_{j=0}^{d(\zeta)} \lambda_i \lambda_j \gamma(u_i, u_j) \right\}.
\]

We remind the reader that \(u_0 = 1\).

The function \(f_\zeta\) can be replaced by some \((1 + d(\zeta))\)-dimensional polynomial of degree \(2n(\zeta)\). Indeed, starting from the expansion
\[
\exp \left\{ \frac{1}{2} \sum_{i=0}^{d(\zeta)} \sum_{j=0}^{d(\zeta)} \lambda_i \lambda_j \gamma(u_i, u_j) \right\} = \sum_{k=0}^{\infty} \frac{1}{2^k k!} \left[ \sum_{i=0}^{d(\zeta)} \sum_{j=0}^{d(\zeta)} \lambda_i \lambda_j \gamma(u_i, u_j) \right]^k
\]

one observes that, upon the action of \(D_\zeta\), the differential order of which is \(2n(\zeta)\), only the term with \(k = n(\zeta)\) survives. The terms of lower degree are washed out by the process of differentiation, whereas the terms of higher degrees cancel when the respective derivatives are evaluated at the origin. Thus, the left-hand side of Eq. (14) is given by the expression
\[
\frac{j_1!(1!)^j \cdots ((2\mu - 2)!)^j 2^{2\mu}}{2^{n(\zeta)n(\zeta)!}} D_\zeta f_\zeta,
\]

where
\[
f_\zeta(\lambda_0, \lambda_1, \ldots, \lambda_{d(\zeta)}) = \int_0^1 du_1 \cdots \int_0^1 du_{d(\zeta)} \left[ \sum_{i=0}^{d(\zeta)} \sum_{j=0}^{d(\zeta)} \lambda_i \lambda_j \gamma(u_i, u_j) \right]^{n(\zeta)}.
\]

In an analogue manner, one demonstrates that the right-hand side of Eq. (14) can be written as
\[
\frac{j_1!(1!)^j \cdots ((2\mu - 2)!)^j 2^{2\mu}}{2^{n(\zeta)n(\zeta)!}} D_\zeta \tilde{f}_\zeta,
\]

where
\[
\tilde{f}_\zeta(\lambda_0, \lambda_1, \ldots, \lambda_{d(\zeta)}) = \int_0^1 d\omega(u_1) \cdots \int_0^1 d\omega(u_{d(\zeta)}) \left[ \sum_{i=0}^{d(\zeta)} \sum_{j=0}^{d(\zeta)} \lambda_i \lambda_j \tilde{\gamma}(u_i, u_j) \right]^{n(\zeta)}.
\]

Comparing Eqs. (22) and (24), we see that Eq. (14) is equivalent to the equality
\[
D_\zeta f_\zeta = D_\zeta \tilde{f}_\zeta
\]

with \(f_\zeta\) and \(\tilde{f}_\zeta\) defined by Eqs. (23) and (25), respectively.
As Eqs. (22) and (24) show, the polynomials \( f_\zeta \) and \( \tilde{f}_\zeta \) depend only on some integers \( d(\zeta) \) and \( n(\zeta) \) that have the property
\[
d(\zeta) + n(\zeta) \leq \mu \leq \nu.
\]
Thus, a sufficient condition that the equality expressed by Eq. (26) holds for all \( \zeta \in J_{2\mu} \) and \( 1 \leq \mu \leq \nu \) is that the equality of polynomials
\[
\tilde{f}_{n,d}(\lambda_0, \lambda_1, \ldots, \lambda_d) \equiv \int_0^1 \omega(u_1) \cdots \int_0^1 \omega(u_d) \left[ \sum_{i=0}^{d} \sum_{j=0}^{d} \lambda_i \lambda_j \gamma(u_i, u_j) \right]^\mu = \int_0^1 du_1 \cdots \int_0^1 du_d \left[ \sum_{i=0}^{d} \sum_{j=0}^{d} \lambda_i \lambda_j \gamma(u_i, u_j) \right]^\mu = f_{n,d}(\lambda_0, \lambda_1, \ldots, \lambda_d)
\]
holds for all integer \( d \) and \( n \) such that \( d + n \leq \nu \). If \( d_1 \leq d_2 \) then the polynomial of dimension \( 1 + d_1 \) is a particular case of the polynomial of dimension \( 1 + d_2 \), with the last \( d_2 - d_1 \) lambda’s set to zero. Thus, it is enough to check the above equality for the cases with \( d + n = \nu \).

Let us show that checking the equality expressed by Eq. (27) for all integers \( n \) and \( d \) such that \( n + d = \nu \) is also a necessary condition. Notice that the polynomials appearing in Eq. (27) are symmetric under the permutation of the variables \( \lambda_1, \lambda_2, \ldots, \lambda_d \). This is due to the symmetry of the covariance matrices \( \gamma(u, \tau) \) and \( \tilde{\gamma}(u, \tau) \). In these conditions, the differentiation functional given by Eq. (20) is the most general differential expression of order \( 2n \), provided that we let the index \( \zeta = \{ j_2, j_3, j_4, \ldots, j_{2\nu} \} \) lie in the set \( S \) that contains all possible indexes for which \( j_2 = 0, j_3 + j_4 + \cdots + j_{2\nu} = d \), and \( j_1 + j_3 + 2j_4 + \cdots + (2\nu - 2)j_{2\nu} = 2n \). Following some permutation of the variables \( \lambda_1, \lambda_2, \ldots, \lambda_d \), any other differential expression can be obtained from and, by the aforementioned symmetry of polynomials, is equal to a differential expression of the type given by Eq. (20) for some \( \gamma \in S \). Therefore, the equality \( D_\zeta f_{n,d} = D_\zeta \tilde{f}_{n,d} \) for all \( \zeta \in S \) implies \( f_{n,d} = \tilde{f}_{n,d} \). It remains to prove that, as \( \zeta \) spans \( J_{2\nu} \), it also spans \( S \). Clearly, any index \( \zeta \in S \) has the property \( j_1 + 2j_2 + 3j_3 + \cdots + 2\nu j_{2\nu} = 2(n + d) = 2\nu \). Therefore, \( S \subset J_{2\nu} \) and the claim of necessity is proved.

We summarize the results of the paper in the following proposition.

**Proposition 1** A short-time approximation of the type given by Eq. (7) has convergence order \( \nu \) if and only if the equality of polynomials expressed by Eq. (27) holds for all integers \( d \) and \( n \) such that \( d + n = \nu \).

The above proposition can be utilized together with the multinomial formula to generate the necessary conditions that the trial covariance matrix must satisfy in order for a short-time approximation to have convergence order \( \nu \). Most likely, more useful statements can be obtained by the mathematically more inclined reader. Nevertheless, we have achieved our goal of formulating the conditions in terms of the covariance matrices alone.

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Appendix

In this appendix, we establish an easily provable theorem that is useful in many situations and provides the basic mathematical background for the discretization of the Feynman–Kac formula. We then utilize the theorem to demonstrate several statements made in the text.

To begin with, let us formalize what we mean by quadrature rules. By a quadrature scheme on \([0, 1]\), we understand a sequence of pairs of vectors of variable lengths \(q_n\), namely the weights \(\{w_{n,1}, w_{n,2}, \ldots, w_{n,q_n}\}\) and the knots \(0 \leq u_{n,1} < u_{n,2} < \cdots < u_{n,q_n} \leq 1\), constructed such that the \(w_{n,i}\)'s are non-negative for all \(1 \leq i \leq q_n\) and \(n \geq 1\) and such that

\[
\lim_{n \to \infty} \sum_{i=1}^{q_n} w_{n,i} h(u_{n,i}) = \int_0^1 h(u) \, du \tag{28}
\]

for all continuous functions \(h : [0, 1] \to \mathbb{R}\). The mid-point, trapezoidal, Simpson, and Gauss–Legendre rules are well-known examples of quadrature schemes. The non-negativity of the weights is a stability requirement that simply says that it is not all right to get a negative answer if a positive function is integrated.

Let \(\{S_n(\omega; u); 0 \leq u \leq 1\}_{n \geq 1}\) be a family of random processes with continuous paths supported on a probability space \((\Omega, \mathcal{F}, P)\) and such that \(S_n(\omega; u)\) converges to some limit \(B^0_u(\omega)\) uniformly in \(u\), \(P\)-almost surely. Assume that the limit \(B^0_u(\omega)\) is equal in distribution to a Brownian bridge. Given an arbitrary quadrature rule, we can define a sequence of approximations to the density matrix by the prescription

\[
\rho_n(x, x'; \beta) = \rho_{fp}(x, x'; \beta) \mathbb{E} \exp \left\{ -\beta \sum_{i=1}^{q_n} w_{n,i} V(x_r(u_{n,i}) + \sigma S_n(\omega; u_{n,i})) \right\}, \tag{29}
\]

for \(n \geq 1\). We shall call the above prescription the standard discretization of the Feynman–Kac formula. The nomenclature is motivated by the following fundamental theorem.

**Theorem 1** If \(V(x)\) is continuous and bounded from below, then the sequence \(\rho_n(x, x'; \beta)\) is bounded by \(C(\beta)\rho_{fp}(x, x'; \beta)\), for some positive \(C(\beta) < \infty\), and

\[
\rho_n(x, x'; \beta) \to \rho(x, x'; \beta), \text{ as } n \to \infty. \tag{30}
\]

**Proof.** Let \(V_0\) be a non-positive lower bound for \(V(x)\). Eq. (28) specialized for \(h(x) = 1\) says that the sequence \(\sum_{i=1}^{q_n} w_{n,i}\) is convergent, thus bounded, say by \(c > 0\). Then the integrand of Eq. (29) is bounded from above by \(C(\beta) = \exp(-c\beta V_0)\) and we have

\[
\rho_n(x, x'; \beta) \leq C(\beta)\rho_{fp}(x, x'; \beta). \tag{31}
\]
The remainder of the theorem follows from the dominated convergence theorem and the Feynman–Kac formula, as soon as we prove that

$$\lim_{n \to \infty} \sum_{i=1}^{q_n} w_{n,i} V[x_r(u_{n,i}) + \sigma S_n(\omega; u_{n,i})] = \int_0^1 V[x_r(u) + \sigma B_u^0(\omega)]du, \text{ a.s.} \quad (32)$$

As such, let \( \omega \in \Omega \) be given. Since \( x_r(u) + \sigma B_u^0(\omega) \) is almost surely continuous in \( u \) — as uniform limit of continuous functions \( x_r(u) + \sigma S_n(\omega; u) \) — it is bounded by some constant \( M > 0 \). In fact, by uniform convergence, we have \( |x_r(u) + \sigma S_n(\omega; u)| \leq 2M \) for large enough \( n \). Since \( V(x) \) is continuous, it is uniformly continuous on the compact set \( |x| \leq 2M \). Consequently, given \( \epsilon > 0 \), there is \( \eta > 0 \) such that \( |V(x) - V(y)| < \epsilon \) whenever \( |x - y| < \eta \). Nevertheless, by uniform convergence,

$$\|[x_r(u) + \sigma S_n(\omega; u)] - [x_r(u) + \sigma B_u^0(\omega)]\| < \eta$$

for large enough \( n \) and so,

$$|V[x_r(u) + \sigma S_n(\omega; u)] - V[x_r(u) + \sigma B_u^0(\omega)]| < \epsilon.$$ 

Since \( \sum_{i=1}^{q_n} w_{n,i} < \epsilon \), it follows that

$$\left| \sum_{i=1}^{q_n} w_{n,i} V[x_r(u_{n,i}) + \sigma S_n(\omega; u_{n,i})] - \sum_{i=1}^{q_n} w_{n,i} V[x_r(u_{n,i}) + \sigma B_{u_{n,i}}^0(\omega)] \right| < \epsilon.$$ 

By Eq. (28), and the continuity of \( V[x_r(u) + \sigma B_u^0(\omega)] \) as a function in \( u \), we also have

$$\left| \int_0^1 V[x_r(u) + \sigma B_u^0(\omega)]du - \sum_{i=1}^{q_n} w_{n,i} V[x_r(u_{n,i}) + \sigma B_{u_{n,i}}^0(\omega)] \right| < \epsilon \quad (34)$$

for \( n \) sufficiently large. Combining Eqs. (33) and (34), we obtain

$$\left| \sum_{i=1}^{q_n} w_{n,i} V[x_r(u_{n,i}) + \sigma S_n(\omega; u_{n,i})] - \int_0^1 V[x_r(u) + \sigma B_u^0(\omega)]du \right| < (1 + c)\epsilon.$$ 

Since \( \epsilon \) is arbitrary, the almost sure convergence appearing in Eq. (32) is demonstrated and the proof of the theorem is concluded.

**Observation.** Because \( \rho_n(x, x'; \beta) \) is bounded by \( C(\beta)\rho_{fp}(x, x'; \beta) \), the dominated convergence theorem and the above theorem also imply convergence in the strong topology, that is,

$$\int_{\mathbb{R}} \rho_n(x, x'; \beta) \psi(x')dx' \to \int_{\mathbb{R}} \rho(x, x'; \beta) \psi(x')dx'$$

for all square integrable \( \psi(x) \). In fact, by choosing the sequence of processes \( S_n(\omega, n) \) to be constant and equal in distribution to \( B_u^0 \), Th. 4 produces various versions of the Trotter convergence theorem for various quadrature rules.

As everywhere else in this paper, in the following, it is understood that the potential \( V(x) \) is continuous and bounded from below.
Corollary 1 If \( \rho_n(x, x'; \beta) \) is defined by Eq. (5), then \( \rho_n(x, x'; \beta) \to \rho(x, x'; \beta) \), as \( n \to \infty \).

Proof. Follows from Th. 1 and the uniform convergence of the random series \( \sum_{k \geq 1} a_k \Lambda_k(u) \) to a Brownian bridge, as guaranteed by the Ito–Nisio theorem. \( \square \)

Corollary 2 If \( \rho_n(x, x'; \beta) \) is defined by Eq. (6), then \( \rho_{2^k-1}(x, x'; \beta) \to \rho(x, x'; \beta) \), as \( k \to \infty \).

Observation. The proof we construct for the subsequence \( n = 2^k - 1 \) is based on a special form of the Lie–Trotter product that is called the Lévy–Ciesielski form. The information from the paragraphs below is taken from Ref. [10]. In fact, for \( n = 2^k - 1 \), any Lie–Trotter product can be put in a Lévy–Ciesielski form, as shown in Ref. [14]. This form has important advantages when it comes to the Monte Carlo implementation, such as fast computation of paths and fast sampling. For this reason, we can restrict our attention to the subsequence \( n = 2^k - 1 \), without any loss of generality for actual applications.

Proof of the corollary. Let \( \{a_{l,j}; 1 \leq l \leq k, 1 \leq j \leq 2l - 1\} \) and \( \{b_{l,j}; 1 \leq l \leq \nu, 1 \leq j \leq 2^l - 1\} \) be two independent sets of i.i.d. standard normal variables. Let \( \{F_{l,j}(u); 1 \leq l \leq \nu, 1 \leq j \leq 2^l - 1\} \) be the system of Schauder functions on the interval \([0, 1]\). The Schauder functions are the primitives of the Haar \( L^2([0, 1]) \) wavelet bases and can be generated by contractions and translations as follows. Let \( F_{1,1}(u) : \mathbb{R} \to \mathbb{R} \) be defined by

\[
F_{1,1}(u) = \begin{cases} 
  u, & u \in [0, 1/2], \\
  1 - u, & u \in (1/2, 1], \\
  0, & \text{elsewhere}. 
\end{cases}
\] (35)

Then,

\[
F_{l,j}(u) = 2^{-(l-1)/2}F_{1,1}(2^l u - j + 1)
\] (36)

for all \( l \geq 1 \) and \( 1 \leq j \leq 2^l - 1 \). Extend the functions \( \{\tilde{\Lambda}_l(u); 1 \leq l \leq \nu\} \) outside the interval \([0, 1]\) by setting them to zero [the same way the first Schauder function \( F_{1,1}(u) \) was extended to the whole real axis in Eq. (35)] and define

\[
G_{l,j}(u) = 2^{-k/2}\tilde{\Lambda}_l(2^k u - j + 1)
\] (37)

for \( 1 \leq l \leq \nu \) and \( 1 \leq j \leq 2^k \).

Let \( d\omega_k(u) \) denote the discrete measure associated with the quadrature scheme specified by the \( nq2^k \) (not necessarily different) quadrature knots

\[
u_{i,j} = 2^{-k}(\theta_i + j - 1), \quad 1 \leq i \leq n_q, \quad 1 \leq j \leq 2^k
\] (38)

and the corresponding weights

\[
w_{i,j} = 2^{-k}w_i.
\] (39)

The new quadrature knots \( \nu'_{i,j} \) are obtained by contractions and translations of the original knots \( \theta_i \).
A problem of matching the covariance of a Brownian motion

With the convention that \(a_{l,2^{l-1}+1} = 0\) and \(b_{l,2^k+1} = 0\) for all \(l \in \mathbb{Z}^k\), we have

\[
\frac{\rho_n(x,x';\beta)}{\rho_{fp}(x,x';\beta)} = \int_{\mathbb{R}} da_{1,1} \ldots da_{k,2^k-1} (2\pi)^{-n/2} \exp\left(-\frac{1}{2} \sum_{l=1}^{k} \sum_{j=1}^{2^{l-1}} a_{l,j}^2\right) \times \\
\times \int_{\mathbb{R}} db_{1,1} \ldots db_{n,2^k} (2\pi)^{-(n+1)n_{\nu}/2} \exp\left(-\frac{1}{2} \sum_{l=1}^{n_{\nu}} \sum_{j=1}^{2^k} b_{l,j}^2\right) \times \\
\times \exp\left\{-\beta \int_{0}^{1} V\left[ x_r(u) + \sigma \sum_{l=1}^{k} a_{l,2^{l-1}+1} F_{l,2^{l-1}+1}(u) + \\
+ \sigma \sum_{l=1}^{n_{\nu}} b_{l,2^k+1} G_{l,2^k+1}(u) \right] \, d\omega_k(u)\right\},
\]

(40)

where \([2^{l-1}u]\) and \([2^k u]\) are the integer parts of \(2^{l-1}u\) and \(2^k u\), respectively.

Let us verify that the sequence of discrete measures \(d\omega_k(u)\) defines a quadrature scheme. If \(h : [0,1] \to \mathbb{R}\) is a continuous functions then, by its uniform continuity, for any \(\epsilon > 0\), there is \(\eta > 0\) such that \(|h(u) - h(\tau)| < \epsilon\) whenever \(|u - \tau| < \eta\).

Now, pick \(k\) large enough that \(1/2^k < \eta\). The Lebesgue integral of \(h\) over \([0,1]\) can be broken in \(2^k\) smaller parts over the intervals \([j-1)/2^k, j/2^k]\), for \(1 \leq j \leq 2^k\).

Since the length of these intervals is smaller than \(\eta\), we have both

\[
\left| \int_{(j-1)/2^k}^{j/2^k} h(u) \, du - 2^{-k}h(j/2^k) \right| < 2^{-k}\epsilon 
\]

(41)

and

\[
\left| \sum_{i=1}^{n_q} w_{i,j} h(u'_{i,j}) - 2^{-k}h(j/2^k) \right| < 2^{-k}\epsilon.
\]

(42)

The latter inequality is true because the points \(u'_{i,j}\) are in the interval \([(j-1)/2^k, j/2^k]\) and because

\[
\sum_{i=1}^{n_q} w_{i,j} = 2^{-k} \sum_{i=1}^{n_q} w_i = 2^{-k}.
\]

From Eqs. (41) and (42), we learn that

\[
\left| \int_{(j-1)/2^k}^{j/2^k} h(u) \, du - \sum_{i=1}^{n_q} w'_{i,j} h(u'_{i,j}) \right| < 2 \cdot 2^{-k}\epsilon
\]

(43)

and, by summing over all \(j\), we obtain

\[
\sum_{j=1}^{2^k} \left| \int_{(j-1)/2^k}^{j/2^k} h(u) \, du - \sum_{i=1}^{n_q} w'_{i,j} h(u'_{i,j}) \right| < 2\epsilon.
\]

(44)
Cristian Predescu

The left-hand side of Eq. (44) is clearly larger than
\[ \left| \sum_{j=1}^{2^k} \int_{(j-1)/2^k}^{j/2^k} h(u)du - \sum_{i=1}^{n_q} w'_{i,j} h\left(u_{i,j}\right) \right| = \left| \int_0^1 h(u)du - \int_0^1 h(u)d\omega_k(u) \right|. \]

Since \( \epsilon \) is arbitrary, it follows that
\[ \lim_{k \to \infty} \int_0^1 h(u)d\omega_k(u) = \int_0^1 h(u)du \]
and since \( h \) is arbitrary, it follows that the sequence of discrete measures \( d\omega_k(u) \) defines a quadrature scheme.

The first part of the random series appearing in Eq. (40) is the Lévy–Ciesielski series, which converges uniformly to a Brownian bridge. The corollary readily follows from the result in the preceding paragraph and Th. 1 if we prove that the tail series
\[ T_k(\bar{b}; u) = \sum_{l=1}^{n_\nu} b_{l,\lfloor 2^k u \rfloor + 1} G_{l,\lfloor 2^k u \rfloor + 1}(u) \] (45)
converges to zero uniformly almost surely, as \( k \to \infty \) (in other words, if our correction term does not ruin the uniform convergence of the Lévy–Ciesielski series). Let \( M > 0 \) be a common bound for the functions \( \{\tilde{A}_k(u); 1 \leq k \leq n_\nu\} \). Then, with the help of Eq. (37), we compute
\[ \left( \max_{0 \leq \varepsilon \leq 1} |T_k(\bar{b}; u)| \right)^4 \leq \frac{M^4}{2^{2k}} \max_{l,j} |b_{l,j}|^4 \leq \frac{M^4}{2^{2k}} \sum_{l=1}^{n_\nu} \sum_{j=1}^{2^k} b_{l,j}^4. \]
Taking the expected value, we get
\[ E\left( \max_{0 \leq \varepsilon \leq 1} |T_k(\bar{b}; u)| \right)^4 \leq \frac{M^4}{2^{2k}} (3n_\nu 2^k) = \frac{3n_\nu M^4}{2^k}. \]
Now, Chebyshev’s inequality produces
\[ P\left( \max_{0 \leq \varepsilon \leq 1} |T_k(\bar{b}; u)| \geq \epsilon \right) \leq \frac{1}{\epsilon^4} E\left( \max_{0 \leq \varepsilon \leq 1} |T_k(\bar{b}; u)| \right)^4 \leq \frac{3n_\nu M^4}{\epsilon^4 2^{2k}} \]
and so,
\[ \sum_{k=1}^{\infty} P\left( \max_{0 \leq \varepsilon \leq 1} |T_k(\bar{b}; u)| \geq \epsilon \right) \leq \frac{3n_\nu M^4}{\epsilon^4} \sum_{k=1}^{\infty} \frac{1}{2^k} = \frac{3n_\nu M^4}{\epsilon^4} < \infty. \]
The first Borel–Cantelli lemma implies that
\[ P\left( \max_{0 \leq \varepsilon \leq 1} |T_k(\bar{b}; u)| \geq \epsilon \text{ i.o.} \right) = 0, \]
A problem of matching the covariance of a Brownian motion

which means that, with probability one, there is a rank $K \geq 1$ such that

$$\max_{0 \leq u \leq 1} |T_k(\bar{b}; u)| < \epsilon, \quad \forall k \geq K.$$ 

Letting $\epsilon$ go to zero through the countable sequence $\epsilon_j = 1/j$, we obtain the almost sure uniform convergence to zero of the tail series $T_k(\bar{b}; u)$. The proof of the corollary is concluded. □

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