

First-arrival-time distributions for a Dirac electron in 1+1 dimensions

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For the special case of freely evolving Dirac electrons in (1+1) dimensions, Feynman checkerboard paths have previously been used to derive Wigner's arrival-time distribution which includes all arrivals. Here, an attempt is made to use these paths to determine the corresponding distribution of first-arrival times. Simple analytic expressions are obtained for the relevant components of the first-arrival propagator. These are used to investigate the relative importance of the first-arrival contribution to the Wigner arrival-time distribution and of the contribution arising from interference between first and later (i.e., second, third, etc.) arrivals. It is found that a distribution of (intrinsic) first-arrival times for a Dirac electron cannot in general be consistently defined using checkerboard paths, not even approximately in the nonrelativistic regime.

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I. INTRODUCTION

In the past decade there has been considerable interest in deriving and understanding arrival-time distributions for quantum particles, using a wide variety of approaches [for recent reviews, see, Refs. [1–3]]. Here we focus on an approach based on Feynman paths [4].

Yamada and Takagi [5,6] applied the consistent histories approach, with Feynman paths as particle histories, to the problem of deriving an intrinsic arrival-time distribution. They considered the special case of a freely evolving nonrelativistic quantum particle in one spatial dimension. In Ref. [5] they showed that within their approach one cannot classify the histories according to the number of times n that a particle arrives at a given spatial point $x=X$ during a specified finite time interval because the amplitude for n arrivals is zero for every finite n . Their qualitative explanation was that a typical Feynman path, being nondifferentiable in time, intersects X an infinite number of times in the given time interval, leaving zero amplitude for any finite value of n . They further suggested that this would also be the case for a quantum particle propagating in three spatial dimensions in the presence of a potential.

The velocity associated with a typical Feynman path for a nonrelativistic electron is infinite at almost every point on it. This is not the case for a relativistic electron—the velocity associated with a Feynman path [4,7] for a relativistic electron is finite at almost every point on it. Hence, such a path will not intersect X an infinite number of times in a given finite time interval and the amplitude for n arrivals need not be zero for every finite n . This is the primary motivation for the following attempt to derive the first-arrival-time distribution for a freely evolving Dirac electron in (1+1) dimensions. In Sec. II, a checkerboard path derivation [8] of Wigner's arrival-time distribution [9] which includes all arrivals is

sketched, primarily to introduce the basic notation and concepts used in the following sections. In Sec. III the first-arrival propagator for this special case is derived. Unfortunately, there is interference between the $n=1$ and $n>1$ contributions to the arrival-time distribution so that the distribution of intrinsic first-arrival times cannot be consistently defined using this approach. Calculated results for the $n=1$ and interference contributions are presented in Sec. IV for two simple cases. Concluding remarks are made in Sec. V.

II. CHECKERBOARD PATHS AND WIGNER'S ARRIVAL-TIME DISTRIBUTION FOR DIRAC ELECTRONS

Consider the (1+1)-dimensional free-electron Dirac equation [10] in the form

$$i\hbar \hat{1} \frac{\partial}{\partial t} \Psi(x,t) = \hat{H} \Psi(x,t) = -i\hbar c \hat{\sigma}_z \frac{\partial}{\partial x} \Psi(x,t) - mc^2 \hat{\sigma}_x \Psi(x,t) \quad (1)$$

with $\Psi(x,t) \equiv [\Psi_+(x,t), \Psi_-(x,t)]^T$ a two-component spinor and

$$\hat{\sigma}_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \hat{\sigma}_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}. \quad (2)$$

The velocity operator

$$\hat{v} \equiv \frac{d\hat{x}}{dt} = \frac{i}{\hbar} [\hat{H}, \hat{x}] = c \hat{\sigma}_z \quad (3)$$

has the orthonormal eigenfunctions

$$\chi_+ = \begin{pmatrix} 1 \\ 0 \end{pmatrix}, \quad \chi_- = \begin{pmatrix} 0 \\ 1 \end{pmatrix} \quad (4)$$

with eigenvalues $+c$ and $-c$, respectively. Hence,

$$\Psi(x,t) = \begin{pmatrix} \Psi_+(x,t) \\ \Psi_-(x,t) \end{pmatrix} = \Psi_+(x,t)\chi_+ + \Psi_-(x,t)\chi_- \quad (5)$$

and the upper and lower components of the wave function $\Psi(x,t)$ are, respectively, the amplitudes at (x,t) for the right-going ($+c$) and left-going ($-c$) velocity eigenstates χ_+ and χ_- . The four components of the retarded 2×2 propagator $K(x_B, t_B; x_A, t_A) \equiv K(B; A)$ of the $(1+1)$ -dimensional free-electron Dirac equation (1) are labeled by velocity directions ($+$ or $-$) at (x_A, t_A) and at (x_B, t_B) and are accordingly defined by

$$\begin{pmatrix} \Psi_+(x_B, t_B) \\ \Psi_-(x_B, t_B) \end{pmatrix} = \int_{-\infty}^{\infty} dx_A \begin{pmatrix} K_{++}(x_B, t_B; x_A, t_A) & K_{+-}(x_B, t_B; x_A, t_A) \\ K_{-+}(x_B, t_B; x_A, t_A) & K_{--}(x_B, t_B; x_A, t_A) \end{pmatrix} \begin{pmatrix} \Psi_+(x_A, t_A) \\ \Psi_-(x_A, t_A) \end{pmatrix}.$$

The left subscript, $+$ or $-$, on K denotes, respectively, a right-going or left-going arrival at x_B at time t_B while the right subscript $+$ or $-$ denotes, respectively a right-going or left-going departure from x_A at time t_A .

Subtracting the Hermitean conjugate of Eq. (1), multiplied from the right by $\Psi(x,t)$, from Eq. (1) multiplied from the left by the Hermitean conjugate of $\Psi(x,t)$ gives the continuity equation

$$\frac{\partial}{\partial t} \Psi^\dagger(x,t) \Psi(x,t) + \frac{\partial}{\partial x} \Psi^\dagger(x,t) c \hat{\sigma}_z \Psi(x,t) = 0. \quad (6)$$

It is assumed throughout the paper that the parameters of the initial wave function are such that Dirac's original identification of

$$\rho(x,t) = \Psi^\dagger(x,t) \Psi(x,t) = |\Psi_+(x,t)|^2 + |\Psi_-(x,t)|^2 \quad (7)$$

and

$$J(x,t) = c \Psi^\dagger(x,t) \hat{\sigma}_z \Psi(x,t) = c [|\Psi_+(x,t)|^2 - |\Psi_-(x,t)|^2] \quad (8)$$

with single-electron probability and probability current densities, respectively, is an adequate approximation.

In Feynman and Hibb's classic book "Quantum Mechanics and Path Integrals" [4] it is stated that the free-electron propagator $K(x_B, t_B; x_A, t_A)$ can be constructed from a model in which a particle going from x_A at time t_A to x_B at time t_B is constrained to move diagonally in space time at constant speed c in checker fashion (i.e., forward in time with spatial increment $\pm \Delta x$ with $\Delta x = c \Delta t = c(t_B - t_A)/N$ for each of N equal time steps $\Delta t > 0$) [11]. Each component of the propagator is obtained as the $N \rightarrow \infty$ limit of the sum over all N -step checkerboard paths joining (x_A, t_A) to (x_B, t_B) , with the first and last steps appropriately fixed, when the weight associated with a path having R (noncompulsary) reversals of direction or corners is taken to be $(imc^2 \Delta t / \hbar)^R$. Jacobson and Schulman [7] regrouped the sum-over-paths into a sum-over- R , i.e.,

$$K_{\beta\alpha}(B; A) = i(mc/2\hbar) \lim_{N \rightarrow \infty} \sum_{R \geq 0} \Phi_{\beta\alpha}(R) \times (imc^2 \Delta t / \hbar)^R \quad (\alpha = \pm; \beta = \pm), \quad (9)$$

where $\Phi_{\beta\alpha}(R)$ is the number of $\beta\alpha$ paths with R noncompulsary reversals. They also evaluated the four checkerboard path integrals, obtaining the following closed-form expressions for the components of the propagator:

$$K_{++}(B; A) = -\frac{ct_{BA} + x_{BA}}{2\lambda_c l_{BA}} J_1 \left(\frac{l_{BA}}{\lambda_c} \right), \quad (10)$$

$$K_{--}(B; A) = -\frac{ct_{BA} - x_{BA}}{2\lambda_c l_{BA}} J_1 \left(\frac{l_{BA}}{\lambda_c} \right), \quad (11)$$

$$K_{+-}(B; A) = K_{-+}(B; A) = \frac{i}{2\lambda_c} J_0 \left(\frac{l_{BA}}{\lambda_c} \right), \quad (12)$$

where $\lambda_c \equiv \hbar/mc$ is the Compton wavelength of the electron, $x_{BA} \equiv (x_B - x_A)$, $t_{BA} \equiv (t_B - t_A)$, and $l_{BA} \equiv c\tau_{BA}$ with $\tau_{BA} \equiv t_{BA} [1 - (v_{BA}/c)^2]^{1/2}$ the proper time for a particle moving with constant velocity $v_{BA} \equiv x_{BA}/t_{BA}$. Jacobson and Schulman also determined the number of reversals, R_0 , that gives the maximum contribution to the sum in Eq. (9). They obtained $R_0 = l_{BA}/\lambda_c$ and also showed that the sum is dominated by terms having R within $\approx R_0^{1/2}$ of R_0 [12]. The picture that emerges [7,13,14] is that in which the particle always moves with speed c and typically travels a distance $\approx (t_{BA}/\tau_{BA})\lambda_c \gg \lambda_c$ between reversals of direction; its motion is Brownian with diffusion constant \hbar/m only on scales much larger than this correlation distance.

Now consider the problem of deriving an expression for the distribution $\Pi(T; X)$ of arrival times T at the spatial point $x = X$ for an ensemble of Dirac electrons all prepared in the same initial state $\Psi(x, 0)$. Following Yamada and Takagi, it is *assumed* that the arrival-time distribution for the fictitious particles of the checkerboard model (with $t_A = 0$, $x_B = X$, and $t_B = T$), should it be well defined, can be identified with the desired distribution for actual electrons. It should be noted

that the expression (7) for $\rho(x,t)$ contains no $+/-$ cross terms arising from interference between paths arriving at (x,t) with right-going (+) and those with left-going (-) velocities. Hence, at least for the particles of the checkerboard model, the probability density $\rho(x,t)$ can be decomposed into two contributions, one associated only with right-going arrivals at (x,t) and the other only with left-going arrivals, $\rho(x,t) = \rho_+(x,t) + \rho_-(x,t)$ with $\rho_{\pm}(x,t) \equiv |\Psi_{\pm}(x,t)|^2$. Now, recall that particles following checkerboard paths move only at speed c and that the time between reversals in their directions of motion is $\approx (t_{BA}/\tau_{BA})\lambda_c/c$ for those paths that make the dominant contribution to the propagator. Hence, for Δt much less than this correlation time, nearly all of the particles in the spatial interval $[X - c\Delta t, X]$ that are right-going at time $T - \Delta t$ should arrive at $x=X$ during the time interval $[T - \Delta t, T]$, giving the dominant contribution to the number of right-going arrivals at X during that time interval. Taking the limit $\Delta t \rightarrow dt = 0^+$ this leads to the prediction that $c\rho_+(X,T)dt$ right-going particles arrive at $x=X$ during the infinitesimal time interval $[T - dt, T]$. A similar argument applies for left-going particles in $[X, X + c\Delta t]$ at time $T - \Delta t$. Hence, the distribution of arrival times T at the spatial point $x=X$ is given by

$$\Pi(T;X) = \Pi_+(T;X) + \Pi_-(T;X);$$

$$\Pi_{\pm}(T;X) = |\Psi_{\pm}(X,T)|^2 \int_0^{T_{max}} dt \rho(X,t), \quad (13)$$

where T_{max} is the maximum arrival time of practical interest. For the special case of free motion in one spatial dimension, the general results presented without derivation or discussion by Wigner [9] simplify to Eq. (13) [8].

III. FIRST-ARRIVAL PROPAGATOR FOR A DIRAC PARTICLE IN (1+1) DIMENSIONS

The fundamental constants \hbar and c are set to 1 in the analysis of this section but restored in the final results.

To begin, suppose that the spatial interval of interest is divided into M pieces each of length $\Delta x = x_{BA}/M$, assuming $x_B > x_A$. Suppose further that a particular path of N time steps consists of P spatial steps of length Δx to the right and Q to the left so that $N = P + Q$ and $M = P - Q$. The resulting space-time grid of path segments available to the particle is illustrated in Fig. 1. Denote by $K_{\beta\alpha}^{(1)}(B;A)$ the component of the propagator (9) associated with first arrivals at $X = x_B$ at time $T = t_B$. It is constructed from only those paths that reach x_B for the *first time* at time t_B . To compute $K_{\beta\alpha}^{(1)}$ we have to count the number of $\beta\alpha$ paths with R noncompulsary reversals for the restricted space-time grid shown in Fig. 1. For simplicity it is assumed that the initial wave function $\Psi(x, t_A = 0)$ is sufficiently well localized to the left of $x_B = X$ that $K_{++}^{(1)}(B;A)$ and $K_{+-}^{(1)}(B;A)$ are the only components of the first-arrival propagator that need to be considered. For $x_A < x_B$ the components $K_{-+}(B;A)$ and $K_{--}(B;A)$ of course have contributions only from multiple-arrival paths.

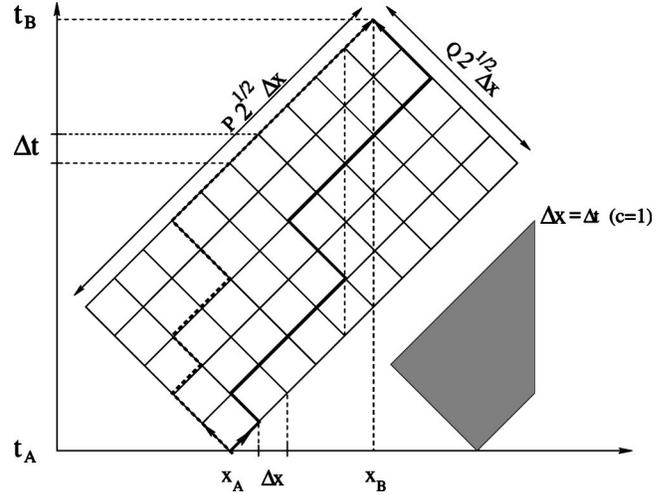


FIG. 1. Checkerboard grid with $\Delta x = \Delta t$ in the $x-t$ plane ($c = 1$). Two paths with five corners, the last of which is compulsory, are shown. The solid-line path with $\alpha = +$ and $\beta = -$ does not contribute to the first-arrival propagator $K^{(1)}(x_B, t_B; x_A, t_A)$ while the dashed-line path with $\alpha = -$ and $\beta = +$ does contribute. The restricted domain for evaluation of the first-arrival propagator is bounded on the right by the vertical dashed line; its shape is shown in grey.

A. Counting the number of first-arrival paths with r corners

Computation of $\Phi_{\beta\alpha}^{(1)}(R)$ involves counting the number of restricted $\beta\alpha$ paths with a given number of corners in a lattice. A path in the lattice of Fig. 2—obtained from Fig. 1 by clockwise rotation through 45° and rescaling by a factor of $(2^{1/2}\Delta x)^{-1}$ —is built up of $(1,0) = \rightarrow$ and $(0,1) = \uparrow$ motions. There are no \leftarrow and \downarrow motions because the paths of Fig. 1 move only forward in time. Denote by a l corner, a point on the path that is reached by a \rightarrow step and is left by a \uparrow step and by a r corner, a point that is reached by a \uparrow step and left by a \rightarrow step. Denote by R_l the number of corners of type l and by R_r the number of corners of type r in a given path. Any path in the $u-v$ lattice with given initial and final points (u_A, v_A) and (u_B, v_B) can be completely specified by either the coordinates of its l corners or by the coordinates of its r corners. Both specifications are needed in the following derivation of the first-arrival propagator.

First, consider the counting problem without any $\beta\alpha$ or first-arrival restrictions on the paths and including compulsory as well as noncompulsary corners. In this simple case,

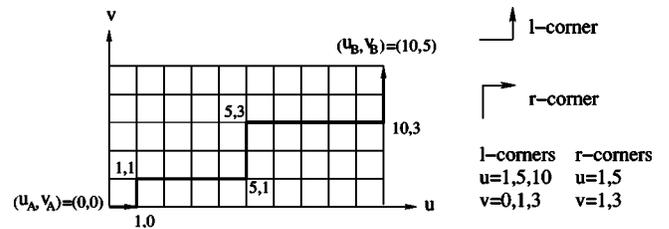


FIG. 2. Example of an unrestricted path that would not be allowed once first-arrival restrictions are introduced. The u and v sequences corresponding to the three l corners and two r corners of the path are indicated.

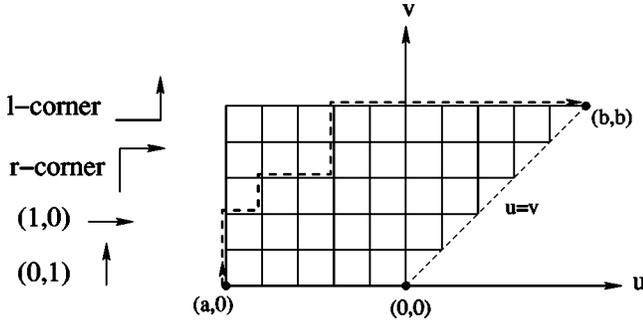


FIG. 3. A restricted path from $(u_A, v_A) = (a, 0)$ to $(u_B, v_B) = (b, b)$ which has two l corners and three r corners.

enumeration of the number of paths in a lattice with a given number of l or r corners is solved in the following manner [15]. To count the number of paths with R_l l -corners one builds two vectors \mathbf{u} and \mathbf{v} that contain the integer u and v coordinates of the R_l corners of such a path:

$$\begin{aligned} \mathbf{u} &= (u_1, u_2, \dots, u_{R_l}), \\ \mathbf{v} &= (v_1, v_2, \dots, v_{R_l}). \end{aligned} \quad (14)$$

These coordinates satisfy the inequalities

$$u_A + 1 \leq u_1 < u_2 < \dots < u_{R_l} \leq u_B, \quad (15)$$

$$v_A \leq v_1 < v_2 < \dots < v_{R_l} \leq v_B - 1,$$

where the sequences are strictly ordered. Denote by $\Phi_l^{wr}(R_l)$ the number of paths with R_l l -corners, where the label “ wr ” is a reminder that the paths are without restriction. This number may be evaluated by observing that there are $u_B - u_A$ integers from which to choose the u coordinates and $v_B - v_A$ from which to choose the v coordinates. The required number is then

$$\Phi_l^{wr}(R_l) = \binom{u_B - u_A}{R_l} \binom{v_B - v_A}{R_l}. \quad (16)$$

Similarly, one obtains

$$\Phi_r^{wr}(R_r) = \binom{u_B - u_A}{R_r} \binom{v_B - v_A}{R_r}. \quad (17)$$

Now consider only those paths which do not *cross* (touching is allowed for the moment) $x = x_B$ before $t = t_B$. When considering such restricted paths in the $u - v$ lattice (see, Fig. 3) it is convenient to choose the bottom-right corner of the accessible region as the origin of the $u - v$ coordinate system so that the region below the diagonal $u = v$ is out-of-bounds.

First consider the case in which the paths are specified by the coordinates of their l corners. This case is the easier of the two because l corners are diagnostic, i.e., a necessary and sufficient condition for a restricted path is that none of its l corners be on the forbidden $u > v$ side of the diagonal. It is thus required to calculate the number of paths with R_l l corners from (u_A, v_A) to (u_B, v_B) with $u_A = a$, $v_A = 0$ and u_B

$= v_B = b$, say, such that $v_i \geq u_i$ for all i . This number, $\Phi_l(R_l; BA)$, is the number (16) minus the number of paths with R_l l -corners such that $u_i > v_i$ for at least one i . To calculate the latter number, take k to be the largest integer such that $u_k > v_k$ and build from (15) the following sequences:

$$\begin{aligned} u_A + 1 &= a + 1 \leq u_1 < u_2 < \dots < u_{k-1} < v_{k+1} < \dots \\ &< v_{R_l} \leq v_B - 1 = b - 1, \end{aligned} \quad (18)$$

$$v_A = 0 \leq v_1 < v_2 < \dots < v_k < u_k < u_{k+1} < \dots < u_{R_l} \leq u_B = b.$$

The sequences are ordered [as can be checked, see Ref. [15]] and there is a one-to-one correspondence between the double sequences (15) and the double sequences (18). The number of all the double sequences (18) is

$$\binom{b - a - 1}{R_l - 1} \binom{b + 1}{R_l + 1}, \quad (19)$$

and therefore $\Phi_l(R_l; BA)$ is given by

$$\Phi_l(R_l; BA) = \binom{b - a}{R_l} \binom{b}{R_l} - \binom{b - a - 1}{R_l - 1} \binom{b + 1}{R_l + 1}. \quad (20)$$

Now consider the computation of the corresponding number of restricted paths from $(a, 0)$ to (b, b) with R_r r -corners, i.e., $\Phi_r(R_r; BA)$. A necessary condition for such a path is, of course, that all of its R_r r corners be on the allowed $v > u$ side of the diagonal. However, this is not a sufficient condition because it is possible for the l corner between two consecutive such r corners to be on the forbidden side of the diagonal. An additional complication is that the first l corner might precede the first r corner and/or the last l corner might follow the last r corner. A simple way to include such l corners in the analysis is to add the end-points $(a, 0) \equiv (u_0, v_0)$ and $(b, b) \equiv (u_{R_r+1}, v_{R_r+1})$ to the set of R_r r corners to obtain the set of $R_r + 2$ points $\{(u_0, v_0), (u_1, v_1), \dots, (u_{R_r}, v_{R_r}), (u_{R_r+1}, v_{R_r+1})\}$. Now, for $1 \leq i \leq R_r - 1$, (u_{i+1}, v_i) is the (diagnostic) l corner between the consecutive r -corners (u_i, v_i) and (u_{i+1}, v_{i+1}) . Hence, $\Phi_r(R_r; BA)$ is the number (17) minus the number of paths with R_r r corners such that $u_{i+1} > v_i$ for at least one i , with $0 \leq i \leq R_r$, to allow for paths for which the first and/or last corner is an l corner. To calculate the latter number, take k to be the smallest integer such that $u_{k+1} > v_k$ and construct the sequences

$$\begin{aligned} a = u_0 &\leq u_1 < u_2 < \dots < u_k < v_{k+1} < \dots < v_{R_r} \leq v_{R_r+1} = b, \\ 1 = v_0 + 1 &\leq v_1 < v_2 < \dots < v_k < u_{k+1} < \dots \\ &< u_{R_r} \leq u_{R_r+1} - 1 = b - 1. \end{aligned} \quad (21)$$

The total number of these double sequences is

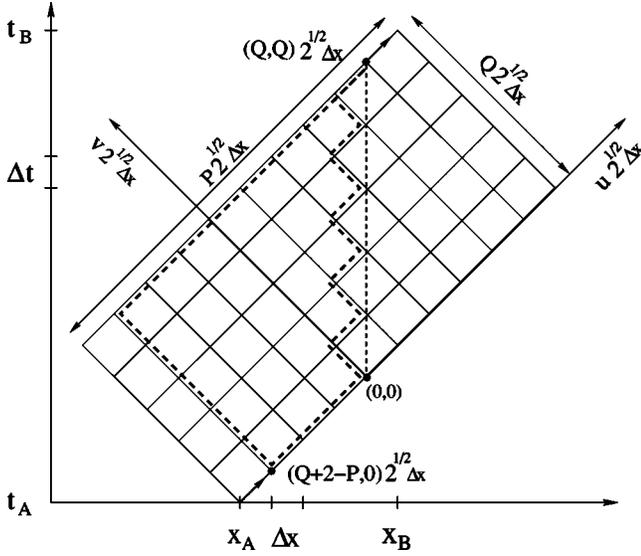


FIG. 4. The dashed line demarcates the grid to be considered in the enumeration problem related to $K_{++}^{(1)}$.

$$\binom{b-a+1}{R_r} \binom{b-1}{R_r}, \quad (22)$$

and therefore $\Phi_r(R_r; BA)$ is given by

$$\Phi_r(R_r; BA) = \binom{b-a}{R_r} \binom{b}{R_r} - \binom{b-a+1}{R_r} \binom{b-1}{R_r}. \quad (23)$$

Finally, consider the desired $\beta\alpha$ first-arrival paths, which are not allowed to *touch* $x=x_B$ before $t=t_B$. The various notations $A, (x_A, t_A)$ and $(u_A, v_A)2^{1/2}\Delta x$ are reserved for the first point of a checkerboard path and $B, (x_B, t_B)$ and $(u_B, v_B)2^{1/2}\Delta x$ for the last point. Denote by $A_\alpha, (x_A + \alpha\Delta x, t_A + \Delta t)$ and $(a_\alpha, 0)2^{1/2}\Delta x$ the point on a $\beta\alpha$ path at time $t=t_A + \Delta t$ and by $B_\beta, (x_B - \beta\Delta x, t_B - \Delta t)$, $(b_\beta, b_\beta)2^{1/2}\Delta x$ the point at time $t=t_B - \Delta t$.

It is important to note that those paths which may touch but do not cross the diagonal $u=v$, extending from $(0,0)$ to (b_β, b_β) , do not touch $x=x_B$ before $t=t_B$ and hence are first-arrival paths. Hence, the above expressions for $\Phi_l(R_l; BA)$ and $\Phi_r(R_r; BA)$ with A, B, a , and b replaced by $A_\alpha, B_\beta, a_\alpha$ and b_β , respectively, can be used to evaluate $\Phi_{\beta\alpha}^{(1)}$ and hence $K_{\beta\alpha}^{(1)}$. It is only necessary, for a given choice of α and β , to determine a_α, b_β and the relation between R and R_l or R_r , keeping in mind that R includes only noncompulsary corners.

For $\beta\alpha=++$ it is clear from Fig. 4 that $(b_+, b_+) = (Q, Q)$, $(u_B, v_B) = (Q+1, Q)$, $(u_A, v_A) = (Q+1-P, 0)$ and $(a_+, 0) = (Q+2-P, 0)$. Then a little thought leads to the conclusion that the number $\Phi_{++}^{(1)}(R)$ is equal to $\Phi_r(R_r; B_+A_+)$ with the identification $R=2R_r-1$. Hence, upon replacing a by $a_+ = Q+2-P$, b by $b_+ = Q$, and R_r by $(R+1)/2$ in Eq. (23), it follows that

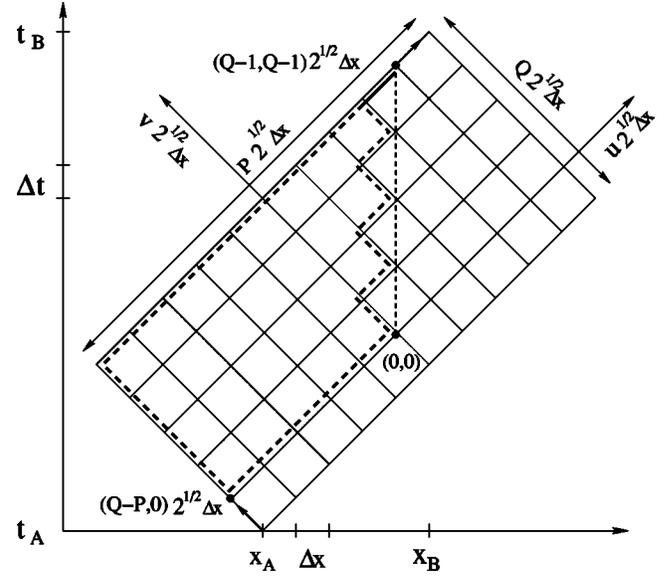


FIG. 5. The dashed line demarcates the grid to be considered in the enumeration problem related to $K_{+-}^{(1)}$.

$$\begin{aligned} \Phi_{++}^{(1)}(R) &= \binom{P-2}{(R+1)/2} \binom{Q}{(R+1)/2} \\ &\quad - \binom{P-1}{(R+1)/2} \binom{Q-1}{(R+1)/2} \\ &= \binom{P-1}{(R+1)/2} \binom{Q-1}{(R-1)/2} \\ &\quad - \binom{P-2}{(R-1)/2} \binom{Q}{(R+1)/2}, \end{aligned} \quad (24)$$

where the identity

$$\binom{n-1}{k} = \binom{n}{k} - \binom{n-1}{k-1}$$

has been used.

For $\beta\alpha=+-$ it is clear from Fig. 5 that $(b_+, b_+) = (Q-1, Q-1)$, $(u_B, v_B) = (Q, Q-1)$, $(a_-, 0) = (Q-P, 0)$, $(u_A, v_A) = (Q-P, -1)$ and $\Phi_{+-}^{(1)}(R)$ is equal to $\Phi_l(R_l; B_+A_-)$ with $R=2R_l$. Hence, upon replacing a by $a_- = Q-P$, b by $b_+ = Q-1$ and R_l by $R/2$ in Eq. (20), it follows that

$$\Phi_{+-}^{(1)}(R) = \binom{P-1}{R/2} \binom{Q-1}{R/2} - \binom{P-2}{(R-2)/2} \binom{Q}{(R+2)/2}, \quad (25)$$

with $\Phi_{+-}^{(1)}(1) = 1$.

B. Evaluation of the propagators

The first-arrival-time propagators are expressed as

$$K_{++}^{(1)}(B;A) = \frac{i}{2\lambda_c} \lim_{N \rightarrow \infty} \sum_{\text{odd} R \geq 1} \Phi_{++}^{(1)}(R)(im\Delta t)^R, \quad (26)$$

$$K_{+-}^{(1)}(B;A) = \frac{i}{2\lambda_c} \lim_{N \rightarrow \infty} \sum_{\text{even} R \geq 0} \Phi_{+-}^{(1)}(R)(im\Delta t)^R.$$

C. The case of $K_{++}^{(1)}$

In this case we have from Eq. (26) and using the approximation $\binom{n}{m} \approx n^m/m!$, that becomes exact as $n \rightarrow \infty$,

$$\begin{aligned} K_{++}^{(1)}(B;A) &= \frac{i}{2\lambda_c} \lim_{N \rightarrow \infty} \sum_{\text{odd} R \geq 1} (im\Delta t)^R \left(\frac{P^{(R+1)/2}}{[(R+1)/2]!} \frac{Q^{(R-1)/2}}{[(R-1)/2]!} - \frac{P^{(R-1)/2}}{[(R-1)/2]!} \frac{Q^{(R+1)/2}}{[(R+1)/2]!} \right) \\ &= \frac{i}{2\lambda_c} \lim_{N \rightarrow \infty} \sum_{\text{odd} R \geq 1} (im\Delta t)^R (P-Q) \frac{(PQ)^{(R-1)/2}}{[(R+1)/2]! [(R-1)/2]!}. \end{aligned} \quad (27)$$

This expression may be transformed [16] to

$$K_{++}^{(1)}(B;A) = i \frac{x_{BA}}{\lambda_c l_{BA}} \lim_{N \rightarrow \infty} \sum_{\text{odd} R \geq 1} \left(\frac{il_{BA}}{2\lambda_c} \right)^R \frac{1}{[(R+1)/2]! [(R-1)/2]!} = - \frac{x_{BA}}{\lambda_c l_{BA}} \frac{l_{BA}}{2\lambda_c} \sum_{k=0}^{\infty} (-)^k \left(\frac{l_{BA}}{2\lambda_c} \right)^{2k} \frac{1}{k!(k+1)!}. \quad (28)$$

In the last line R has been replaced by $2k+1$ and the limit $N \rightarrow \infty$ taken. Comparison with the power series representation of the Bessel function,

$$J_n(z) = \left(\frac{z}{2} \right)^n \sum_{k=0}^{\infty} \frac{(-)^k}{k!(k+n)!} \left(\frac{z}{2} \right)^{2k}, \quad (29)$$

immediately gives

$$K_{++}^{(1)}(B;A) = - \frac{x_{BA}}{\lambda_c l_{BA}} J_1 \left(\frac{l_{BA}}{\lambda_c} \right). \quad (30)$$

D. The case of $K_{+-}^{(1)}$

In this case we start with the expression

$$\begin{aligned} K_{+-}^{(1)}(B;A) &= \frac{i}{2\lambda_c} \lim_{N \rightarrow \infty} \sum_{\text{even} R \geq 0} (im\Delta t)^R \left(\frac{P^{R/2}}{(R/2)!} \frac{Q^{R/2}}{(R/2)!} \right. \\ &\quad \left. - \frac{P^{(R-2)/2}}{[(R-2)/2]!} \frac{Q^{(R+2)/2}}{[(R+2)/2]!} \right). \end{aligned} \quad (31)$$

Steps analogous to those above lead to

$$K_{+-}^{(1)}(B;A) = \frac{i}{2\lambda_c} \left[J_0 \left(\frac{l_{BA}}{\lambda_c} \right) + \frac{c-v_{BA}}{c+v_{BA}} J_2 \left(\frac{l_{BA}}{\lambda_c} \right) \right]. \quad (32)$$

An interesting equality emerges from the above results, namely,

$$K_{++}(B;A) - K_{--}(B;A) = K_{+-}^{(1)}(B;A). \quad (33)$$

In fact all paths contributing to the $K_{--}(B;A)$ component of the propagator touch the line $x=x_B$ at least twice and therefore this set of paths is complementary to that contributing to $K_{++}^{(1)}(B;A)$ in the limit of $N \rightarrow \infty$.

Finally,

$$\begin{aligned} K_{++}^{(2,3,\dots)}(B;A) &= K_{++}(B;A) - K_{++}^{(1)}(B;A) \\ &= K_{--}(B;A) \\ &= \frac{x_{BA} - ct_{BA}}{2\lambda_c l_{BA}} J_1 \left(\frac{l_{BA}}{\lambda_c} \right); \\ K_{+-}^{(2,3,\dots)}(B;A) &= K_{+-}(B;A) - K_{+-}^{(1)}(B;A) \\ &= - \frac{i}{2\lambda_c} \frac{c-v_{BA}}{c+v_{BA}} J_2 \left(\frac{l_{BA}}{\lambda_c} \right). \end{aligned} \quad (34)$$

IV. INTERFERENCE BETWEEN FIRST AND LATER ARRIVALS

The decomposition $K_{\beta\alpha}(B;A) = K_{\beta\alpha}^{(1)}(B;A) + K_{\beta\alpha}^{(2,3,\dots)}(B;A)$ according to first and later (second, third, etc.) arrivals of a particle at x_B at t_B leads immediately to the corresponding decomposition $\Psi_{\beta}(x_B, t_B) = \Psi_{\beta}^{(1)}(x_B, t_B) + \Psi_{\beta}^{(2,3,\dots)}(x_B, t_B)$ for the $\beta = \pm$ components of the wave function $\Psi(x_B, t_B)$. Substitution of the latter expression, with $t_A=0$, $x_B=X$, and $t_B=T$, into the result (13) for the arrival time distribution gives

$$\Pi(T;X) = \Pi^{(1)}(T;X) + \Pi^{(2,3,\dots)}(T;X) + \Pi^{(1 \times 2,3,\dots)}(T;X). \quad (35)$$

For the special case considered here in which the initial probability density $\rho(x_A, 0)$ is negligible for $x_A \geq X$,

$$\Pi^{(1)}(T; X) = C \left| \int_{-\infty}^X dx_A [K_{++}^{(1)}(X, T; x_A, 0) \Psi_+(x_A, 0) + K_{+-}^{(1)}(X, T; x_A, 0) \Psi_-(x_A, 0)] \right|^2 \quad (36)$$

is the contribution of first arrivals,

$$\begin{aligned} \Pi^{(2,3,\dots)}(T; X) = & C \left| \int_{-\infty}^X dx_A [K_{++}^{(2,3,\dots)}(X, T; x_A, 0) \Psi_+(x_A, 0) + K_{+-}^{(2,3,\dots)}(X, T; x_A, 0) \Psi_-(x_A, 0)] \right|^2 \\ & + C \left| \int_{-\infty}^X dx_A [K_{-+}^{(2,3,\dots)}(X, T; x_A, 0) \Psi_+(x_A, 0) + K_{--}^{(2,3,\dots)}(X, T; x_A, 0) \Psi_-(x_A, 0)] \right|^2 \end{aligned} \quad (37)$$

is the contribution of later arrivals and

$$\begin{aligned} \Pi^{(1 \times 2,3,\dots)}(T; X) = & 2C \Re \left\{ \int_{-\infty}^X dx_A [K_{++}^{(1)}(X, T; x_A, 0) \Psi_+(x_A, 0) + K_{+-}^{(1)}(X, T; x_A, 0) \Psi_-(x_A, 0)]^* \right. \\ & \left. \times \int_{-\infty}^X dx_A [K_{++}^{(2,3,\dots)}(X, T; x_A, 0) \Psi_+(x_A, 0) + K_{+-}^{(2,3,\dots)}(X, T; x_A, 0) \Psi_-(x_A, 0)] \right\} \end{aligned} \quad (38)$$

is the contribution due to interference between first and later arrivals. [C is the normalization factor appearing in (13)]. Of particular interest here is the magnitude of the interference contribution relative to the first-arrival contribution in the regime $v_{BA} \ll c$.

First, however, briefly consider the regime in which v_{BA} is so close to c that the correlation distance $\lambda_c [1 - (v_{BA}/c)^2]^{-1/2}$ for reversal of direction is sufficiently large that for a typical checkerboard path there is insufficient time for more than one arrival at x_B . To be more quantitative, assume that the initial amplitude $\Psi_-(x_A, t_A)$ of the $-c$ velocity eigenstate is completely negligible with respect to the initial amplitude $\Psi_+(x_A, t_A)$ of the $+c$ eigenstate so that one need consider only $K_{++}(B; A)$ and $K_{-+}(B; A)$. Also, for $(x_B, t_B) = (X, T)$ and $(x_A, t_A) = (x_A, 0)$ assume that $x_{BA} = X - x_A$ is very close to $ct_{BA} = cT$ for those values of T for which $\Pi(T; X)$ is non-negligible and for those values of x_A for which $\rho(x_A, 0)$ is non-negligible. In this regime, $K_{++}^{(1)}(B; A)/K_{++}(B; A) \approx 1 - \delta/2$ and $K_{-+}^{(2,3,\dots)}(B; A)/K_{++}(B; A) \approx \delta/2$, where $v_{BA} \equiv (1 - \delta)c$ with $\delta \ll 1$. In addition, $|K_{-+}(B; A)/K_{++}(B; A)| \approx (l_{BA}/2ct_{BA}) J_0(l_{BA}/\lambda_c) / J_1(l_{BA}/\lambda_c)$ with $l_{BA} \approx (2\delta)^{1/2} ct_{BA}$. If δ is sufficiently small that $l_{BA} \ll \lambda_c$ then, using the leading term in Eq. (29) for $n=0$ and for $n=1$, $|K_{-+}(B; A)/K_{++}(B; A)| \approx \lambda_c / x_{BA} = \lambda_c / (X - x_A)$ which is typically very much less than 1 for an initial wave packet $\Psi(x_A, 0)$ that is well-localized away from $x=X$. Hence, at least to the extent that the concepts of single-particle probability and probability current densities are still meaningful in the regime in which v_{BA} is very close to c , the interference term is very small and $\Pi(T; X) = \Pi^{(1)}(T; X)$ to a good approximation. Strictly speaking, however, for the special case

under consideration the first-arrival-time distribution is well-defined only in the limit $v_{BA} \rightarrow c$.

Now, consider the nonrelativistic regime. With the definitions $\Psi_{\pm}(x, t) \equiv \phi_{\pm}(x, t) \exp(-imc^2 t/\hbar)$, $\phi(x, t) \equiv \phi_+(x, t) + \phi_-(x, t)$ and $\Delta\phi(x, t) \equiv \phi_+(x, t) - \phi_-(x, t)$, the (1+1)-dimensional free-electron Dirac equation (1) can be written as

$$i\hbar \partial\phi(x, t)/\partial t = -i\hbar c \partial\Delta\phi(x, t)/\partial x - 2mc^2 \phi(x, t),$$

$$i\hbar \partial\Delta\phi(x, t)/\partial t = -i\hbar c \partial\phi(x, t)/\partial x. \quad (39)$$

If $i\hbar \partial\phi/\partial t$ is negligible with respect to $2mc^2 \phi$ (with c fixed at its actual value, not set equal to infinity) then ϕ can be replaced by $-(i\hbar/2mc) \partial\Delta\phi/\partial x$ in the second equation of Eq. (39) to obtain the Schrödinger equation $i\hbar \partial\Delta\phi(x, t)/\partial t = -(\hbar^2/2m) \partial^2 \Delta\phi(x, t)/\partial x^2$. If one further assumes that $|\Delta\phi(x, t)|^2 \gg |\phi(x, t)|^2$ and identifies $2^{-1/2} \Delta\phi(x, t)$ with the Schrödinger wave function $\psi_S(x, t)$ then the expressions (7) and (8) immediately lead to the desired nonrelativistic expressions, $\rho_S(x, t) = |\psi_S(x, t)|^2$ and $J_S(x, t) = (\hbar/m) \Im[\psi_S^*(x, t) \partial\psi_S(x, t)/\partial x]$, respectively, for the nonrelativistic probability and probability current densities. Consistent with these considerations is the following simple choice of initial ($t=0$) wave function $\Psi(x_A, 0)$ for the nonrelativistic regime: $\Psi_+(x_A, 0) = -g\Psi_0(x_A)$ and $\Psi_-(x_A, 0) = (1-g^2)^{1/2} \Psi_0(x_A)$ where g is a real constant very close to $2^{-1/2}$ (see below) and $\Psi_0(x) = (2\pi)^{-1/4} (\Delta x)^{-1/2} \exp[-(\Delta x)^{-2}(x-x_0)^2 + ik_0 x]$ is a minimum-uncertainty-product gaussian with initial centroid x_0 , initial variance Δx , mean

wave vector k_0 and variance $\Delta k = 1/2\Delta x$. In the numerical calculations presented below the constant g is chosen so that the characteristic velocity $v(x_A, 0) \equiv J(x_A, 0)/\rho(x_A, 0) = (2g^2 - 1)c \equiv v$ (independent of x_A) is equal to $v_0 \equiv \hbar k_0/m$ with $v_0 \ll c$.

Now, from Eqs. (10), (11) and (30) it immediately follows

$$\frac{K_{+-}^{(1)}(B;A)}{K_{+--}^{(1)}(B;A)} = \frac{J_0(l_{BA}/\lambda_c) + [(c - v_{BA})/(c + v_{BA})]J_2(l_{BA}/\lambda_c)}{J_0(l_{BA}/\lambda_c)} \approx \frac{J_0(l_{BA}/\lambda_c) + J_2(l_{BA}/\lambda_c) - (2v_{BA}/c)J_2(l_{BA}/\lambda_c)}{J_0(l_{BA}/\lambda_c)}. \quad (41)$$

In the regime $v_{BA}/c \ll 1$ under consideration, $l_{BA}/\lambda_c \approx ct_{BA}/\lambda_c = (c/v_{BA})(x_{BA}/\lambda_c) \gg x_{BA}/\lambda_c \gg 1$. Using the leading two terms in Hankel's asymptotic expansions [17] of $J_0(z)$ and $J_2(z)$ for large argument z , i.e.,

$$J_0(z) \approx \left(\frac{2}{\pi z}\right)^{1/2} \left[\cos\left(z - \frac{\pi}{4}\right) + \frac{1}{8z} \sin\left(z - \frac{\pi}{4}\right) \right] (|z| \gg 1),$$

$$J_2(z) \approx \left(\frac{2}{\pi z}\right)^{1/2} \left[\cos\left(z - \frac{5\pi}{4}\right) - \frac{15}{8z} \sin\left(z - \frac{5\pi}{4}\right) \right] (|z| \gg 1) \quad (42)$$

gives

$$\frac{K_{+-}^{(1)}(B;A)}{K_{+--}^{(1)}(B;A)} \approx 2 \left(\frac{v_{BA}}{c}\right) \times \frac{\cos\left(\frac{l_{BA}}{\lambda_c} - \frac{\pi}{4}\right) + \frac{\lambda_c}{16x_{BA}} \sin\left(\frac{l_{BA}}{\lambda_c} - \frac{\pi}{4}\right)}{\cos\left(\frac{l_{BA}}{\lambda_c} - \frac{\pi}{4}\right) + \frac{\lambda_c}{8l_{BA}} \sin\left(\frac{l_{BA}}{\lambda_c} - \frac{\pi}{4}\right)}. \quad (43)$$

For $l_{BA}/\lambda_c = (4n+3)\pi/4$ with n an integer, the right-hand side of Eq. (43) is 1. However, in a well-designed arrival-time experiment for the wave function under discussion one would arrange that $(X - x_0) \gg \Delta x$ so that $\rho(x_A, 0)$ is completely negligible for $x_A \geq X$ and also that $\Delta x \gg \lambda_c$ so that there is negligible probability of generating particle-antiparticle pairs. Hence, λ_c/x_{BA} would be extremely small over the important range of x_{BA} . Hence, the set of x_{BA} values where the right-hand side of Eq. (43) is not close to $2v_{BA}/c$ is of small measure and can be ignored when considering integrals over x_A , provided that v_{BA}/c is not itself extremely small (a rough estimate requires that $v_{BA}/c \gg \lambda_c/16\pi x_{BA}$).

that

$$K_{++}^{(1)}(B;A)/K_{+-}^{(1)}(B;A) = 2v_{BA}/(c \pm v_{BA}) \approx 2v_{BA}/c \quad (40)$$

and from Eq. (12) and (32) it follows that

Taking into account that $-\Psi_-(x_A, 0) \approx \Psi_+(x_A, 0)$ for $|v(x_A, 0)| \ll c$ and that the terms involving $K_{--}^{(1)}$ and $K_{+-}^{(1)}$ are negligible when $|X - x_0| \gg \Delta x$ then leads directly to the estimates

$$\Pi^{(1)}(T;X) \approx (1/2)(2v/c)^2 \Pi(T;X) = 2(v/c)^2 \Pi(T;X), \quad (44)$$

$$\Pi^{(1 \times 2, 3, \dots)}(T;X) \approx 2(1/2)(2v/c) \Pi(T;X) = 2(v/c) \Pi(T;X) \quad (45)$$

for the Gaussian wave function under consideration. It should be noted that $v_{BA} = (X - x_A)/T$ has been approximated by v which is consistent with $(X - x_0) \gg \Delta x$ in the absence of significant wave packet spreading. Figures 6 and

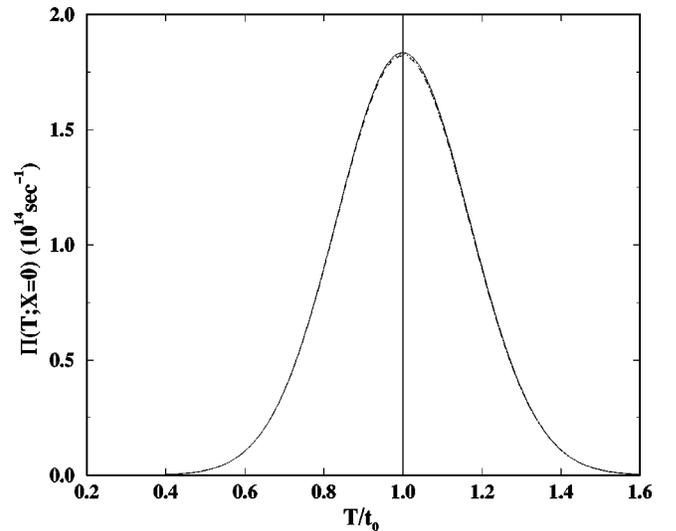


FIG. 6. Arrival-time distribution $\Pi(T;0)$ (solid curve), scaled first-arrival contribution $\Pi^{(1)}(T;0)/2(v/c)^2$ (dotted line) and scaled interference contribution $\Pi^{(1 \times 2, 3, \dots)}(T;0)/2(v/c)$ (dashed line) for the initial gaussian wave function described in the text. $k_0 = 1.00 \text{ \AA}^{-1}$, $\Delta k = 0.02 \text{ \AA}^{-1}$ and $x_0 = -6\Delta x$ with $\Delta x = 1/2\Delta k$; $v/c = v_0/c = 3.862 \times 10^{-3}$; $t_0 \equiv |x_0|/v_0 = 1.296 \times 10^{-14} \text{ sec}$; and $T_{max} = 2t_0$.

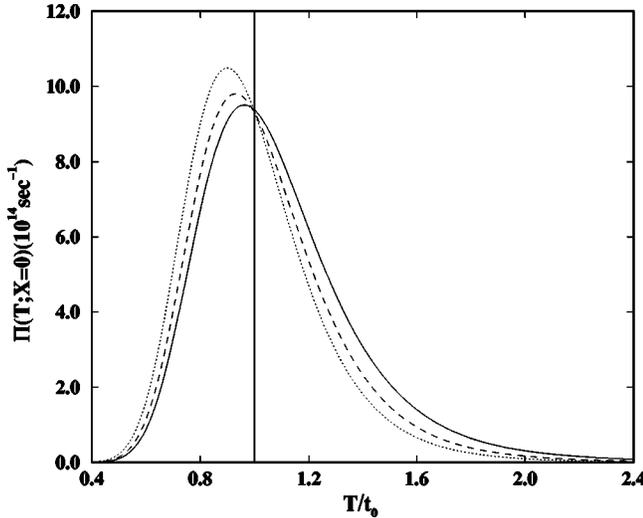


FIG. 7. Arrival-time distribution $\Pi(T;0)$ (solid line), scaled first-arrival contribution $\Pi^{(1)}(T;0)/2(v/c)^2$ (dotted line) and scaled interference contribution $\Pi^{(1 \times 2,3,\dots)}(T,0)/2(v/c)$ (dashed line) for the initial gaussian wave function described in the text. $k_0 = 1.00 \text{ \AA}^{-1}$, $\Delta k = 0.2 \text{ \AA}^{-1}$ and $x_0 = -8\Delta x$ with $\Delta x = 1/2\Delta k$; $v/c = v_0/c = 3.862 \times 10^{-3}$; $t_0 = 1.728 \times 10^{-15} \text{ sec}$; and $T_{max} = 3t_0$.

7 show results for $\Pi(T;0)$, $\Pi^{(1)}(T;0)/2(v/c)^2$ and $\Pi^{(1 \times 2,3,\dots)}(T;0)/2(v/c)$ obtained by numerical evaluation of Eqs. (36) to (38) for gaussian wave packets with $\Delta k/k_0 = 0.02$ and 0.2 , respectively. In the former case the above estimates are excellent approximations; in the latter case, even though wave packet spreading is more important, the estimates still provide good approximations for the very large differences in overall scale between the three quantities.

V. CONCLUDING REMARKS

In summing up, it is interesting to make a qualitative comparison of the results of the Feynman path and Bohm trajectory approaches for investigating arrival times for Dirac electrons.

In Bohmian mechanics [18–20] an electron is postulated to be an actually existing point-like particle and an accompanying wave which guides its motion. For a Dirac electron in the presence of a potential $V(\vec{r},t)$, the time evolution of the guiding wave $\Psi(\vec{r},t)$ is described by the three-dimensional (3D) Dirac equation and the trajectory of the pointlike particle is determined by the equation-of-motion $d\vec{r}(t)/dt = [\vec{J}(\vec{r},t)/\rho(\vec{r},t)]|_{\vec{r}=\vec{r}(t)}$. It is further postulated that, for an ensemble of electrons all prepared in the same initial state $\Psi(\vec{r},0)$, the probability of such a particle *having* initial position $\vec{r}^{(0)} \equiv \vec{r}(t=0)$ is given by $\rho(\vec{r}^{(0)},0)$. The various properties stated below for the intrinsic arrival times of the pointlike particles of Bohm's theory follow readily from the fact that, for a given initial wave function, trajectories with different starting points $\vec{r}^{(0)}$ never intersect or even touch each other [21].

Now, the expression for the intrinsic 1D arrival-time dis-

tribution obtained with either approach can be cast in the same form as in classical mechanics, namely,

$$\Pi(T;X) = \Pi_+(T;X) + \Pi_-(T;X);$$

$$\Pi_{\pm}(T;X) \equiv \pm J_{\pm}(X,T) / \int_0^{T_{max}} dt [J_+(X,t) - J_-(X,t)], \quad (46)$$

where J_+ and J_- , respectively, are the right-going and left-going components of the probability current density J . The decomposition $J = J_+ + J_-$ is not uniquely defined. The decomposition associated with the fictitious particles of the checkerboard path approach, which move only at the speed of light c , is $J_{\pm} = \pm c|\Psi_{\pm}|^2$, while that associated with the (assumed) actual particles of the Bohm trajectory approach, each of which moves at a variable speed that cannot exceed c , is $J_{\pm} = J\Theta[\pm J]$, where Θ is the unit step function. The former decomposition leads to an arrival-time distribution $\Pi(T;X)$ proportional to the probability density $\rho(X,T)$ while the latter leads to one proportional to the absolute value of the probability current density, i.e., $|J(X,T)|$. Moreover, unless one or other (or both) of the two components of $\Psi(X,t)$ is zero for $T_1 \leq t \leq T_2$, there are more arrivals—many more if $T_2 - T_1$ is much larger than the Jacobson-Schulman correlation time—of the fictitious particles at X during that time interval than there are of the supposed actual particles of Bohm's theory.

Given the probability current density $J(x,t)$ and using the noncrossing property of Bohm trajectories it is straightforward to decompose the intrinsic arrival-time distribution into contributions from first arrivals, from second arrivals, etc. with no interference terms between different orders of arrival [22]. In marked contrast to this, the decomposition based on Feynman checkerboard paths in general contains a nonzero interference term between first and later (i.e., second, third, etc.) arrivals so that from the calculation one cannot extract a well-defined intrinsic first-arrival-time distribution. In the nonrelativistic regime this interference term can be very large compared to the first-arrival term. Because of this and the extremely small correlation length for reversal of direction ($\approx \lambda_c$ which is only about 10^{-3} of the diameter of an atom!), suppression of the interference term by decoherence [23] within a time interval much less than λ_c/c in duration immediately following the instant of first arrival would be very difficult, if not impossible, in a practical arrival-time measurement. Unless this can be achieved, assuming that the first-arrival times of the fictitious particles of the checkerboard model are directly relevant to the arrival times measured in a time-of-flight experiment on actual electrons is not justified.

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