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Anomalous paths in quantum mechanical path-integrals

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ABSTRACT

We investigate modifications of the discrete-time lattice action, for a quantum mechanical particle in one spatial dimension, that vanish in the naïve continuum limit but which, nevertheless, induce nontrivial effects due to quantum fluctuations. These effects are seen to modify the geometry of the paths contributing to the path-integral describing the time evolution of the particle, which we investigate through numerical simulations. In particular, we demonstrate the existence of a modified lattice action resulting in paths with any fractal dimension, d_f , between one and two. We argue that $d_f = 2$ is a critical value, and we exhibit a type of lattice modification where the fluctuations in the position of the particle becomes independent of the time step, in which case the paths are interpreted as superdiffusive Lévy flights. We also consider the jaggedness of the paths, and show that this gives an independent classification of lattice theories.

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1. Introduction

The path-integral representation of the amplitude $\langle x', t' | x, t \rangle$ for a quantum mechanical particle of mass *m* moving in a local potential *V*(*x*) is usually written as a limit of a multi-dimensional integral [1]:

$$Z \equiv \langle x', t' | x, t \rangle = \lim_{N \to \infty} \mathcal{N} \int dx_1 \dots dx_{N-1} e^{-S_N},$$
(1)

where we have changed to imaginary time $(t \rightarrow -it)$ and set $\hbar = 1$. Here $\mathcal{N} = (m/2\pi a)^{N/2}$ and S_N is the discrete-time action which should approach the classical continuum action *S* as the lattice constant $a \equiv (t_f - t_i)/N$ goes to zero, *i.e.*,

$$\lim_{N \to \infty} S_N = S = \int_{t_i}^{t_f} dt \left[\frac{1}{2} \dot{x}^2 + V(x) \right].$$
 (2)

We have chosen units such that the mass m of the particle is one. The particular choice

$$S_N \equiv \sum_{k=0}^{N-1} S_k = \sum_{k=0}^{N-1} a \left[\frac{1}{2} \left(\frac{\Delta x_k}{a} \right)^2 + V(x_k) \right],$$
(3)

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where $\Delta x_k \equiv x_{k+1} - x_k$, with a time-step $dt \rightarrow \Delta t \equiv a$, is referred to as the naïve discretization of the classical action *S*, and has, for example, been used in modeling time as a discrete and dynamical variable [2]. The choice of Eq. (3) is, however, by no means unique and the ambiguity of the discretization has been investigated previously by, *e.g.*, Klauder et al. in Ref. [3]. As an interesting example, it has also been shown that adding terms proportional to $a\Delta x_k^{2n}$, as $a \rightarrow 0$ (n = 1, 2, ...), to each term S_k in the sum in Eq. (3) permits a radical speedup of the convergence in Monte Carlo simulations [5]. Classically, one expects $\Delta x_k/a \rightarrow \dot{x}$ to be well-defined as $a \rightarrow 0$ and thus $S_k = \mathcal{O}(a)$, and, as was noted in Ref. [5], one would have $a\Delta x_k^{2n} \rightarrow a^{2n+1}\dot{x}^{2n} = \mathcal{O}(a^{2n+1})$, which clearly vanish in the $a \rightarrow 0$ limit. We will refer to these considerations as the "naïve continuum limit" in the following.

As was pointed out in Ref. [3], and in a related framework in Ref. [4], the argumentation above is, however, not true for quantum mechanical paths, as one expects $\Delta x_k = \mathcal{O}(\sqrt{a})$ in accordance with the Itô calculus for a Wiener process, and thus the action then contains terms S_k of order one. Modifications as those considered in Ref. [5] still vanish, but only as fast as $\mathcal{O}(a^{n+1})$. This implies no difficulty for the numerical speedup procedure, but in general, it is clear that one must take care when modifying the action in the presence of quantum fluctuations.

We now wish to expand on the work from Ref. [3] and proceed to study precisely those modifications to the discrete action that vanish in the naïve limit, but might induce non-trivial effects when quantum fluctuations are taken into account. We will





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show that not only can non-vanishing local potentials be induced by such alterations, as was shown in Ref. [3], but the situation is further complicated in that the size of quantum fluctuations can be changed under the modified lattice theory, such that no naïve assumptions on the continuum limit can be made. This can be seen to manifest itself in the geometrical properties of the paths contributing to the path-integral, and as we will see shortly, can generate both sub- and super-diffusive behaviour.

2. Geometry of path-integral trajectories

We now quickly review two useful measures that will be used to quantify the geometry of relevant paths in the path-integral. The geometry of path-integral trajectories has been investigated previously, in particular by Kröger et al. in Ref. [6], where a fractal dimension was defined and found both analytically and numerically for local and velocity-dependent potentials. More recently a complementary property termed "jaggedness" was identified by Bogojevic et al. in Ref. [7]. Both of these measures signify the relevance of different paths as to what degree they contribute to the total path-integral.

To define the fractal dimension, d_f , for path-integral trajectories, we recall that the fractal dimension for a classical path can be defined in the following way: We first define a length of the path, $L(\epsilon)$, as obtained with some fundamental resolution ϵ . This can, for example, be done by making use of a minimal covering of the path with "balls" of diameter ϵ such that $L(\epsilon) = N(\epsilon) \times \epsilon$, where $N(\epsilon)$ is the number of balls. A fractal dimension can then be defined as the unique number d_f such that $L(\epsilon) \sim \epsilon^{1-d_f}$ as $\epsilon \rightarrow 0$ [12]. For path-integral trajectories, a total length can be defined as $\langle L \rangle = \langle \sum_k |\Delta x_k| \rangle$, and the role of ϵ will be played by the expected absolute change in position, $\langle |\Delta x_k| \rangle$, over one small time step $\Delta t \simeq a$. Here $\langle \cdot \rangle$ denotes the quantum-mechanical average using the probability distribution obtained from Eq. (1). For a typical value, $|\Delta x|$, of $\langle |\Delta x_k| \rangle$, say $|\Delta x| \simeq (\Delta t)^{1/\gamma}$, we then have that $\langle L \rangle \simeq N |\Delta x| \simeq T |\Delta x|^{1-\gamma}$ since $N \simeq T / \Delta t$, with $T = t_f - t_i$. We then conclude that $d_f = \gamma$. The fractal dimension can therefore be obtained through a scaling with the number of lattice sites N, as $N \to \infty$, with $T = N \Delta t \simeq Na$ held fixed, *i.e.*,

$$\langle L \rangle \sim N^{1-1/d_f},\tag{4}$$

for sufficiently large *N*. This is also the definition made use of in Ref. [6], and is a measure of how the increments $\langle |\Delta x_k| \rangle$ scale with the time step $\Delta t \simeq a$. In the spirit of anomalous-diffusion considerations (see *e.g.* Ref. [13]), we will refer to those paths with a fractal dimension $d_f < 2$, as defined above, as sub-diffusive, reflecting that they spread in space at a slower than normal rate. Similarly those paths with $d_f > 2$ are referred to as super-diffusive, which then corresponds to Lévy flights (see *e.g.* Ref. [14]).

A remark on the physical interpretation of d_f is in order before we proceed. The length $\langle L \rangle$ defined above is not necessarily an experimentally observable length. It gives us, however, an insight into the nature of how the geometry of those paths with a nonzero measure change under modification of the lattice action. The definition of a fractal dimension for the *physical* path of a quantum mechanical particle must necessarily involve considerations of a measuring apparatus, as was done by Abbott and Wise [8]. Inclusion of quantum measurements in a path-integral framework has been discussed in the literature (see *e.g.* Ref. [9]), but will not be considered in this work.

It is well known that the paths contributing to the pathintegral, Eq. (1), are continuous but non-differentiable. Indeed, using a partial integration, Feynman and Hibbs [1] showed that for any observable F the identity

$$\left(\frac{\delta F}{\delta x_k}\right) = \left(F\frac{\delta S}{\delta x_k}\right) \tag{5}$$

holds. In the case $F = x_k$ this leads to

$$\left\langle \Delta x_k^2 \right\rangle = \mathcal{O}(a),\tag{6}$$

for the lattice action Eq. (3) and for sufficiently small *a*, and where we from now on assume that expressions like $\langle x_k \, dV(x_k)/dx_k \rangle$ are finite. Hence, we expect $\langle |\Delta x_k| \rangle \propto 1/\sqrt{N}$ and $\langle L \rangle \propto \sqrt{N}$ corresponding to a fractal dimension of $d_f = 2$, which has been confirmed numerically in Ref. [6].

The second measure we will use to describe the relevant paths in the path-integral, is the "jaggedness", *J*, defined in Ref. [7], which counts the number of maxima and minima of a given path:

$$J = \frac{1}{N-1} \sum_{k=0}^{N-2} \frac{1}{2} \Big[1 - \langle \operatorname{sgn}(\Delta x_k \Delta x_{k+1}) \rangle \Big], \tag{7}$$

with $J \in [0, 1]$. It is a measure of the correlation between Δx_k and Δx_{k+1} with J = 1/2 + O(a) for completely uncorrelated increments. We therefore expect the jaggedness to be invariant under modifications only altering nearest neighbor interactions on the lattice. Below we will consider the average value of J for sub- and super-diffusive paths.

3. Sub-diffusive paths

Sub-diffusive paths, as defined here, were discovered to be the contributing paths in the presence of a velocity dependent potential, $V_0|v|^{\alpha}$, in Ref. [6]. We will here consider a similar modification, that in fact *vanish* in the naïve continuum limit, yet changes the geometry of the paths when quantum fluctuations are taken into account:

$$S_k \to S_k + ga^{\xi} \left| \frac{\Delta x_k}{a} \right|^{lpha},$$
 (8)

where g is a coupling constant, $\xi \ge 1$ and $\alpha \ge 0$. The last term is identical to the modification considered in Ref. [6] for $\xi = 1$, but naïvely vanishes for any $\xi > 1$. Due to quantum fluctuations, however, Eq. (6) must be replaced by

$$\frac{1}{a} \langle \Delta x_k^2 \rangle + g \alpha a^{\xi - \alpha} \langle |\Delta x_k|^{\alpha} \rangle = \mathcal{O}(1), \tag{9}$$

showing that for $\alpha > 2\xi$ the last term dominates, and we expect $\langle |\Delta x_k| \rangle \propto a^{(\alpha-\xi)/\alpha}$, corresponding to a fractal dimension of $d_f = \alpha/(\alpha - \xi)$. For $\alpha \leq 2\xi$ we still have $d_f = 2$ showing that 2ξ is a critical point for the fractal dimension as a function of α . For $\xi = 1$ this reproduces the results from [6]. In Fig. 1 we show how the length $\langle L \rangle$ scales with the number of lattice sites *N* for various α and $\xi = 2$. The results are produced numerically by standard Monte Carlo methods [6,10,11]. From this scaling one can find the fractal dimension according to Eq. (4). In Fig. 2 we have extracted the fractal dimension as a function of α numerically for $\xi = 1$, 2 and 3. We see that the numerical results fit well to the expected values of $d_f = 2$ for $\alpha \leq 2\xi$ and $d_f = \alpha/(\alpha - \xi)$ for $\alpha > 2\xi$, shown as solid lines in the figure.

4. Super-diffusive paths

Consider now modifications of the form

$$S_k \to f(S_k),\tag{10}$$

for some analytical function, f(x), with the constraint f(x) = x as $x \to 0$, in order to reproduce the classical limit. This constitutes



Fig. 1. Scaling of the average length $\langle L \rangle$ as a function of lattice sites, *N*, for the lattice action given in Eq. (8), with $\xi = 2$ and various α . $\log(\langle L \rangle)$ was fitted to $\beta \log(N) + b$. The values for α are 3, 4, 5 and 6, starting from the top line and descending. Statistical error bars are not visible in the figure.



Fig. 2. The fractal dimension, d_f , as a function of α , for $\xi = 1, 2$ and 3, for the action defined in Eq. (8). The dots are numerical results, and the solid lines represent the expected theoretical values according to $d_f = \alpha/(\alpha - \xi)$, with $\xi = 3$ (the top line), $\xi = 2$ (the middle curve), and $\xi = 1$ (the lowest curve).

a large class of local modifications—*i.e.*, only influencing nearestneighbor couplings on the lattice—that have the same naïve $a \rightarrow 0$ limit.

We will also assume, if required, that there exists some large distance infrared cutoff so that the integral

$$\psi(x_{k+1}) = N \int dx_k \, e^{-f(S_k)} \psi(x_k), \tag{11}$$

exists, describing the evolution of the wave-function $\psi(x)$ over a small time step, a, under the modified lattice action. For reasons of simplicity, we assume the infrared regularization $\psi(x_k) = 0$ for $|x_{i+k} - x_k| \ge L$, for some sufficiently large *L*.

Through a straightforward renormalization procedure (see Appendix A) we are able to write down an effective action that is equivalent to the modification, Eq. (10), in the continuum limit. Remarkably, the so obtained effective action can formally be written in the naïve form of Eq. (3), *i.e.*

$$S_N \to S_N^{\text{eff}} = a \sum_{k=0}^{N-1} \left[\frac{1}{2s[f]} \left(\frac{\Delta x_k}{a} \right)^2 + g[f] V(x_k) \right],$$
 (12)

where

$$s[f] = \frac{\int_{\Omega} dy \, y^2 \exp(-f(\frac{y^2}{2}))}{\int_{\Omega} dy \exp(-f(\frac{y^2}{2}))},$$
(13)

and

$$g[f] = \frac{\int_{\Omega} dy \, f'(\frac{y^2}{2}) \exp(-f(\frac{y^2}{2}))}{\int_{\Omega} dy \exp(-f(\frac{y^2}{2}))}.$$
 (14)

Here the integrals run over the domain Ω , given by $-L/\sqrt{a} < y < L/\sqrt{a}$. Hence the integrals become independent of the infrared cutoff, *L*, in the $a \rightarrow 0$ limit. As discussed in Appendix A, this can be interpreted as a renormalization of the particle's mass and potential, and will in general be finite or infinite in the limit $a \rightarrow 0$, depending on the form of *f*. With the modification of Eq. (10), Eq. (6) must, however, be replaced by

$$\left\langle f'(S_k)\Delta x_k^2\right\rangle = \mathcal{O}(a),\tag{15}$$

potentially changing how $\langle |\Delta x_k| \rangle$ scales with *a* and thus the fractal dimension d_f as defined above. Similarly, for the equivalent effective counterpart of Eq. (12), we see that the scaling can be written

$$\Delta x_k^2 = s[f]a, \tag{16}$$

and therefore all modifications to the short-time scaling are contained in the functional s[f]. For a function f(x) that is bounded, however, s[f] diverges like L^2/a as a goes to zero, and therefore $\langle \Delta x_k^2 \rangle \simeq L^2$ in terms of the infrared cutoff L. In this case we therefore expect that the particle can make arbitrarily large jumps, independent of a. This behavior can be interpreted, at least formally, as an infinite fractal dimension for the particle's path since $\langle \Delta x_k^2 \rangle \simeq a^{2/d_f}$, and is typical for *any* such f. Such paths are analogous to Poisson paths, such as appear in Ref. [15], which involve paths with continuous segments joined by jumps whose magnitude is drawn from a well defined distribution at time intervals, again, with a suitable distribution.

We illustrate these features in terms of the following family of lattice modifications, defined through Eq. (10),

$$f \equiv f_{\gamma}(x) = \begin{cases} (1+x)^{\gamma} & \gamma > 0\\ -(1+x)^{\gamma} & \gamma < 0. \end{cases}$$
(17)

Here $f_{\gamma}(x) \simeq 1 + \gamma x$ as $x \to 0$ (the scaling factor γ and constant term is irrelevant for our discussion). As γ approaches zero from above, $s[f_{\gamma}]$ becomes larger, and is infinite in the limit $\gamma \to 0$. Since $s[f_{\gamma}]$ implies a rescaling of a, as can be seen in Eq. (16), the exceedingly large values of $s[f_{\gamma}]$ for small γ means we need a correspondingly large number of lattice sites to approach the continuum limit. In any case, as long as $s[f_{\gamma}]$ implies a finite rescaling, we expect the fractal dimension to be invariant. For $\gamma < 0$, however, the integral $s[f_{\gamma}]$ does not exist as a approaches zero.

In Fig. 3 we show example paths for a free particle, V(x) = 0, and four different γ , generated by standard Monte Carlo methods. The paths exhibit larger jumps for smaller γ . For $\gamma = -1$ the path has a radically different geometry. In Fig. 4 the length $\langle L \rangle$ is plotted for varying number of lattice sites, *N*, for the same values of γ . The scaling $\langle L \rangle \propto N^{\beta}$ was found to be $\beta = 0.499 \pm 0.001$, $\beta = 0.495 \pm 0.004$ and $\beta = 0.997 \pm 0.003$ for the respective cases $\gamma = 2$, $\gamma = 1$, $\gamma = 0.5$ and $\gamma = -1$ (the errors are mean square errors from the linear regression). The corresponding fractal dimensions, as defined in Eq. (4), are consistent with $d_f = 2$ for the $\gamma > 0$ cases and $d_f = \infty$ for $\gamma = -1$.

The behavior for negative γ is in fact typical for *any* modification of the form Eq. (10) with a bounded f(x), and f(x) = x as $x \to 0$. In Fig. 4 we also include results for the modifications $f(x) = \tanh(x)$ and $f(x) = \sin(x)$. The scaling was found to be $\beta = 1.006 \pm 0.011$ and $\beta = 0.982 \pm 0.007$ respectively, corresponding to an infinite fractal dimension in both cases.

In Fig. 5 we show how β scales with γ for the modifications in Eq. (17). As γ becomes small and positive, there are numerical difficulties due to the necessity of a large number of lattice sites. We here show results for positive γ no smaller than $\gamma = 0.3$. The results are consistent with $\beta = 1$ and $d_f = \infty$ for $\gamma < 0$ and $\beta =$ 0.5, and $d_f = 2$ for $\gamma > 0$, and points towards critical behaviour at $\gamma = 0$, in the limit $N \to \infty$.



Fig. 3. Example paths for the lattice modification defined through Eq. (17), for different γ , showing the various behaviour. Top left is for $\gamma = 2.0$, top right for $\gamma = 1.0$, bottom left for $\gamma = 0.5$ and bottom right for $\gamma = -1$, using dimensionless units.



Fig. 4. Scaling of the average length $\langle L \rangle$ as a function of lattice sites *N*. $\log(\langle L \rangle)$ was fitted to $\beta \log(N) + b$. The $\gamma = -1$ and "tanh" action coincide at the top line, the second line is for the "sin" action, the third for $\gamma = 0.5$, the forth for $\gamma = 1.0$ and the fifth for $\gamma = 2.0$.

We have also calculated the jaggedness for sub- and superdiffusive actions. In the sub-diffusive case, *i.e.* actions of the form given in Eq. (8), we find results consistent with J = 1/2 as expected, since there are no correlations between increments Δx_k and Δx_{k+1} introduced through the modification. This highlights the fact that a classification in terms of jaggedness is independent of a classification in terms of fractal dimension, as was stressed in Ref. [7]. Indeed, even when the paths have a fractal dimension close to one, they are not at all smooth and still fall in to the same jaggedness class, with J = 1/2.

For the super-diffusive case there is, however, a subtlety involved in that the particle will always be subject to the infrared boundary effects. In practice, for a finite number of Monte Carlo samples stored on a computer, the particle's position is always confined to some interval for all times, say $-L/2 < x_k < L/2$. If the probability density for the particle's position at time $t_k \simeq ka$, $p(x_k) = |\psi(x_k)|^2$, becomes independent of the position at prior



Fig. 5. $\beta = (d_f - 1)/d_f$ as a function of γ for the modification Eq. (17). For negative γ , $\beta = 1.0$ corresponds to $d_f = \infty$, and for positive γ , $\beta = 0.5$ to $d_f = 2$.



Fig. 6. Typical distributions for the jaggedness with N = 512 and a = 1/N. The leftmost Gaussian is centered at 0.5 with width 0.022, and well approximates the case of the naïve action Eq. (3), and the action given in Eq. (8) with $\xi = 1$ and $\alpha = 10$, for which the numerical results represented as dots are nearly indistinguishable. The middle Gaussian is centered at 0.625, and the coinciding dots are numerical results for the action given through Eq. (17), with $\gamma = -1$ and the particle's position restrained to a box of width one. The rightmost Gaussian is centered at 0.667, and the dots are numerical results for a uniform distribution of the particles position at each time step.

times, such as is the case for the super-diffusive paths considered here, the conditional probability $p_{\text{peak}}(x_k)$ for a "peak" at x_k , where a "peak" is defined as a point x_k such that Δx_{k-1} and Δx_k have opposite signs, is just

$$p_{\text{peak}}(x_k) \equiv P((x_{k-1} < x_k \text{ and } x_{k+1} < x_k))$$

or $(x_{k-1} > x_k \text{ and } x_{k+1} > x_k)$, (18)

that is,

$$p_{\text{peak}}(x_k) = P(x_{k-1} < x_k)P(x_{k+1} < x_k) + P(x_{k-1} > x_k)P(x_{k+1} > x_k).$$
(19)

Consider now, as an example, the case of a uniform distribution on the interval, *i.e.* $p(x_k) = 1/L$ for $-L/2 < x_k < L/2$, and zero otherwise. One easily finds that $p_{\text{peak}}(x_k) = (1/2 + x_k/L)^2 + (1/2 - x_k/L)^2$. One might think that for large *L* the probability for a peak should be close to 1/2, but since there is no restriction on the particle's position, it can be close to the boundary for *any L*. The expected number of peaks then becomes $\int_{-L/2}^{L/2} p_{\text{peak}}(x_k) p(x_k) dx_k = 2/3 \simeq 0.667$, which is, of course, precisely the jaggedness. For super-diffusive actions we find, in numerical simulations, that the jaggedness takes values in between the value for the naïve action and the value for a uniform distribution as just discussed. In Fig. 6 we show some typical example distributions p(J) of the jaggedness. We compare the sub-diffusive case with the usual naïve action, and find that the distributions are nearly indistinguishable and very well approximated by a Gaussian centered at 0.5. We also show a distribution for a super-diffusive action, and for a uniform distribution of the particle's position at each time step, for comparison.

5. Conclusions

To conclude, we have shown that lattice actions, that approach the classical action in the naïve continuum limit, can display highly anomalous behaviour when quantum fluctuations are taken into account. Not only can non-vanishing local potentials be induced by such lattice modifications, as was shown by Klauder et al. in Ref. [3], but non-local effects can appear in that the geometry of the paths is changed. We have demonstrated modified lattice theories where the paths in the path-integral, with measure greater than zero, exhibit both sub-diffusive and super-diffusive behaviour. We find it noticeable that under certain assumptions, a large class of modified actions can, through a renormalization procedure, always be written formally on the naïve discretized form. Alternative views on the notion of fractal dimensions in quantum physics has been discussed in the literature as in, e.g., Ref. [16] which, however, is closely related to the notion of fractional derivatives [13] and therefore different from the local deformations of the lattice actions as consider in the present Letter. Finally, we remark, as was argued already in Ref. [3], that the lattice corrections considered in the present Letter do not effect the physics of the continuum limit in the field-theoretical case, at least not for asymptotically free gauge field theories.

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Appendix A. Renormalizations induced by modified lattice actions

For the convenience of the readers we give a derivation of the Schrödinger equation for the modified mechanics defined in Eq. (10). Consider the evolution of the wave function over a small time step a:

$$\psi_{k+1}(x_{k+1}) = N \int_{-L+x_{k+1}}^{L+x_{k+1}} dx_k \exp(-f(S_{k+1,k}))\psi_k(x_k), \quad (A.1)$$

where *N* is a normalization constant and we have restricted the particles movement to the interval $-L < x_{i+1} - x_i < L$ to ensure the integral always is finite. Introducing the variables *x* and *y* through $x = x_{k+1}$ and $x_k = x + \sqrt{ay}$, and by Taylor expanding a sufficiently smooth potential $V(x_k)$, $f(x_k)$ and $\psi_k(x_k)$, dropping terms of order $O(a^2)$, we obtain

 $\psi_{k+1}(x)$

$$= N\sqrt{a} \int_{\Omega} dy \exp\left\{-f\left(\frac{y^2}{2}\right)\right\} \psi_k(x)$$

+ $\frac{1}{2}N\sqrt{aa} \int_{\Omega} dy y^2 \exp\left\{-f\left(\frac{y^2}{2}\right)\right\} \psi_k''(x)$
- $N\sqrt{aa}V(x) \int_{\Omega} dy f'\left(\frac{y^2}{2}\right) \exp\left\{-f\left(\frac{y^2}{2}\right)\right\} \psi_k(x),$ (A.2)

with a domain of integration Ω as given by $-L/\sqrt{a} < y < L/\sqrt{a}$. We now choose the normalization constant such that

$$N\sqrt{a} \int_{\Omega} \mathrm{d}y \, \exp\left\{-f\left(\frac{y^2}{2}\right)\right\} = 1. \tag{A.3}$$

Then

$$\psi_{k+1}(x) = \psi_k(x) + \frac{a}{2}s[f]\psi_k''(x) - aV(x)g[f]\psi_k(x),$$
(A.4)

where s[f] and g[f] are given in Eqs. (13) and (14). We now obtain the following imaginary-time Schrödinger equation

$$\lim_{a \to 0} \frac{\psi_{k+1}(x) - \psi_k(x)}{a} = \frac{\partial \psi(x, t)}{\partial t}$$
$$= \frac{1}{2} s[f] \psi''(x, t) - g[f] V(x) \psi(x, t), \quad (A.5)$$

i.e.,

$$\frac{\partial \psi(\mathbf{x},t)}{\partial t} = \frac{s[f]}{2} \psi''(\mathbf{x},t) - g[f] V(\mathbf{x}) \psi(\mathbf{x},t).$$
(A.6)

Introducing a mass m and \hbar again, we see that s[f] and g[f] constitutes a renormalization of the mass and potential respectively. One can also use this wave equation to show that the imaginary-time commutation relation $[x, p] = \hbar$ still holds in the discretized theory when we use $m_R \dot{x}$ for the momentum and the bare mass m has been replaced by the renormalized mass $m_R = m/s[f]$ (see Sections 7–5 in Ref. [1]). Since \hbar is unrenormalized we can make use of units such that $\hbar = 1$.

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