Relativistic wave equations: an operational approach

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Abstract
The use of operator methods of an algebraic nature is shown to be a very powerful tool to deal with different forms of relativistic wave equations. The methods provide either exact or approximate solutions for various forms of differential equations, such as relativistic Schrödinger, Klein–Gordon, and Dirac. We discuss the free-particle hypotheses and those relevant to particles subject to non-trivial potentials. In the latter case we will show how the proposed method leads to easily implementable numerical algorithms.

Keywords: relativistic wave equations, operational methods, evolution operator

(Some figures may appear in colour only in the online journal)

1. Introduction
Operation techniques embedded with the formalism of the evolution operator have played a central role in the study of time-dependent problems in quantum mechanics. The associated technicalities have attracted the attention of physicists [1] and mathematicians [2], who provided tools of crucial importance to deal with the study of the time-dependent Schrödinger equation. Expansions of the evolution operator like those developed by Stueckelberg,

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Feynman, and Dyson on the physicists’ side and by Magnus and Fer on the other have paved the way to the formalism underlying the modern diagrammatic procedures [3–5]. Evolution problems in relativistic quantum mechanics have raised new questions mainly associated with the pseudo-differential nature of the involved operators. These methods, neither popular nor well appreciated by physicists in the past, are now getting increasing attention for a more sound formulation of the underlying mathematical foundations and for the wealth of phenomenology they can describe [6]. In particular, the operational approach has been shown to be effective in solving pseudo-differential equations as well as those containing fractional derivatives and related to anomalous transport phenomena. The latter not only include the so-called fractional dynamics adopted to investigate anomalous diffusion and similar effects in complex physical systems [7], but also appear useful to model problems in life sciences, engineering, economics, and even studies of human mobility [8–14].

In this paper we will focus our attention on the equations of relativistic quantum mechanics. Our aim is to push forward previous investigations of solutions to the relativistic evolution equation involving fractional derivatives, namely, the relativistic Schrödinger (called also Salpeter’s) equation [15, 16, 18] and to compare them with operationally obtained solutions to the Dirac-type and Klein–Gordon equations. The main motivation of this research is that fundamental relativistic wave equations give results whose interpretation, based on Hamiltonian mechanics concepts as well as explanations of arising paradoxes (like Zitterbewegung or the Klein paradox), still remains controversial [17] and needs clarification. In our opinion the use of fractional derivatives or of pseudo-differential operators to treat relativistic quantum problems allows a more complete understanding of longstanding problems associated with nonlocal evolution problems. These problems, once considered academic, have in recent years acquired a more concrete nature (mainly in the context of the so-called quantum simulation), namely, the possibility of ‘simulating’ quantum nonrelativistic effects in condensed matter and atomic physics [19].

We begin by recalling results coming from the study of the relativistic heat equation (as proposed in [15]), whose solutions provide us an example of the method. Section 3 is devoted to the analysis of the free relativistic Schrödinger equation and the unexpected behaviour of its solutions. In section 4 we consider the Dirac factorization of the relativistic Schrödinger equation and compare such obtained results with those of section 3. The Klein–Gordon equation, together with the relation of its solutions to the Dirac case, is studied in section 5, while in section 6 we present the analysis of the simplest ‘interacting model’, namely, the relativistic particle (Schrödinger’s and Dirac’s) driven by a linear potential. In what follows, we restrict ourselves to 1+1 spacetime models. This simplifies the presentation and preserves readability of the results, but there is no reason to assume that general properties of solutions will not be shared with the realistic 1+3 situation.

2. The relativistic heat equation

In reference [15] the following equation has been defined as the relativistic heat equation

\[ \partial_t F(x, t) = -\sqrt{d^2 - k^2} \partial_x^2 F(x, t), \]  

\[ F(x, 0) = f(x), \]  

where \( t \) has the dimension of time and \( d \) and \( k \) have the dimensions of the inverse time and of a length divided by time, respectively. Introducing dimensionless variables \( \tau = dt \) and \( \xi = x/dk \) we can cast equation (1) in the form
From the mathematical point of view, equations (1) and (3) are evolution equations with the distinctive feature that they contain the square root of a differential operator. The formal solution of equation (3) can be written as

\[ F(\xi, \tau) = e^{-\tau \sqrt{1 - \bar{\partial}^2}} f(\xi). \]  

Although the use of the exponential operators is quite common, the handling of exponentials containing the square root of a differential operator (as well as higher order roots) has been rarely treated. Nevertheless, some analytical tools have been developed for this purpose in the past. In the specific case of equation (4), a suitable representation, which allows evaluation of the action of ‘square rooted’ operator on the function \( f(\xi) \), can be realized through the Doetsch formula [20]. The latter allows explicit evaluation of the r.h.s. of equation (4) according to the following integral transform

\[
F(\xi, \tau) = \left[ \int_0^\infty g_{1/2}(\eta) e^{-\eta \sqrt{\eta^2 - \bar{\partial}^2}} d\eta \right] f(\xi), \tag{5}
\]

where \( D(\xi, \eta^2) = e^{\eta^2 \bar{\partial}^2} f(\xi) \). Note two peculiar features of equation (6):

- The solution of the relativistic heat equation (1) depends, via an appropriate transform, on the solution of the ‘classical’ diffusion problem, as \( D(\xi, \tau) \) is the solution of the ordinary heat equation

\[
\partial_\tau D(\xi, \tau) = \partial^2_\xi D(\xi, \tau), \tag{7}
\]

\[
D(\xi, 0) = f(\xi), \tag{8}
\]

usually expressible in terms of the Gauss–Weierstrass transform

\[
D(\xi, \tau) = e^{\frac{\tau}{\pi}} \int_{-\infty}^{\infty} e^{-\frac{\xi^2}{\tau}} f(\sigma) d\sigma, \tag{9}
\]

or written, using the Fourier transform, as

\[
D(\xi, \tau) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{i\xi q} \tilde{f}(q) dq, \tag{10}
\]

with \( \tilde{f}(q) = (2\pi)^{-1/2} \int_{-\infty}^{\infty} e^{-\frac{q^2}{4}} f(\xi) d\xi \).

- The integral kernel

\[
g_{1/2}(\eta) = \frac{1}{2\sqrt{\pi \eta}} e^{-\eta^{1/2} d\eta} \tag{11}
\]

is a generic example of the so-called Lévy stable distributions or \( \alpha \)-stable distributions. The latter, denoted as \( g_{\alpha}(\eta) \), are for \( \eta \geq 0 \) and for \( 0 < \alpha < 1 \), given uniquely by the inverse Laplace transform of \( e^{-\eta^\alpha} \), i.e., \( e^{-\eta^\alpha} = \int_{0}^{\infty} e^{-\eta g_{\alpha}(\eta)} d\eta \) [21–23].

Let us now consider a solution of the relativistic heat equation for two specific initial functions. First, we assume that the initial condition of equation (8) is a Gaussian:
where \( \beta > 0 \). We can therefore write the relevant solution as

\[
D_1(\beta; \xi, \tau) = \frac{1}{2\sqrt{\pi\beta}} \int_{-\infty}^{\infty} e^{-\beta(\xi^2 + \tau^2 + \xi\tau)} d\xi,
\]

(13)

recognized as the so-called Glaisher identity [24]. Equation (13), if substituted into equation (6), leads to

\[
F_1(\beta; \xi, \tau) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{-\beta\sqrt{\xi^2 + \tau^2}} e^{i\xi\tau} d\xi,
\]

(14)

the same as the result of formal manipulation on equations (4) and (12).

The second example of the initial condition under consideration is

\[
f_2(\beta; \xi) = \frac{1}{2K_1(\beta)} \int_{-\infty}^{\infty} e^{-\beta\sqrt{\xi^2 + 1 + \xi^2}} d\xi,
\]

(15)

where \( \beta > 0 \) and \( K_1(\beta) \) is the modified Bessel function of the second kind, i.e., the Macdonald function [25]. Equation (15) is essentially the Fourier transform of

\[
\tilde{f}_2(\beta; k) = \frac{1}{2K_1(\beta)} e^{-\beta\sqrt{1 + k^2}}
\]

(16)

normalized in the \( L^1 \)-norm: \( \int_{-\infty}^{\infty} \tilde{f}_2(\beta; k) dk = 1 \). The solution of the relativistic heat equation with equation (15) as the initial function, calculated as shown before, reads

\[
F_2(\beta; \xi, \tau) = \frac{e^{-\tau\sqrt{1 + \xi^2}}}{2K_1(\beta)} \int_{-\infty}^{\infty} e^{-\beta\sqrt{1 + \xi^2 + \xi\tau}} d\xi
\]

(17)

\[
= \frac{1}{2K_1(\beta)} \int_{-\infty}^{\infty} e^{-\beta(\xi + \tau)\sqrt{1 + \xi^2 + 1}} d\xi.
\]

The use of the identity (see [25, 26])

\[
\int_{-\infty}^{\infty} e^{-a\sqrt{\xi^2 + \xi^2}} \cos(bx) dx = \frac{2a q}{\sqrt{a^2 + b^2}} K_1\left( q\sqrt{a^2 + b^2} \right)
\]

(18)

allows us to cast equation (17) in the form

\[
F_2(\beta; \xi, \tau) = \frac{\beta + \tau}{2K_1(\beta)} \frac{K_1\left( \sqrt{(\beta + \tau)^2 + \xi^2} \right)}{\sqrt{(\beta + \tau)^2 + \xi^2}},
\]

(19)

which may be viewed as the identity

\[
e^{-\tau\sqrt{1 + \xi^2}} \frac{\beta}{2K_1(\beta)} \frac{K_1\left( \sqrt{\beta^2 + \xi^2} \right)}{\sqrt{\beta^2 + \xi^2}} = \frac{\beta + \tau}{2K_1(\beta)} \frac{K_1\left( \sqrt{(\beta + \tau)^2 + \xi^2} \right)}{\sqrt{(\beta + \tau)^2 + \xi^2}}
\]

(20)

and interpreted as a generalization of the Glaisher formula for the Macdonald function.

Concluding this section, we would like to note that naming equation (1) as a relativistic heat equation may lead to some misconception and should be used carefully, or even in
quotes. The concept of relativistic diffusion needs clarification, at least from the mathematical point of view. Namely, equation (1) does not represent a diffusion process in the strict sense, because the norm of the relevant solution is not preserved during the evolution and decays with $e^{-\tau}$. Equations (14) and (19) are properly normalized at any time if multiplied by $e^{\tau}$ and $K_1(\beta)e^{\beta\tau}/\pi$, respectively. In such a case they may be considered as the evolution of a distribution, which satisfies the equation $\partial_t F_i(\xi, \tau) = -[(1 - \partial_\xi^2)^2 - 1]F_i(\xi, \tau), i = 1, 2$ and exhibits the correct nonrelativistic limit in the sense that $-[(1 - \partial_\xi^2)^2 - 1] \approx \frac{1}{2} \partial_\xi^2 - \frac{1}{8} \partial_\xi^4 + \ldots$.

3. The relativistic Schrödinger equation (Salpeter equation)

The example of relativistic evolution, which will be considered here, is the free relativistic Schrödinger equation, called also the Salpeter equation, which has the form

$$i\hbar \partial_t \Psi = c\sqrt{\beta^2 + m^2c^2}\Psi.$$  (21)

Written in dimensionless variables, it reads

$$i\partial_t \Psi(q, \tau) = \sqrt{1 + \hat{q}^2} \Psi(q, \tau),$$  (22)

where $\tau = ct/\lambda_c$, $\hat{q} = \lambda_c\hat{p}/\hbar$, and $\lambda_c = \hbar/(mc)$, and then

$$i\partial_\tau \Psi(\xi, \tau) = \sqrt{1 - \partial_\xi^2} \Psi(\xi, \tau),$$

$$\Psi(\xi, \tau) \big|_{\tau=0} = \Psi_0(\xi)$$  (23)

with $\hat{q} = i\partial_\xi$, $\xi = x/\lambda_c$, originating from the standard quantization rule $\hat{p} = i\partial_\xi$.

In analogy to the equation (6) we can express the formal solution of equation (23) in terms of a Gaussian wave function $\Psi(\xi, \tau)$ and the Lévy distribution $g_{1/2}(\eta)$

$$\Psi(\xi, \tau) = e^{-i\tau\sqrt{1-\partial_\xi^2}} \Psi_0(\xi)$$

$$= \int_0^\infty g_{1/2}(\eta)e^{\eta^2}D(\xi, -\eta\tau^2)d\eta.$$  (24)

Taking $D(\xi, -\eta\tau^2)$ in the Fourier space (see equation (10) with appropriately changed variables), we can show that equation (24) is equivalent to the solution of equation (22). It is seen from

$$\Psi(\xi, \tau) = \int_0^\infty g_{1/2}(\eta)e^{\eta^2}\left[\int_{-\infty}^{\infty} e^{i\eta^2\xi^2} D(q) \frac{dq}{\sqrt{2\pi}}\right]d\eta$$

$$= \int_{-\infty}^{\infty} e^{\eta^2}\left[\int_0^\infty g_{1/2}(\eta)e^{\eta^2(1+\eta^2)}d\eta\right] \Psi_0(q) \frac{dq}{\sqrt{2\pi}}$$

$$= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{i\eta^2} e^{-i\tau\eta^2} \Psi_0(q) dq.$$  (25)

The inverse Fourier transform of equation (25) defines $\tilde{\Psi}(q, \tau)$, which is a formal solution of equation (22).

Let us now solve the relativistic Schrödinger equation for an initial Gaussian packet given in equation (12), i.e., for the initial condition $e^{-i\beta^2}$ taken in equation (25). Thus we have

$$\Psi(\beta; \xi, \tau) = \frac{N}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{-\rho^2 - i\tau(1+\rho^2)} d\rho$$  (26)

with the normalization constant $N = [2\beta/(\pi\lambda_c^2)]^{1/4}$ calculated in $L^2[\mathbb{C}, d\xi]$ space.
A time-dependent solution of equation (23) obtained along the same lines as before for
\[ \Psi_0(q) = \frac{1}{2K_0(2\beta)} e^{-\beta \sqrt{1+q^2}} \] reconstructs the result presented in [18]; see equation (4.18) there. Namely, we get
\[ \Psi_0(\beta; \xi, \tau) = \beta e^{\frac{i \xi}{\beta}} K_1\left(\sqrt{\beta^2 + \xi^2}\right) \frac{K_1\left(\sqrt{(\beta + i \tau)^2 + \xi^2}\right)}{\sqrt{\beta^2 + \xi^2}}. \] (27)

A comparison between nonrelativistic and relativistic solutions reveals interesting behaviour, illustrated in figure 1. The snapshots, taken at different times, show that the wave packet spreading is 'slower' for the relativistic case than for its nonrelativistic counterpart. Furthermore, the relativistic solutions reveal, at longer times, the wave-function deformation determining their departure from the Gaussian-like initial form.

Numerical calculations of \( \Psi_0(\beta; \xi, \tau) \) done for various values of \( \beta \) and \( \tau \) show that there exists a non-trivial relation between \( \beta \) and \( \tau \), which implies an appearance, or non-appearance, of two peaks. The two-peaks evolution pattern appears for reasonable values of \( \tau \) if \( \beta \) is small enough; if not, the evolution follows the standard spreading. As a quantity indicating what is going on for a given \( \tau \), the value of the second derivative \( D_2^{(i)}(\beta, \tau) \) can be used. For \( \tau = 0 \) it is evident that \( D_2^{(i)}(\beta, 0) < 0 \) for both equations (26) and (27). But if \( \tau \) grows, it may happen that \( D_2^{(i)}(\beta, \tau) \) changes its sign, which means that the point \( \xi = 0 \) becomes the local minimum. This depends on the parameter \( \beta \), measuring the width of the initial wave packet, i.e., the initial localization, as well. From numerical calculation of \( D_2^{(i)}(\beta, \tau) \) for fixed \( \tau \), we find there exist \( \beta_0 \) such that for \( \beta > \beta_0 \) we observe two peaks. For instance, \( D_2^{(1)}(\beta, 7) \approx 0 \) for \( \beta_0 = 0.689 \), and \( D_2^{(2)}(\beta, 7) \approx 0 \) for \( \beta_0 = 1.411 \). All this allows us to make a conjecture: The two-peaks evolution pattern is possible only if the initial wave packet is localized sufficiently strongly, \( \beta < \beta_0 \), and the 'critical' localization \( \beta_0 \) is related to the Compton wavelength of the particle.

To justify the previous assumption we will consider another solution of equation (23), which is obtained for the initial condition defined in the Fourier space as follows
\[ \Psi_0^{(3)}(\beta; q) = \frac{N |q|}{\sqrt{1+q^2}} e^{-\beta \sqrt{1+q^2}}. \] (28)

After a simple calculation, applying equation (2.5.39.6) on p. 456 of [26], equation (25), with the initial condition equation (28), reads
\[ \Psi_3(\beta; \xi, \tau) \equiv N \left( \sqrt{a^2 + \xi^2} \right)^{1/2} e^{-\sqrt{a^2 + \xi^2}}. \] (29)

where \( a = \beta + i \tau \). The normalization constant \( N \) calculated in \( L^2 \) is given as
\[ N = \sqrt{\pi/\lambda_0} [K_0(2\beta) + \frac{\pi}{2} - \pi K_0(2\beta) L_{-1}(2\beta) - \pi \beta K_0(2\beta) L_0(2\beta)]^{-1/2}, \] where \( L_n(u) \) is the modified Struve function. Now we can find the exact relation between \( \beta \) and \( \tau \) for which two peaks appear. Namely,
Figure 1. (a) Comparison, for $\beta = 0.2$ and $\tau = 0.2$, of the relativistic solutions $|\Psi(\beta; \xi, \tau)|, i = 1, 2$ (equations (27) and (26) for lines I and II, respectively) and nonrelativistic solutions $F_i(\beta; \xi, \tau), i = 1, 2$ (equations (19) and (14) for lines III and IV, respectively). (b) The same as in (a) but for $\beta = 0.2$ and $\tau = 2$. 
from which it is immediately seen that
\[ D^2(\beta, \tau) = \frac{N}{2\sqrt{\pi}} \left( \beta^2 + \tau^2 \right)^{3/2} \left( \tau^2 (3 - 4\beta) - \beta^2 (3 + 4\beta) \right). \] (30)

The role of the $\beta$ parameter appearing in equation (31) is not trivial. It is an arbitrary constant associated with the intrinsic nonlocality of the problem under study, and it defines the scale of delocalization of the particle within its Compton wavelength.

Figure 2. The three various behaviours of $|\Psi(\beta; \xi, \tau)|^2$ for $\beta = 1/4$ and $\tau = 0.2$ (blue line), $\tau = \tau_c = \sqrt{2}/4$ (red line), and $\tau = 1$ (green line).

The waves packets whose evolution has been reported in figures 1 are initially localized within a Compton wavelength. In the case of an electron-neutrino assuming a mass of 2 eV, we are talking about an uncertainty in position of the order of a few microns. The case of massless particles could be obtained from equation (31) by going to the limit $\lambda_c \to \infty$. However, keeping such a limit without any further caution does not provide physical results. Equation (31) yields indeed a completely delocalized function unless we assume in such a
limit \( \beta \lambda_c \) remains a constant. Since \( \beta \) is an arbitrary scale factor, we can define \( b = \beta \lambda_c \). By such a redefinition, the dependence on the Compton wavelength is a physical dependence and no longer a scale factor. We can keep the limit safely and, using the well-known asymptotic property \( K_1(z) |_{z \to \infty} \approx 1/z \), we obtain the same result as quoted in [27]:

\[
\Psi_2(x, t) = \frac{2\hbar}{\pi} \frac{b + ic \ell}{\sqrt{(b + ic \lambda)^2 + x^2}}.
\] (32)

It is now worthwhile to recast equation (32) in the form

\[
\Psi(x, t) = \frac{b}{2\pi} \left[ \frac{1}{b + ic(x + ct)} + \frac{1}{b - ic(x - ct)} \right]
\] (33)

and note that it can be viewed as the solution of the equation [27]

\[
i\hbar \partial_t \Psi = c\sqrt{-\nabla^2_x} \Psi,
\] (34)

\[
\Psi(x, 0) = \frac{b}{2\pi} \left( \frac{1}{b + ix} + \frac{1}{b - ix} \right).
\] (35)

which is the massless limit of the relativistic Schrödinger equation. Equation (34) does not depend explicitly on the Planck constant; such a dependence is, however, implicitly contained in the arbitrary constant \( b \). We find also that the r.m.s. value of the particle position is indeed provided by \( \sqrt{\langle x(t)^2 \rangle} = \sqrt{b^2 + (ct)^2} \), which is consistent with the constraint \( \sqrt{\langle p(t)^2 \rangle} > \hbar / (2\sqrt{b^2 + (ct)^2}) \) (for additional comments see [15]). The emergence of two peaks, when the particle is localized according to the previous discussion, is a distinctive feature of the nonlocal nature inherent to the relativistic Schrödinger equation. This aspect of the problem may acquire a deeper physical nature when studied in the presence of a potential, as will be discussed in a forthcoming paper.

4. The Dirac factorization and the one-dimensional quantum relativistic equation

This section is devoted to the Dirac factorization of the one-dimensional relativistic Schrödinger equation. This problem has recently raised a significant amount of interest, because it may provide an important toy model to benchmark a genuine quantum relativistic effect, like the Zitterbewegung, in the so-called quantum simulation of a Dirac equation [28]. The equation can be written in the two-component form [29] as follows

\[
i\hbar \partial_t \Psi = c\hat{a}\hat{p} \Psi + \hat{p}mc^2 \Psi.
\] (36)

In the dimensionless coordinates \( \tau \) and \( \xi \) it can be expressed by

\[
i\hat{a} \partial_\tau \Psi = -i\hat{a} \partial_\xi \Psi + \hat{p} \Psi,
\] (37)

where \( \hat{a} \) and \( \hat{p} \) are Pauli-like matrices satisfying the identities

\[
\hat{a}^2 = \hat{p}^2 = \hat{1} = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}, \quad \hat{a} \hat{p} + \hat{p} \hat{a} = 0
\] (38)
and numerically given as
\[
\hat{a} = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \quad \text{and} \quad \hat{b} = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}.
\]

We have denoted by \( \Psi = \begin{pmatrix} \psi \\ \bar{\psi} \end{pmatrix} \) the two-component vector, replacing the wave function \( \Psi \).

The vector \( \Psi \) should not be confused with a spinor, and equation (37) is not the relativistic counterpart of the Pauli equation, since we are not dealing with a problem including the spin degrees of freedom. The reason for using two-dimensional matrices is that we want equation (37) to describe negative and positive energy states.

The solution of equation (37) can be formally obtained, using the evolution operator method, as
\[
\Psi(\xi, \tau) = \tilde{U}(\tau) \Psi(\xi, 0)
\]
with the evolution operator written in the form of the \( 2 \times 2 \) matrix [30]
\[
\tilde{U}(\tau) = \exp \left( \tau \begin{pmatrix} i & \partial_\xi \\ \partial_\xi & -i \end{pmatrix} \right) = \begin{pmatrix} i\tilde{A}(\tau) + \tilde{B}(\tau) & \tilde{A}(\tau)\partial_\xi \\ \tilde{A}(\tau)\partial_\xi & -i\tilde{A}(\tau) + \tilde{B}(\tau) \end{pmatrix},
\]
where \( \tilde{A}(\tau) = \sin(\sqrt{\Delta} \tau)/\sqrt{\Delta}, \tilde{B}(\tau) = \cos(\sqrt{\Delta} \tau), \) and \( \Delta = 1 - \partial_\xi^2 \). This evolution operator is unitary and, therefore, the Dirac wave function remains normalized at any time, namely
\[
\int_{-\infty}^{\infty} \left| \Psi(\xi, \tau) \right|^2 d\xi = \int_{-\infty}^{\infty} \left( \left| \psi_+(\xi, \tau) \right|^2 + \left| \psi_-(\xi, \tau) \right|^2 \right) d\xi = 1.
\]

The use of the Fourier transform method allows one to obtain the solution of the free-particle Dirac equation: from equation (39) with \( \tilde{U}(\tau) \) given in equation (40), we get the evolution of the two components in the form
\[
\psi_+(\xi, \tau) = \int_{-\infty}^{\infty} \left\{ \frac{\sin\left(\sqrt{1 + k^2} \tau\right)}{\sqrt{1 + k^2}} \left[ \tilde{\psi}_+(k, 0) + k\tilde{\psi}_-(k, 0) \right] \right\} e^{i k \xi} \frac{dk}{\sqrt{2\pi}},
\]
\[
\psi_-(\xi, \tau) = \int_{-\infty}^{\infty} \left\{ \frac{\sin\left(\sqrt{1 + k^2} \tau\right)}{\sqrt{1 + k^2}} \left[ k\tilde{\psi}_+(k, 0) - \tilde{\psi}_-(k, 0) \right] \right\} e^{i k \xi} \frac{dk}{\sqrt{2\pi}},
\]
where \( \tilde{\psi}_+(k, 0) \) are the Fourier transforms of \( \psi_+(\xi, 0) \). In figure 3 we present the snapshots of the time evolution for \( \left| \Psi(\xi, \tau) \right|^2 \), calculated using equations (41)–(43), with the initial wave packet being a Gaussian containing both positive and negative energy components.

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\[ \Psi_1(k, 0) = e^{-\beta k^2} \begin{pmatrix} 1 \\ 1 \end{pmatrix} \] (44)

The presence of the negative energy term yields the behaviour with some features not exhibited by the Salpeter equation. These effects are due to the superposition between the two states, which induces also the characteristic trembling motion of the packet average position and yields the 'Zitterbewegung' [31], namely, an oscillating motion around a linear trajectory. We remark that using equation (42) with the initial condition possessing only the positive energy term, which is given by a Gaussian function, we get the evolution, which is not dissimilar from the one described with the relativistic Schrödinger equation and from the behaviour obtained in [15], using different solution techniques.

Following [27] we can also find an exact solution for the one-dimensional massive Dirac equation with the initial condition

\[ \Psi_2(k, 0) = e^{-\beta \sqrt{1+k^2}} \begin{pmatrix} \psi_e(k, 0) \\ \psi_\mu(k, 0) \end{pmatrix} \] (45)

According to equations (18) and (42), and taking the integral representation of \((1 + k^2)^{-1/2}\), i.e., \(\int_0^\infty e^{-\sqrt{1+k^2}} ds\), we obtain

\[ \psi_e(\xi, \tau) = \frac{1}{\sqrt{2\pi}} \left[ K_0(c^*) + \frac{\beta - i\tau}{c^*} K_1(c^*) \right] + \frac{1}{\sqrt{2\pi}} \left[ -K_0(c) + \frac{\beta + i\tau}{c} K_1(c) \right] \] (46)

and

\[ \psi_\mu(\xi, \tau) = \frac{1}{\sqrt{2\pi}} \left[ \frac{\xi}{c} K_1(c) - \frac{\xi}{c^*} K_1(c^*) \right], \] (47)

where \(c = \sqrt{(\beta + i\tau)^2 + \xi^2}\) and \(c^*\) is its complex conjugation.

5. One-dimensional Klein–Gordon equation

The free-particle Klein-Gordon (FP-KG) equation [1] has played a fundamental role in the description of relativistic spin 0 particles. It is well known that according to the Feshback–Villars transformation [32], FP-KG is transformed into a two-component Schrödinger-like equation, which exhibits a Hamiltonian with unconventional properties. In this section we treat this transformation from a different perspective. We will look for a formal solution of the FP-KG equation by treating it as a D'Alembert equation, and next we will show that the solution of the one-dimensional FP-KG can be expressed in terms of those of the Dirac equation.

Let FP-KG be written in the dimensionless coordinates \(\xi\) and \(\tau\) introduced in section 3 as follows

\[ \partial_\xi^2 \Psi(\xi, \tau) = -\left(1 - \partial_\xi^2\right) \Psi(\xi, \tau) \] (48)

with the initial conditions

\[ \Psi(\xi, 0) = \Psi_0(\xi), \quad \left[ \partial_\xi \Psi(\xi, \tau) \right]_{\tau=0} = \Psi_1(\xi). \] (49)

Using the formal procedure presented in the previous sections to solve equations (48) and (49), we can cast our problem in the form of an ordinary second-order equation with constant coefficients, namely,
which resembles that of a harmonic oscillator. Therefore, we postulate that the solution of FP-KG can be written in the form

\[ \psi_\xi (\tau, \xi) = C_1 e^{i\hat{D}\tau} \psi_\xi (\xi) + C_2 e^{-i\hat{D}\tau} \psi_\xi (\xi), \]  

where \( C_1, C_2 \) are linked to the initial conditions of equation (49) by the identities

\[ C_1 + C_2 = \psi_\xi (\xi), \quad C_1 - C_2 = -i\hat{D}^{-1}\psi_\xi (\xi). \]  

Thus, inserting the explicit formulas on \( C_1 \) and \( C_2 \) into equation (51) and making simple symbolic manipulations, we get

\[ \psi_\xi (\tau, \xi) = \frac{1}{2} \left[ \psi_{K-S}(\tau, \xi) \psi_{K-S}(\xi, -\tau) + \psi_{K-S}(\xi, -\tau) \psi_{K-S}(\tau, \xi) \right] \]

\[ \quad = \cos (\tau D) \psi_0 (\xi) + \sin (\tau D) \hat{D}^{-1}\psi_1 (\xi). \]  

The function \( i\psi_{K-S}(\xi, \tau) \) is the solution of the (forward) free relativistic Schrödinger particle with initial condition \( \psi_0 (\xi) - i\hat{D}^{-1}\psi_1 (\xi) \), and \( i\psi_{K-S}(\xi, -\tau) \) is its backward counterpart, with initial condition \( \psi_0 (\xi) + i\hat{D}^{-1}\psi_1 (\xi) \).

The evolution of the solution of the FP-KG equation is given in figure 4, where we have used an initially Macdonald distribution, namely, for \( \psi_0 (k) = e^{i\beta\sqrt{1+k^2}} \) and \( \hat{D}(\xi) = 0 \). Then, equation (54) gives

\[ \psi_{K-G}(\xi, \tau) = \left\{ \frac{\beta - i\tau}{\sqrt{2\pi c}} K_1(c^*) + \frac{\beta + i\tau}{\sqrt{2\pi c}} K_1(c) \right\}, \]  

where \( c \) and its complex conjugation \( c^* \) are defined below equation (47). This is the effect of the interference with the backward time solution, which yields a fairly rich evolution, according to which the wave function undergoes a free propagation. The use of a Gaussian does not provide any significant difference.

The evolution operator associated with equation (48), related to the two-component Schrödinger-like equation, should be written as the \( 2 \times 2 \) matrix:

\[ \begin{pmatrix} a_1 & a_2 \\ a_3 & a_4 \end{pmatrix} = \begin{pmatrix} 1 & i \sin(\tau) \\ -i \cos(\tau) & 1 \end{pmatrix}. \]
\[ \hat{U}_{KG}(\tau) = \begin{pmatrix} \cos \left( \tau \hat{D} \right) & \sin \left( \tau \hat{D} \right) \hat{D}^{-1} \\ -\sin \left( \tau \hat{D} \right) & \cos \left( \tau \hat{D} \right) \end{pmatrix}, \tag{56} \]

whose determinant is equal to one. The inverse operator of \( \hat{U}_{KG}(\tau) \) exists, and it fulfills the relation \( [\hat{U}_{KG}(\tau)]^{-1} = \hat{U}_{KG}(-\tau) \). In consequence, equation (48) can be rewritten as

\[ \begin{align*}
\Psi_0(\xi, \tau) \\
\Psi_1(\xi, \tau)
\end{align*} = \hat{U}_{KG} \begin{align*}
\Psi_0(\xi) \\
\Psi_1(\xi)
\end{align*}, \tag{57} \]

where the equation on the upper component \( \Psi_0(\xi, \tau) \) is given in equation (54), whereas \( \Psi_i(\xi, \tau) = \partial t \Psi_{KG}(\xi, \tau) \). Moreover, the conserved quantity associated with such an evolution operator is

\[ \int_{-\infty}^{\infty} (|\Psi_0(\xi)|^2 + |\Psi_1(\xi)|^2) d\xi, \tag{59} \]

which cannot be understood in terms of probability; rather, it has been constructed in analogy with the energy of a harmonic oscillator.

Let us observe that in Fourier space equation (54) reads

\[ \tilde{\Psi}_{KG}(k, \tau) = \cos [\tau \omega(k)] \tilde{\Psi}_0(k) + \frac{\sin [\tau \omega(k)]}{\omega(k)} \tilde{\Psi}_1(k), \tag{58} \]

where \( \omega(k) = \sqrt{1 + k^2} \) and \( \tilde{\Psi}_i(k), i = 0, 1 \) are the initial condition written in the Fourier space. The analogy with a harmonic oscillator is made more clear by noting that in the Fourier space, the quantity

\[ \int_{-\infty}^{\infty} \left[ \partial_t \Psi_{KG}(k, \tau) \right]^2 + \left[ \omega(k) \Psi_{KG}(k, \tau) \right]^2 dk \]

is a constant of motion.

Finally, we point out that it is also possible express the solution of our problem in terms of the free-particle Dirac solution by setting

\[ \hat{D} = -i\alpha \partial_\xi + \beta. \tag{60} \]
6. Quantum relativistic particles in a linear potential

6.1. The Salpeter equation with a linear potential

Let us consider the Salpeter equation with a potential function \( V(\xi) \)

\[
\partial_{\xi} \Psi(\xi, \tau) = \sqrt{1 - \partial_{\xi}^2} \Psi(\xi, \tau) + \mu_0 V(\xi),
\]

\( \Psi(\xi, 0) = \Psi_0(\xi), \)  \hspace{2cm} (61)

which in the Fourier space reads

\[
i \partial_{\tau} \tilde{\Psi}(k, \tau) = \sqrt{1 + k^2} \tilde{\Psi}(k, \tau) + \mu_0 \tilde{V}(k, \tau),
\]

where \( \tilde{V} = V(i\partial_\xi) \) and \( \mu_0 \) is a dimensionless constant. We note that, in a general case, the presence of an operator \( \tilde{V} \), which doesn’t commute with the kinetic term, hampers the possibility of a straightforward analytical solution for equation (62). It can, however, be obtained for \( V(\xi) \) being a linear function, for which a physically meaningful example is when a relativistic, charged particle interacts with a static constant electric field. In this case, equation (62) can be rewritten in the form

\[
i \partial_{\tau} \tilde{\Psi}(k, \tau) = \sqrt{1 + k^2} \tilde{\Psi}(k, \tau) + i\mu_0 \partial_{\tau} \tilde{\Psi}(k, \tau).
\]

To solve equation (63), first we eliminate the derivative by the use of the transformation

\[
\tilde{\Psi}(k, \tau) = e^{i\mu_0 \partial_\tau} \tilde{\Phi}(k, \tau),
\]

which yields a modified equation:

\[
i \partial_{\tau} \tilde{\Phi}(k, \tau) = \tilde{K} \tilde{\Phi}(k, \tau)
\]

\( \tilde{K}(k, \tau) = e^{-i\mu_0 \partial_\tau} \sqrt{1 + k^2} e^{i\mu_0 \partial_\tau}. \)  \hspace{2cm} (65)

The use of the procedure touched upon in [16] allows us to cast its formal solution in the form

\[
\tilde{\Phi}(k, \tau) = \exp\left[-i \int_0^\tau \tilde{K}(k, \tau') d\tau' \right] \tilde{\Phi}_0(k)
\]

where \( \tilde{\Phi}_0(k) \equiv \tilde{\Psi}_0(k) \). The solution can finally be obtained in the form

\[
\tilde{\Psi}(k, \tau) = e^{i\mu_0 \partial_\tau} \tilde{\Phi}(k, \tau) = \tilde{\Phi}(k + \mu_0 \tau, \tau)
\]

\[
= e^{-i \int_0^\tau \frac{k + \mu_0 \tau'}{\sqrt{1 + (k + \mu_0 \tau')^2}} d\tau'} \tilde{\Phi}_0(k + \mu_0 \tau).
\]

In figure 5 we report a comparison between the solutions for the free-particle case and that of equation (61) with a linear potential put in. In both cases we have assumed as initial wave function the function given in equation (16) with \( \beta = 1 \). It is important to emphasize that the presence of the potential provides a suppression of the emergence of the bi-modal behaviour by inducing a kind of external localization.

6.2. The Dirac equation with a linear potential

The one-dimensional Dirac equation with a linear potential has the form

\[
i \partial_{\tau} \Psi = -i\partial_{x} \Psi + \beta \Psi + \xi \partial_{_{\parallel}} \Psi,
\]

\( \Psi_1 \) represents the wave function of the particle.

In figure 5 we report a comparison between the solutions for the free-particle case and that of equation (61) with a linear potential put in. In both cases we have assumed as initial wave function the function given in equation (16) with \( \beta = 1 \). It is important to emphasize that the presence of the potential provides a suppression of the emergence of the bi-modal behaviour by inducing a kind of external localization.
where $\mu_0 = \mu_0 \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$. The solution of the problem is slightly more complicated than in the case of the Salpeter equation, since the Hamiltonian is now expressed as a linear combination of Pauli matrices. The formal solution of equation (69) can be given by using the operational method with the evolution operator defined as

$$\hat{U} = e^{i\hat{h} \tau + i\hat{\rho}_0 + i\hat{\rho}_1}.$$  

(70)

Although the problem is amenable to an exact treatment, we use the Zassenhaus formula at the second order in the commutator expansion [33] and write the operator $\hat{U}$ in the following disentangled form

Figure 5. (a) Comparison between evolution of the modulus of a wave function in the presence of linear potential (I: blue curve; equation (68) with $\Psi(k)$ given in equation (16)) and evolution of the wave function of a free-particle (II: red curve; equation (27)) at $\tau = 4.3$; (b) the same as (a) but $\tau = 2.3$. 

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\[ \hat{U} = e^{i \hat{\mathcal{H}}_0 t + i \delta \hat{e} \hat{\mathcal{H}}_0 \frac{\partial}{\partial \xi} \hat{\mathcal{H}}_0 \frac{\partial}{\partial \xi}}. \] (71)

which is valid at the order \( o(t^3) \). The repeated application of the evolution operator (71) on the initial state yields the solution in the form of the following recursion

\[ \Psi_n(\xi, \tau) = \begin{pmatrix} e^{i(\xi + \mu_\xi \zeta)} & 0 \\ 0 & e^{-i(\xi - \mu_\xi \zeta)} \end{pmatrix} \begin{pmatrix} \cosh(\tau \sigma_\xi) & \sinh(\tau \sigma_\xi) \\ \sinh(\tau \sigma_\xi) & \cosh(\tau \sigma_\xi) \end{pmatrix} \Psi_{n-1}(\xi, \tau), \] (72)

where

\[ \Psi_0(\xi, \tau) = \begin{pmatrix} \psi_+^0(\xi) \\ \psi_-^0(\xi) \end{pmatrix}, \] (73)

\[ \Psi_j(\xi, \tau) = \begin{pmatrix} \psi_+^j(\xi) + \psi_-^j(\xi) e^{i(\xi - \mu_\xi \zeta)} \\ \psi_-^j(\xi) + \psi_+^j(\xi) e^{-i(\xi + \mu_\xi \zeta)} \end{pmatrix}, \] (74)

with

\[ \psi_\pm^j(\xi) = \frac{1}{2} \left[ \psi_\pm(\xi + \tau) \pm \psi_\pm(\xi - \tau) \right]. \] (75)

The previous iteration can be extended to any type of potential, and specific relevant examples will be discussed in a paper specially dedicated to this problem.

7. Conclusion

In this paper we have presented a number of issues of a practical nature to deal with problems concerning relativistic equations and other fractional evolution problems often encountered in physics. We have left open many points, especially the link of the present treatment with previous and well-established methods: in particular, the use of the Green function technique for the treatment of the Klein–Gordon equation or the method associated with the use of the Wightman solution [34]. In this last case the solution of the free-particle Klein–Gordon is written in terms of Wightman (and anti-Wightman as well) functions expressed as combinations of Hankel and Macdonald functions of the first kind. The relevant link with our formalism and with other types of solutions (Pauli–Jordan, Feynman, Stueckelberg) will be the topic of a forthcoming investigation. We have touched on many aspects of a relevant formalism, and we have fixed the main mathematical apparatus, which can be exploited to discuss a number of problems in relativistic nonlocal quantum mechanics.

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