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Scattering in strong electromagnetic fields: Transverse size effects in time-dependent basis light-front quantization

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The framework of “time-dependent basis light-front quantization” (tBLFQ) offers a nonperturbative approach to scattering problems in external fields, based on Fock space truncation. Here we extend tBLFQ to include spatio-temporal field inhomogeneities in multiple spacetime directions. This extension is necessary for the proper modeling of e.g. intense laser fields. We focus on the example of nonlinear Compton scattering of an electron on an axicon-type laser, with an emphasis on the transverse structure of the beam. We analyze the impact of field intensity and particle energy, as well as basis truncation effects, on the radiation spectrum of the scattered electron.

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I. INTRODUCTION

The theoretical predictions of perturbative QED, such as electron $g-2$ [1,2] and the Lamb shift [3], have been experimentally verified to extreme precision. However, there are many problems for which standard perturbative methods are insufficient, such as bound states [4], or break down completely, such as strongly coupled systems. Thus other methods are required to access *nonperturbative* physics.

In the context of external field problems, perhaps the most famous nonperturbative effect is the spontaneous conversion of a sufficiently strong electric field into electron-positron pairs, or Schwinger pair production, the probability of which is a nonperturbative function of the coupling [5]. It may be possible to observe Schwinger pair creation using the ultra-intense electromagnetic fields of future lasers [6,7], and there is currently an international effort to develop both the theoretical and experimental tools necessary to investigate this and other quantum effects in strong (external) fields, such as vacuum birefringence; for a review see [8].

When investigating nonperturbative phenomena, exactly solvable systems offer intuition, but one often turns to numerical schemes which do not rely on perturbative approximations. The lattice is a well-known example, and real-time lattice techniques can now be used to analyze e.g. string breaking and the Schwinger effect [9]. Within

laser-plasma physics, a common tool is the simulation of processes using particle-in-cell (PIC) codes, in which Monte Carlo routines based on QED calculations are included to simulate (tree level) quantum effects; for a review see [10]. The advantages of such approaches are that extremely complicated electron-positron-photon interactions can be simulated, with arbitrarily complex laser fields. The disadvantage, though, is that such methods are firmly rooted in classical physics, and the approximations behind the inclusion of quantum effects are known to break down at both low [11–13] and high energy [14,15]. Here we consider a different approach; it is not our goal to compete with e.g. PIC schemes, but to complement them with an alternative, and fully quantum mechanical, framework.

Any numerical implementation of QFT requires a cutoff in order to render problems finite-dimensional. Because the set of all products of one-particle states gives a possible basis of Hilbert space [16], one can try to solve problems using a finite truncation of this basis, corresponding to a truncation in particle number [17,18]. This is the main idea of basis light-front quantization (BLFQ) [19]; one solves the Schrödinger equation in a truncated Fock space, rather than a perturbative expansion. Thus BLFQ is a nonperturbative approach. An extension of this framework which is suitable for investigating e.g. QFT scattering processes in *time-dependent* background fields is time-dependent basis light-front quantization (tBLFQ) [20,21]. The idea of tBLFQ can also be used in nuclear physics [22].

Previously the emission of photons from an electron scattering off an intense background field (called nonlinear

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Compton scattering) [23] has been studied in tBLFQ, but in the approximation that the background has essentially univariate spacetime dependence [20,24]. However, intense laser fields for example are strongly spatially focused, and so further progress in high-intensity laser-matter interactions requires the consideration of complex background field geometries. Motivated by this, and by a desire to improve the overall scope of tBLFQ, we here extend the framework to include effects due to multidimensional field inhomogeneities.

This paper is organized as follows. We review the main ideas of BLFQ and tBLFQ in Sec. II. In Sec. III we use tBLFQ to study the simpler system of a single electron in a static, but position-dependent, background field (what will be the “center-of-mass” motion of the system). In Sec. IV we reintroduce the dynamical photon fields and study photon emission from the electron; observables are defined and their time-evolution is studied. In Sec. V, we add a time dependence to the background. We present our conclusions and outlook in Sec. VI. The appendixes include conventions and other details behind the calculations in the text. A comparison of our results with perturbation theory, in the limit of weak background fields, is also included in the appendixes.

II. BACKGROUND

In the light-front formalism, field theories are quantized at fixed light-front time $x^+ = x^0 + x^3$, following [25], rather than at fixed x^0 as in the equal-time formalism; see [26,27] for reviews. The remaining coordinates are $x^- = x^0 - x^3$ (longitudinal) and $x^\perp = \{x^1, x^2\}$ (transverse).

Due to the larger number of *kinematic* Lorentz generators, over *dynamical* generators, in light-front coordinates, wave functions in one frame can be more easily mapped to those in other frames. Furthermore, light-front kinematics forbids massive particles from contributing to the vacuum, therefore making the trivial Fock vacuum the vacuum of the full theory. (This holds only if one neglects zero-mode contributions, which it is not always safe to do [28–35].)

One consequence of this is that the physical cases of interest here are easily expandable in simple Fock modes. For example, if the vacuum received its most significant contribution from some higher Fock sector, take for example $|e^+e^- \rangle$, then the Fock expansion of the physical electron would “begin” with $|e \rangle_{\text{phys}} \sim |eee^+ \rangle$. The triviality of the vacuum means though that the first contributions take the form $|e \rangle_{\text{phys}} = a|e \rangle + b|e\gamma \rangle$. It is the simplicity of the Fock expansion that makes the Hamiltonian formalism feasible in QFT.

A. Time-evolution in light-front field theory

In the Hamiltonian formalism of QFT, the time-evolution of a system is governed by the Schrödinger equation, which in light-front quantization takes the form

$$i \frac{\partial}{\partial x^+} |\Psi; x^+ \rangle_S = \frac{1}{2} P_S^-(x^+) |\Psi; x^+ \rangle_S, \quad (1)$$

in which the Schrödinger picture Hamiltonian P_S^- contains two parts,

$$P_S^-(x^+) = P_{\text{QED}}^- + V_S(x^+), \quad (2)$$

where P_{QED}^- is the full light-front Hamiltonian of, here, QED, and V_S consists of additional interaction terms introduced by a background field, to be specified explicitly below. If the effects introduced by this background are the primary source of interest, then it is convenient to work in an interaction picture in which the “free” states are eigenstates of the full QED Hamiltonian P_{QED}^- , and where the only nontrivial time-evolution is induced by the new interaction V_S . It is of course not possible to do this exactly for QED, and hence a numerical approximation will be introduced below. In principle, though, in this interaction picture, the Schrödinger equation becomes

$$i \frac{\partial}{\partial x^+} |\Psi; x^+ \rangle_I = \frac{1}{2} V_I(x^+) |\Psi; x^+ \rangle_I, \quad (3)$$

the formal solution to which is

$$|\Psi; x^+ \rangle_I = \mathcal{T}_+ \exp \left(-\frac{i}{2} \int_0^{x^+} dx'^+ V_I(x'^+) \right) |\Psi; 0 \rangle_I. \quad (4)$$

States and operators in the two pictures are as usual related by

$$\begin{aligned} |\Psi; x^+ \rangle_I &= e^{\frac{i}{2} P_{\text{QED}}^- x^+} |\Psi; x^+ \rangle_S, \\ A_I(x^+) &= e^{\frac{i}{2} P_{\text{QED}}^- x^+} A_S(x^+) e^{-\frac{i}{2} P_{\text{QED}}^- x^+}. \end{aligned} \quad (5)$$

In this interaction picture, we work in a basis of eigenstates of P_{QED}^- . This basis, called the “tBLFQ basis” $|\beta \rangle$ (more details will be presented in the next section), will simplify the operator exponentials in the definition of the operators in Eq. (5). In this basis the matrix elements of the interaction Hamiltonian V_I become

$$\langle \beta' | V_I | \beta \rangle = \langle \beta' | V_S | \beta \rangle \exp \left[\frac{i}{2} (P_{\beta'}^- - P_{\beta}^-) x^+ \right]. \quad (6)$$

It is then straightforward to evolve the quantum state according to Eq. (4), by decomposing the time-evolution operator into many small steps of light-front time x^+ with a step size δx^+

$$\begin{aligned} &\mathcal{T}_+ \exp \left(-\frac{i}{2} \int_0^{x^+} dx'^+ V_I(x'^+) \right) \\ &\rightarrow \left[1 - \frac{i}{2} V_I(x_n^+) \delta x^+ \right] \cdots \left[1 - \frac{i}{2} V_I(x_1^+) \delta x^+ \right]. \end{aligned} \quad (7)$$

This discretization is the Euler scheme; it is however not usually a good choice. The most significant shortcoming of this scheme is its poor stability (which means that the norm of the state vector $|\Psi; x^+\rangle$ changes as time evolves). Instead, we will adopt the second-order difference scheme MSD2, which has been proved to have better stability [36]. This scheme relates the state at $x^+ + \delta x^+$ to those at both x^+ and $x^+ - \delta x^+$ (rather than just that at x^+ as in the Euler scheme), thus:

$$\begin{aligned} |\Psi; x^+ + \delta x^+\rangle_I &= |\Psi; x^+ - \delta x^+\rangle_I + (e^{-iV_I \delta x^+/2} - e^{iV_I \delta x^+/2}) |\Psi; x^+\rangle_I \\ &\approx |\Psi; x^+ - \delta x^+\rangle_I - iV_I |\Psi; x^+\rangle_I. \end{aligned} \quad (8)$$

Once the quantum state at each time step is obtained, it is straightforward to construct observables from it.

B. Basis light-front quantization

As mentioned above, it is convenient to evolve the system in the tBLFQ basis, which by definition comprises the eigenstates of P_{QED}^- . In such a basis, the complicated exponentials of the operator in Eq. (5) reduce to phase factors, which greatly simplifies the computation. The eigenstates themselves are constructed in BLFQ, which is a Hamiltonian formalism incorporating the advantages of light-front dynamics. The main idea is to solve, numerically and in a Fock space truncation, the time-independent Schrödinger equation

$$P_{\text{QED}}^- |\beta\rangle = P_\beta^- |\beta\rangle. \quad (9)$$

BLFQ has the advantage of being able to solve bound state problems involving positronium [37] and hadron structures [38–53]. In this paper we follow previous work on the *physical* electron eigenstates in BLFQ [54]. A sector-dependent renormalization [55,56] of the electron mass is performed, improving over [20]. We review the main steps here; more details may be found in [20].

Equation (9) is infinite-dimensional. To reduce the equation to a finite-dimensional problem a truncation of the basis should be implemented. We are in this paper interested in transitions between the physical electron state and its excitations, which are scattering states, due to the interaction introduced by a background field. For simplicity we retain only the first two Fock sectors $|e\rangle$ and $|e\gamma\rangle$, which are enough to give a description of the photon emitted from the electron excited by the background field. In this approximation, the physical electron and photon states have the form, with $\#$ indicating some coefficients,

$$|e\rangle_{\text{phys}} = \#|e\rangle + \#|e\gamma\rangle, \quad |\gamma\rangle_{\text{phys}} = \#|\gamma\rangle. \quad (10)$$

This truncation of the Fock space implicitly assumes that higher Fock sectors give decreasing contributions to the

low-lying eigenstates, in which we are most interested; one motivation for this is the success of perturbation theory in QED, and further details of this approximation to the physical states will be discussed in Sec. IV C.

We now characterize the single-particle Fock sector states themselves. These carry four quantum numbers, ultimately corresponding to three momentum components and a spin or helicity. The first quantum number k labels the longitudinal momentum p^+ of the particle. We compactify the longitudinal direction x^- on a circle of length $2L$ and impose (anti)periodic boundary conditions on (fermions) bosons. The longitudinal momentum p^+ therefore takes the discrete values

$$p^+ = \frac{2\pi}{L} k, \quad (11)$$

where the dimensionless quantity $k = 1, 2, 3, \dots$ for bosons (neglecting the zero mode) and $k = \frac{1}{2}, \frac{3}{2}, \frac{5}{2}, \dots$ for fermions. The next two quantum numbers n and m are those of a 2D harmonic oscillator (2D-HO) in the transverse plane, of mass M and frequency Ω . These numbers therefore encode the transverse momenta. The 2D-HO eigenstates have the corresponding eigenvalues

$$E_{n,m} = (2n + |m| + 1)\Omega. \quad (12)$$

Note though that the only parameter of the 2D-HO that enters the eigenstates is the scale parameter $b := \sqrt{M\Omega}$; for details see Appendix C. The fourth and final quantum number λ is the light-front helicity of the particle. We write $\bar{\alpha} = \{k, n, m, \lambda\}$ as a shorthand for the four quantum numbers and define the corresponding single-particle state as $|\bar{\alpha}\rangle$ [20]. These states are chosen so as to preserve as many symmetries of the Hamiltonian as possible, and thus simplify calculations in BLFQ. Compared with a standard plane-wave basis, the 2D-HO basis preserves rotational symmetry in the transverse plane, even in the finitely truncated BLFQ basis. The shortcoming of the 2D-HO basis states is that they are not eigenstates of momentum and consequently it becomes difficult to separate the relative motion from the center-of-mass motion. We will however identify a method to resolve this, in Sec. IV D. For more details of the symmetries of the QED Hamiltonian and the basis, see [20,57]. To construct N -particle states $|\bar{\alpha}_N\rangle$, we simply take the direct product of single-particle states $|\bar{\alpha}_N\rangle = \otimes |\bar{\alpha}\rangle$. The working BLFQ basis $|\alpha\rangle$ is then the direct sum of all single- and multiparticle states retained within after the Fock truncation.

Even with the restriction in Fock number, the parameters in $\bar{\alpha}$ (aside from the helicity) are still unbounded and must also be truncated. To impose this, we introduce two parameters K_{total} and N_{max} with which to truncate the BLFQ basis in the longitudinal and transverse directions

respectively. For the longitudinal degrees of freedom, K_{total} is defined by

$$K_{\text{total}} = \sum_i k_i, \quad (13)$$

where the sum runs over all single-particle states. If the total longitudinal momentum is conserved, as is the case in this paper, then only a single K_{total} is enough to describe the system. Furthermore, because the light-front wave functions are functions of longitudinal momentum fractions $x_i = k_i/K_{\text{total}}$, rather than longitudinal momenta themselves, larger K_{total} only provides a finer description of the system.

In the transverse plane, we define the total transverse quantum number for a BLFQ basis state $|\alpha\rangle$ as

$$N_\alpha = \sum_i 2n_i + |m_i| + 1,$$

where the sum again runs over all particles in the state. All the retained basis states satisfy

$$N_\alpha \leq N_{\text{max}}. \quad (14)$$

Physically, N_{max} limits the total ‘‘energy’’ of 2D-HO states summed over all particles. N_{max} is specified globally across all Fock sectors to ensure that the transverse motion in different Fock sectors is truncated at the same energy. As a result, N_{max} determines both the UV and IR cutoffs for the transverse basis; for details, see [58,59].

At this stage we have a finite basis in which to work. Working in the BLFQ basis $|\alpha\rangle$ we can diagonalize the Hamiltonian in order to find the QED eigenstates $|\beta\rangle$ and eigenvalues P_β^- satisfying the time-independent Schrödinger equation (9). [It is straightforward to obtain the matrix elements in the BLFQ basis by using the commutation relations of the creation and annihilation operators of the Fock states; see the Appendix, Eq. (B9).] From there, matrix elements $\langle\beta'|V_I|\beta\rangle$ as in Eq. (6) are easily calculated in terms of the wave function $\langle\alpha|\beta\rangle$; for more details see Appendix D.

III. PARTICLE IN AN EXTERNAL FIELD

In this section, we consider the simplest limit of our system, namely single particles in an external field. In the language of tBLFQ, this means that we retain only the single-electron sector $|e\rangle$ in our calculations, neglecting the electron-photon sector $|e\gamma\rangle$ (and others). In doing so, we effectively turn off dynamical photon generation and absorption. These will be reinstated in Sec. IV.

A general background field generates four new interaction terms in the full light-front Hamiltonian of QED, given in full in Eq. (A2). Three of these vanish if we can represent the background by a potential having only a

single, longitudinal, component \mathcal{A}^- . We make this simplifying assumption from here on; note that there remains a large amount of freedom in the choice of the background field. Our choice is motivated by the desire to go beyond previous tBLFQ results [20] and include field dependence on the transverse coordinates x^\perp . Consider then an ‘‘axicon’’ laser beam propagating in the z -direction. The beam has a radially polarized electric field and an azimuthally polarized magnetic field, both transverse to the propagation direction. This implies the vanishing of the electric field on the symmetry, or z , axis [60,61]. Typically one takes the transverse fields to go like $\sim|x^\perp| \exp -|x^\perp|^2$. Axicon beams also have longitudinal fields (i.e. fields pointing in the z -direction) with essentially the same temporal (or x^+) dependence as the transverse fields, but which are suppressed relative to the transverse by a small focusing parameter. Thus a toy model of an axicon beam can be given by ignoring the longitudinal fields, and taking the potential to be

$$e\mathcal{A}^\mu(x) = \delta_-^\mu \frac{m_e a_0}{\sqrt{\pi}} \exp\left(-\frac{1}{2} b_l^2 x^\perp x^\perp\right) f(x^+), \quad (15)$$

in which $f(x^+)$ encodes the time dependence of the field; a_0 is a dimensionless amplitude, and b_l is a width. We will set $f \equiv 1$ to begin with, reintroducing explicit time dependence in Sec. V. Figure 1 shows the resulting, transverse, electric and magnetic fields. These are radially and azimuthally polarized respectively (as desired), orthogonal and of equal magnitude at every point. The transverse field profile as a function of b_l is also shown.

A further motivation for this choice of background is that its symmetries match well with those of the BLFQ basis; the potential (15) is proportional to the lowest order 2D-HO eigenstate Φ_{00} in coordinate space (recall $f \equiv 1$ for now and see Appendix C for details on the basis),

$$e\mathcal{A}^\mu(x) = \delta_-^\mu \frac{m_e a_0}{b_l} \Phi_{00}^{b_l}(x^\perp). \quad (16)$$

The classical physics of a particle in the model axicon field above is straightforward. Consider a particle near the symmetry axis. Depending on its charge, or equivalently the sign of a_0 , the particle is either repelled from or attracted to the axis. In the former case, see the right-hand panel of Fig. 1, the particle will be accelerated into the weak-field region and then drift outward at some acquired velocity. Particles of opposite charge will be attracted to the symmetry axis, which is also a weak field region, but will overshoot, and may then be attracted back again, oscillating around the axis. This intuition will help us analyze the results of the tBLFQ calculation, below.

A. Time-evolution in tBLFQ

We now give the first tBLFQ calculation. We begin with a single-electron state and consider its evolution in the

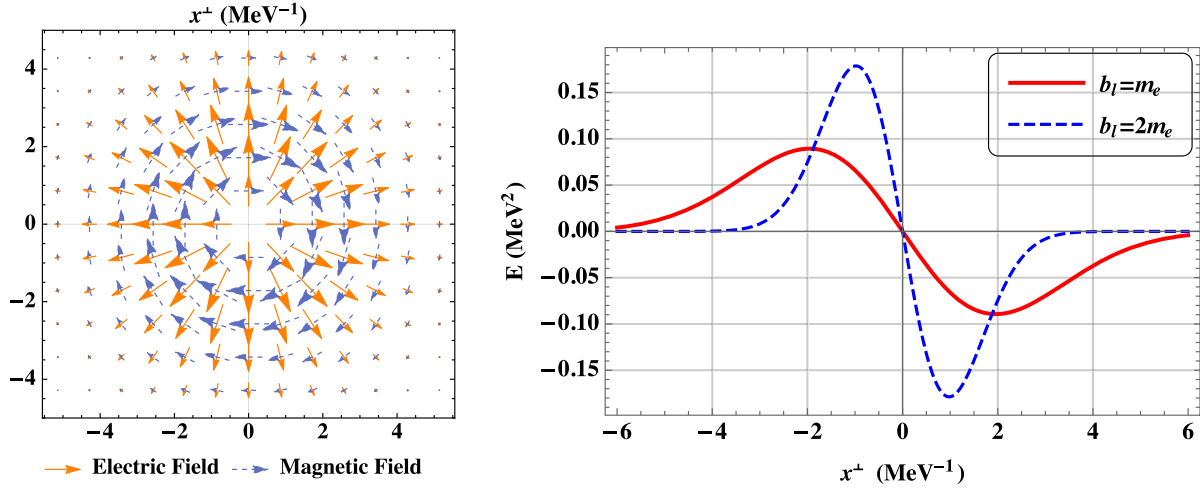


FIG. 1. Left: transverse electric and magnetic fields of the background, shown as solid and dashed lines respectively, as a function of the transverse coordinates. Parameters: $a_0 = 1$, $b_l = m_e$. The two fields are of the same magnitude and are perpendicular at every point. Right: transverse electric field for different widths b_l . As b_l increases the field becomes narrower in position space, but reaches higher peak values.

axicon background above. We take the longitudinal dependence of the initial state to be a plane wave of momentum p^+ . As the background has no longitudinal position dependence, longitudinal momentum p^+ is conserved (both classically and quantum mechanically). Thus we need only consider the physics of the transverse directions. For the transverse degrees of freedom, we take the initial wave function to be a normalized Gaussian wave packet, also proportional to the HO eigenstate $\Phi_{00}^b(x^\perp)$. We will take the initial state width b_i to be equal to electron mass m_e because it is the only mass scale in QED. Hence the electron is localized to within $\sim m_e$ in transverse momentum space and $\sim 1/m_e$ in transverse position space.

We evolve the wave function in time numerically by solving the Schrödinger equation (4) using the MSD2 scheme described above. This gives the wave function at each time step. In our calculation, the width of the 2D-HO basis in the transverse plane b is chosen to coincide with the width of the initial state $b_i = m_e$, in order to improve convergence. The mod-square of the wave function then gives the particle probability distribution in either coordinate or momentum space, as desired. (See [62] for a comprehensive discussion of the position operator in light-front quantization, and [63] for an application to radiation reaction in the front form.)

Figure 2 shows the time-evolution of the wave packet for $a_0 > 0$. (The background field Hamiltonian conserves angular momentum L_z , and our initial state is rotationally symmetric; hence it is enough to pick a slice through the transverse plane.) As time passes, the wave packet is dispersed and repelled from the origin, with the initial peak of high probability density being pushed to large transverse positions. This is the analog of the classical repulsion described above. We can also make a more

quantitative check against the classical theory, as follows. The classical equations of motion reduce, in the transverse directions, to the coupled ordinary differential equations

$$\ddot{x}^\perp = \frac{m_e a_0 b_l^2}{2p^+ \sqrt{\pi}} x^\perp \exp\left(-\frac{1}{2} b_l^2 x^\perp x^\perp\right), \quad (17)$$

where a dot indicates an x^+ derivative. We numerically integrate this equation for a large number of initial conditions sampled from the same Gaussian position and momentum space distributions as define the wave packet in the quantum calculation. The idea is then to track the position of the peak of the time-evolved distribution, expecting that this would be reproduced to some

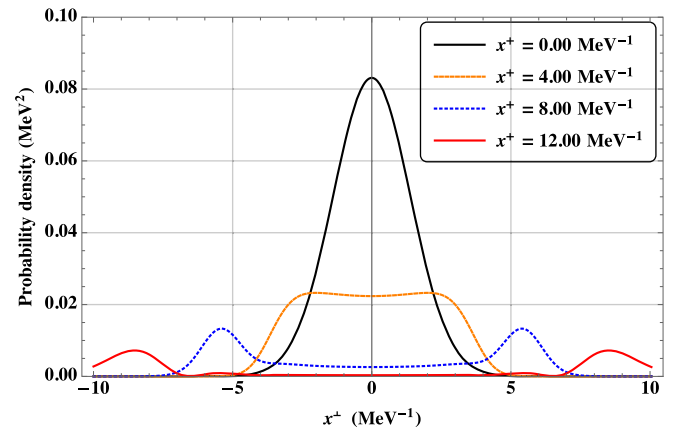


FIG. 2. A single electron initially in a HO basis state is evolved through time in the axicon background with $a_0 > 0$. The wave packet is dispersed from the origin, effectively feeling a repulsive electromagnetic field. Parameters: $N_{\max} = 200$, $K_{\text{total}} = 10.5$, $b = m_e$, $b_l = m_e$, $a_0 = 10$, $b_i = m_e$, $p^+ = 10.5$ MeV.

approximation by the peak of the quantum wave packet, where the probability density is highest. We find that the peak of the distribution shifts, in position space, out to $|x^\perp| \approx 5.5 \text{ MeV}^{-1}$ at $x^+ = 8 \text{ MeV}^{-1}$ and to $|x^\perp| \approx 8 \text{ MeV}^{-1}$ at $x^+ = 12 \text{ MeV}^{-1}$, in good agreement with the peak positions in the tBLFQ calculation; see Fig. 2. This is a very simple check, but it gives us confidence that we are capturing the correct physics both qualitatively and quantitatively.

In the quantum theory there is of course wave packet spreading even without an external field. We therefore continue to the case of $a_0 < 0$, all other parameters constant. This amounts to changing the sign of the electromagnetic field or, in this system, exchanging particles for antiparticles. The results are shown as a series of snapshots in Fig. 3. In this case the initial wave packet is first compressed to the origin, with the probability density increasing there, before expanding back out, and being recompressed toward the origin. This competition between wave packet focusing and spreading [61] naturally reflects the expected classical dynamics. The behavior initially appears reminiscent of that of a harmonic oscillator in a potential trap. However, the Gaussian falloff of our background means that there is no true trapping, and for large times the probability density around the origin falls to zero.

IV. PHOTON EMISSION

In the section above we discussed the time-evolution of a single electron in a background field, including only the $|e\rangle$ Fock state in the tBLFQ calculations. This represents the center-of-mass (CM) motion (in that it resembles the CM motion in nonrelativistic mechanics, as will be explained below). In this section, we add the second simplest Fock sector $|e\gamma\rangle$, with which we can study the photon emission and absorption. We are mainly interested in the production (stimulated by the background) of a physical photon from a physical electron, and the relative motion between them. However, the CM spectrum is continuous and typically mixed with the relative motion, making the full energy spectrum difficult to handle, even in a discrete basis. Fortunately, in BLFQ, it is possible to factorize the CM and the relative motion. By adopting the ‘‘Lagrange multiplier’’ method [64,65], we can isolate and study only the relative excitation.

As discussed in Sec. II, time-evolution is naturally studied in the tBLFQ basis $|\beta\rangle$ of QED eigenstates defined in Eq. (9). Calculations would be made even simpler if this basis had the property of factorization, such that the CM and relative motion could be separated, for then it would be possible to perform the calculation in the space of ‘‘relative’’ variables.

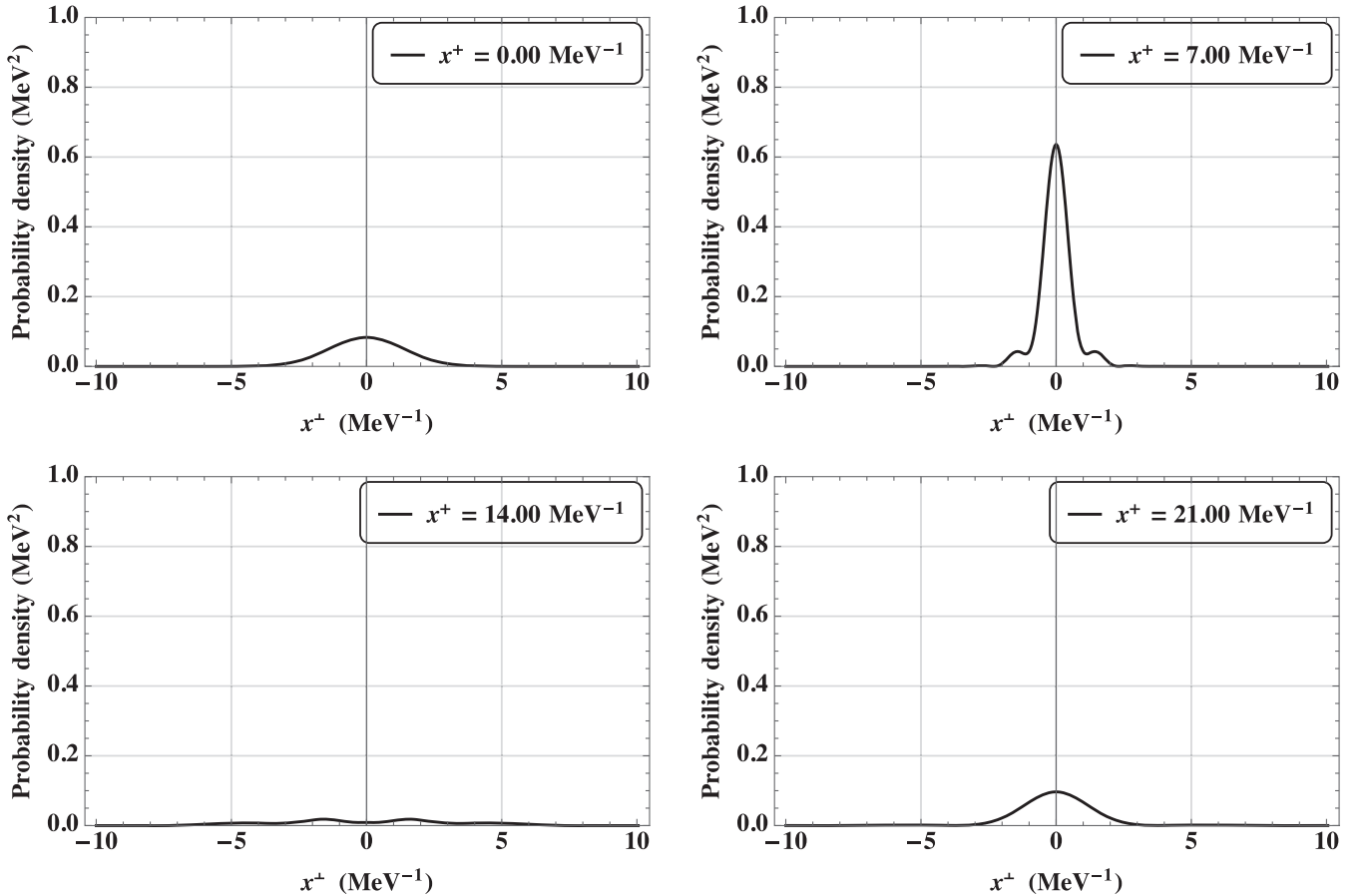


FIG. 3. For $a_0 < 0$, the same initial state as in Fig. 2 is first attracted to, and then shows oscillations around, the origin. Parameters: $N_{\text{max}} = 200$, $K_{\text{total}} = 10.5$, $b = m_e$, $b_l = m_e$, $a_0 = -10$, $b_i = m_e$, $p^+ = 10.5 \text{ MeV}$.

In the absence of a background (as in BLFQ), the wave function can indeed be factorized—the problem we must confront is that this may not be true in a background field (as in tBLFQ). Indeed the CM motion is mixed with the relative motion, in which we are mainly interested.

In what follows, we will in Secs. IV A and IV B describe the factorization in BLFQ, and then argue that we can *approximately* evolve (using tBLFQ) the system inside its CM frame, for simplicity. In Sec. IV C we define some useful observables which encode the physics of interest before going on to study their time dependence in Sec. IV D.

A. Factorization in the tBLFQ basis

In the light-front framework, the transverse boosts have the same algebra as the Galilean boosts, which implies that

$$\begin{aligned} \text{in the } |e\rangle \text{ sector;} & \quad \Psi(p_1^\perp, k_1, s_1) = \langle p_1^\perp, k_1, s_1 | \beta \rangle = \phi_{\text{CM}}(p_1^\perp) \psi_{\text{rel}}(k_1, s_1), \\ \text{in the } |e\gamma\rangle \text{ sector;} & \quad \Psi(p_1^\perp, p_2^\perp, k_1, k_2, s_1, s_2) = \langle p_1^\perp, p_2^\perp, k_1, k_2, s_1, s_2 | \beta \rangle = \phi_{\text{CM}}(P^\perp) \psi_{\text{rel}}(q^\perp, k_1, k_2, s_1, s_2), \end{aligned} \quad (18)$$

where 1 (2) represents the electron (photon); $P^\perp = p_1^\perp + p_2^\perp$ is the CM momentum and $q^\perp = x_2 p_1^\perp - x_1 p_2^\perp$ is the relative momentum in the transverse directions.

In general, this factorization will not be preserved when a finite (truncated) basis is used. There is however a method by which to retain exact factorization in finite truncation [57], as follows. We introduce a “modified CM Hamiltonian” H_{CM} defined by [57,66]

$$H_{\text{CM}} = \left(\sum_i p_i^\perp \right)^2 + b^4 \left(\sum_i x_i x_i^\perp \right)^2 = (P^\perp)^2 + b^4 (R^\perp)^2. \quad (19)$$

We then subtract from the QED Hamiltonian P_{QED}^- the real CM motion $(P^\perp)^2/P^+$, and replace it with H_{CM} , as so:

$$P_{\text{QED}}^- \rightarrow P_{\text{QED}}^- - \frac{(P^\perp)^2}{P^+} + \frac{\lambda}{P^+} (H_{\text{CM}} - 2b^2), \quad (20)$$

in which the term $-2b^2$ subtracts the zero-point energy, and $\lambda > 0$ is a Lagrange multiplier [64,65]. Unlike the original CM Hamiltonian, the modified H_{CM} commutes with P_{QED}^- in finite *truncated* BLFQ bases. As a result, exact factorization of the wave function is achieved, despite the finite truncation [57]. Furthermore, by modifying the CM Hamiltonian as above, we lift the *nonzero* CM excitations by $2\lambda(2N + |M|)b^2/P^+$ *without* altering the part of the spectrum with $N = M = 0$, i.e. the part without CM excitation. By choosing a sufficiently large λ , we lift all the CM excitations beyond the highest relative excitation of interest and thus we obtain the desired excitation spectrum of the relative motion.

Now, the modified CM Hamiltonian commutes with the full Hamiltonian; hence their common eigenstates and the

the CM motion can be decoupled from the relative motion as in nonrelativistic many-body systems [27], with the longitudinal momentum fraction $x_i = p_i^+/P^+$ playing the role of the nonrelativistic mass. Thus the CM position is defined as $R^\perp = \sum_i x_i x_i^\perp$, where x_i^\perp is the single-particle coordinate. These features allow us to factorize the CM motion and the relative motion as in nonrelativistic quantum mechanics.

The QED Hamiltonian P_{QED}^- (not including terms related to the background field) commutes with the CM (total) momentum P^\perp as well as the CM Hamiltonian $(P^\perp)^2/P^+$. They therefore have simultaneous eigenstates, and the corresponding wave function can be factorized:

CM motion have definite N and M . The two-particle wave function can thus be factorized as

$$\Psi(p_1^\perp, p_2^\perp, k_1, k_2, s_1, s_2) = \tilde{\Phi}_{NM}^b(P^\perp) \psi_{\text{rel}}(q^\perp, k_1, k_2, s_1, s_2), \quad (21)$$

where $\tilde{\Phi}_{NM}^b$ is the CM motion wave function and a HO eigenstate (see Appendix C) with $b = m_e = 0.511$ MeV; ψ_{rel} is the relative wave function, which we will ultimately use to construct observables. This relative wave function can be expanded in the HO basis as

$$\psi_{\text{rel}}(q^\perp, k_1, k_2, s_1, s_2) = \sum_{nm} f_{nmk_1k_2s_1s_2} \tilde{\Phi}_{nm}^{b_b}(q^\perp), \quad (22)$$

where the width $b_b := \sqrt{x_1 x_2} b$, and an x dependence is introduced for the exact factorization in finite truncated bases [57]. The x_j are the longitudinal momentum fractions of the electron, $x_1 = k_1/(k_1 + k_2)$, and the photon, $x_2 = 1 - x_1$. The coefficients f can be obtained by the Talmi-Moshinsky transformation [57,67,68], which is typically used to separate the CM motion and the relative motion in the HO basis.

B. Relative motion in a background field

In the presence of a background field, the full Hamiltonian $P^-(x^+) = P_{\text{QED}}^- + V(x^+)$ may not commute with the CM Hamiltonian $(P^\perp)^2/P^+$. In this case there is no longer an exact factorization of the wave function as in Eq. (18). This means that the relative motion will be affected by the CM motion and thus cannot be studied separately.

As presented in Fig. 4, however, the CM motion evolves much faster than the relative motion (in that the momentum

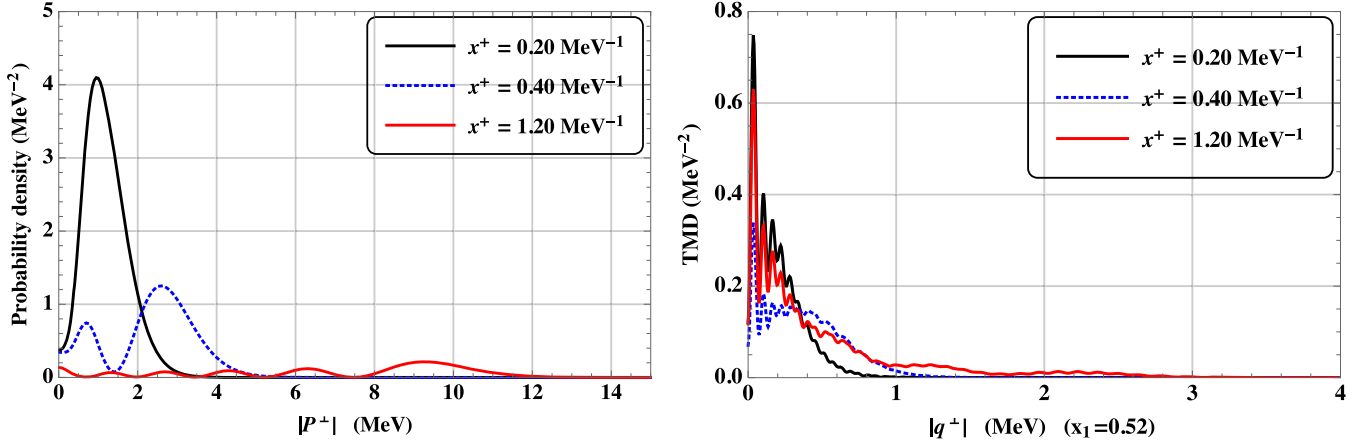


FIG. 4. Time-evolutions of the CM momentum probability density and the (relative) transverse momentum distribution (TMD), presented in the left and right panels respectively. The CM momentum increases at a much faster rate than the relative momentum as time passes. More details of TMDs will be discussed in the text. In the left panel $N_{\max} = 400$ and the initial state is a normalized Gaussian wave packet with a width $b = 0.511$ MeV; in the right panel $N_{\max} = 100$ and the initial state is the physical electron state. Other parameters are the same: $K_{\text{total}} = 10.5$, $b_l = m_e$, $a_0 = 200$, $P^+ = 10.5$ MeV.

distribution of the CM motion expands to places with higher momentum more quickly) in a field with $a_0 = 200$. We argue that their time-evolutions occur at different timescales and thus make an approximation that the wave function of the system can be factorized into a CM part and a relative part even in this background field and study them separately. We make a further approximation by showing as follows that we can neglect the change of the background field the relative motion “feels” induced by the CM motion. At every instant, we can always expand the factorized wave function of the system in terms of components with definite CM momenta. For every component with a definite CM momentum, its relative motion is effectively evolved in a CM frame in a background field that would be obtained by the boost, which takes the CM momentum to zero, of the background field (15). Now, because the background (15) has only a longitudinal component, it is left invariant under a boost in the transverse directions,

$$\begin{aligned} \mathcal{A}^+ &\rightarrow \mathcal{A}^+, & \mathcal{A}^\perp &\rightarrow \mathcal{A}^\perp + \mathcal{A}^+ C^\perp, \\ \mathcal{A}^- &\rightarrow \mathcal{A}^- + \mathcal{A}^\perp \cdot C^\perp + \mathcal{A}^+ C^{\perp 2}, \end{aligned} \quad (23)$$

where the C^\perp are dimensionless numbers that depend on the boost. What we should consider is the change of field caused by the displacement of the field from the origin in the CM frame. For now, we neglect this displacement of the background field. This approximation is valid at the beginning of the time-evolution during which the displacement is small. We can then simply evolve the system inside the subspace with no CM motion, that is the subspace $N = M = 0$, with the original background field in Eq. (16). This is equivalent to saying that we only evolve the relative wave function ψ_{rel} in Eq. (21).

In the future work, we plan to expand our basis by including nonzero N and M , i.e. nonzero CM excitations. Thereby we can have an accurate description of the CM motion as well as the relative motion coherently and simultaneously.

C. Observables

We retain, from the whole Fock space, only the $|e\rangle$ and $|e\gamma\rangle$ Fock sectors. In the latter, we can study photon emission and the relative motion between the electron and the emitted photon. However, the $|e\gamma\rangle$ sector also contains nonphysical photons, namely those “bare” photons which contribute to the physical electron. These are not real photons that are emitted and can be measured by a detector. However, we find in our calculations that the excited states of P_{QED}^- are dominated by the $|e\gamma\rangle$ sector; the overlap of excited states with the $|e\rangle$ sector, mod-squared, being less than 10^{-3} at $N_{\max} = 100$, $K_{\text{total}} = 10$. This implies that the electron and the photon are weakly coupled, and therefore we approximate the excited states (of P_{QED}^-) as scattering states of a physical electron and a physical photon. In this paper, we focus on the physical part of the wave function $\bar{\Psi}$ which is obtained by projecting out the physical electron state

$$\begin{aligned} \bar{\Psi}(p_1^\perp, p_2^\perp, k_1, k_2, s_1, s_2; x^+) \\ := \langle p_1^\perp, p_2^\perp, k_1, k_2, s_1, s_2 | (\mathbf{1} - |e\rangle_{\text{phys}}\langle e|_{\text{phys}}) | \Psi; x^+ \rangle_S, \end{aligned} \quad (24)$$

where $|e\rangle_{\text{phys}}$ is the physical electron state; $|\Psi; x^+\rangle_S$ is the Schrödinger picture state vector obtained by using Eqs. (4) and (5). Physically, this eliminates the nonphysical bare photon contribution. It is easiest to achieve the projection in

the $|\beta\rangle$ basis, where it corresponds to simply dropping the lowest component of the wave function.

It would be desirable to obtain the scattering matrix in tBLFQ; however, the systematic construction of “out states” [16] is a topic for future work. In this paper we instead concentrate on two observables at finite time. They are the longitudinal momentum distribution (LMD), which describes the longitudinal motion, and the transverse momentum distribution (TMD), which describes the transverse motion.¹ These are defined in terms of the *physical* part of the relative wave function as

$$\text{LMD}(x_1; x^+) := \sum_{s_1, s_2} \int \frac{d^2 q^\perp}{(2\pi)^2} |\tilde{\psi}_{\text{rel}}(q^\perp, k_1, k_2, s_1, s_2; x^+)|^2, \quad (25)$$

$$x_1 = \frac{k_1}{k_1 + k_2},$$

$$\text{TMD}(q^\perp, x_1; x^+) := \sum_{s_1, s_2} |\tilde{\psi}_{\text{rel}}(q^\perp, k_1, k_2, s_1, s_2; x^+)|^2, \quad (26)$$

in which the relative wave function $\tilde{\psi}_{\text{rel}}$ is obtained by a TM transformation from the two particle wave function (24), as in Eqs. (21) and (22). The meaning of the LMD is that it gives the probability of finding a scattered electron-photon state with an electron longitudinal momentum fraction $x_1 = k_1/(k_1 + k_2)$. The TMD tells us the probability of finding the scattered electron-photon states with relative *transverse* momentum q^\perp , and momentum fraction of the physical electron, x_1 . If, in Eq. (26), we drop the sum over s_1 and s_2 then we obtain a spin-dependent TMD. If the TMD (26) is integrated over q^\perp then we recover the LMD (25).

D. Time-evolution in tBLFQ

In this section we discuss the time-evolutions of the LMD and the TMD, as defined above. We start with the physical electron state, which is the lowest eigenstate of P_{QED}^- *without* the background field. Evolving this state in time, with the background (16) present, we obtain the wave function at each time step. (This is performed numerically using the MSD2 scheme described in Sec. II A.) Equations (25) and (26) then tell us how to extract the time-evolutions of the LMD and the TMD, respectively.

In order not to overcomplicate the presentation and obtain a broad overview of the physics involved, we start by considering a simple quantity, namely the probability to find a physical photon of *any* (longitudinal or transverse) momentum; this is obtained by integrating the LMD over all longitudinal momentum fractions x_1 . Figure 5 shows this probability as a function of time in backgrounds with

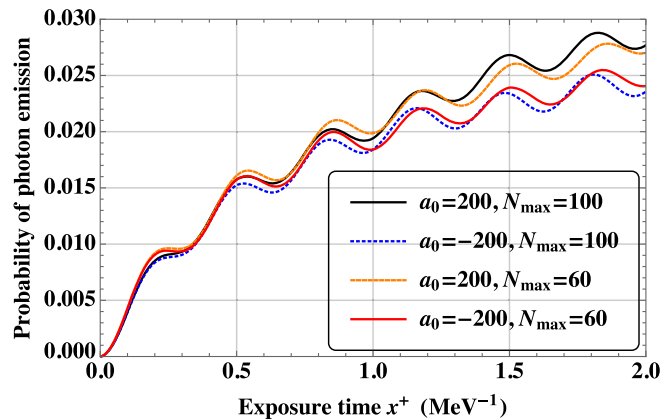


FIG. 5. Probability to find electron-photon states, or rather the probability of photon emission. This probability initially increases with time. The later oscillations are discussed in the text. Parameters: $N_{\text{max}} = 100$ and 60 , $K_{\text{total}} = 10.5$, $b_l = m_e$, $a_0 = \pm 200$, $p^+ = 10.5$ MeV.

$a_0 = \pm 200$, i.e. fields which give, broadly speaking, repulsive and attractive forces to the electron, respectively. Initially, at light-front time $x^+ = 0$, there is only a physical electron, and no physical photons; hence the probability starts from zero. As time passes the probability increases as electrons are raised to excited states by the background. Initially, the probability of photon emission is independent of the sign of a_0 , but at later times the sign of the field begins to have an effect. As can be seen in the figure, all the curves show a similar trend—the probability of photon emission increases, but with small oscillations. This reflects the exchange of energy/momentum back and forth between the electrons and produced photons as time passes. We observe that, numerically, these oscillations come from the matrix element $\langle \text{lowest state} | V | \text{lowest state} \rangle$, which is significantly bigger than the other matrix elements.

We now turn to what are essentially the *differential* photon emission probabilities. We begin by studying, through the TMD, the time-evolution of only the *transverse* motion of the physical electron, fixing $k_1 = 5.5$, i.e. the *longitudinal* momentum fraction $x_1 = 5.5/10.5 \approx 0.52$. We will study how this depends on N_{max} and a_0 . Finally, we will reinstate the longitudinal dependence, allowing us to study the time-evolutions of TMDs with different longitudinal momentum fractions x_1 and the LMD. A comparison of the time-evolution of TMDs obtained in tBLFQ and perturbation theory, in the limit of weak background fields, is discussed in Appendix E.

1. N_{max} dependence

We should of course address how our results depend on the various parameters defining our basis. Here we focus on N_{max} , which is a truncation parameter for transverse degrees of freedom. It limits the energy (12) of the 2D-HO basis. In a basis with a larger N_{max} , we have a higher

¹Since the momentum distribution is more commonly measured in experiments, we will focus on the distribution in momentum space in the remainder of the paper.

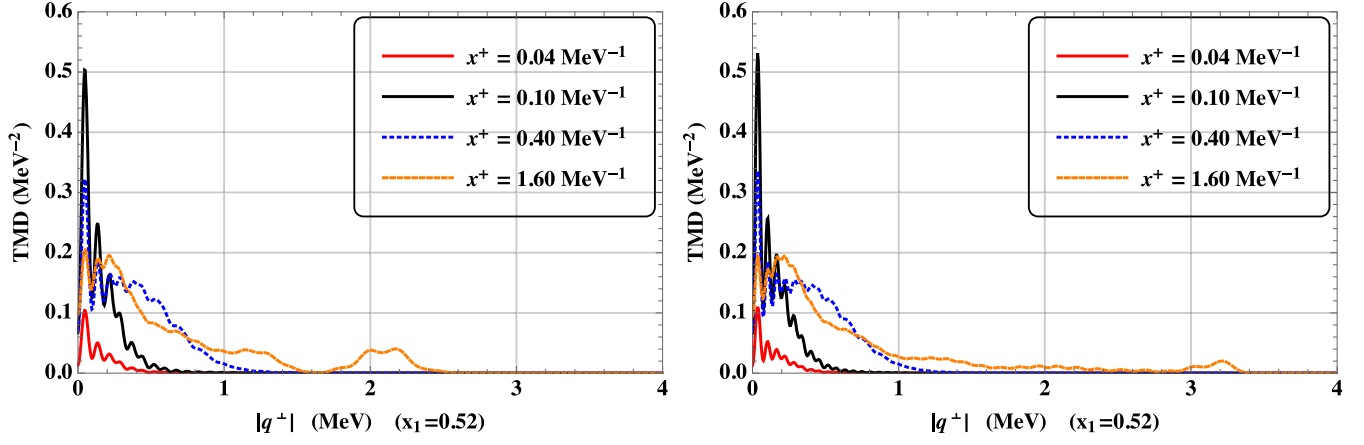


FIG. 6. Time-evolution of the TMD in two bases with (left) $N_{\max} = 60$ and (right) $N_{\max} = 100$. All other parameters are the same in both calculations: $K_{\text{total}} = 10.5$, $b_b = 0.50m_e$, $b_l = m_e$, $a_0 = 200$, $x_1 = 0.52$, $p^+ = 10.5$ MeV. Initially there is only a physical electron. As time passes, the transverse momenta are excited to higher values. For the smaller basis (left-hand panel), the excitations hit the UV cutoff (at around 2.6 MeV) at a time $x^+ \approx 1.2$ MeV $^{-1}$, whereas for the larger basis (right-hand panel), the UV cutoff (now at 3.4 MeV due to the larger N_{\max}) at a later time of $x^+ \approx 1.6$ MeV $^{-1}$. Overall, there is little difference between the curves until they hit the UV cutoff. Beyond the corresponding “cutoff” times, the curves are no longer reliable.

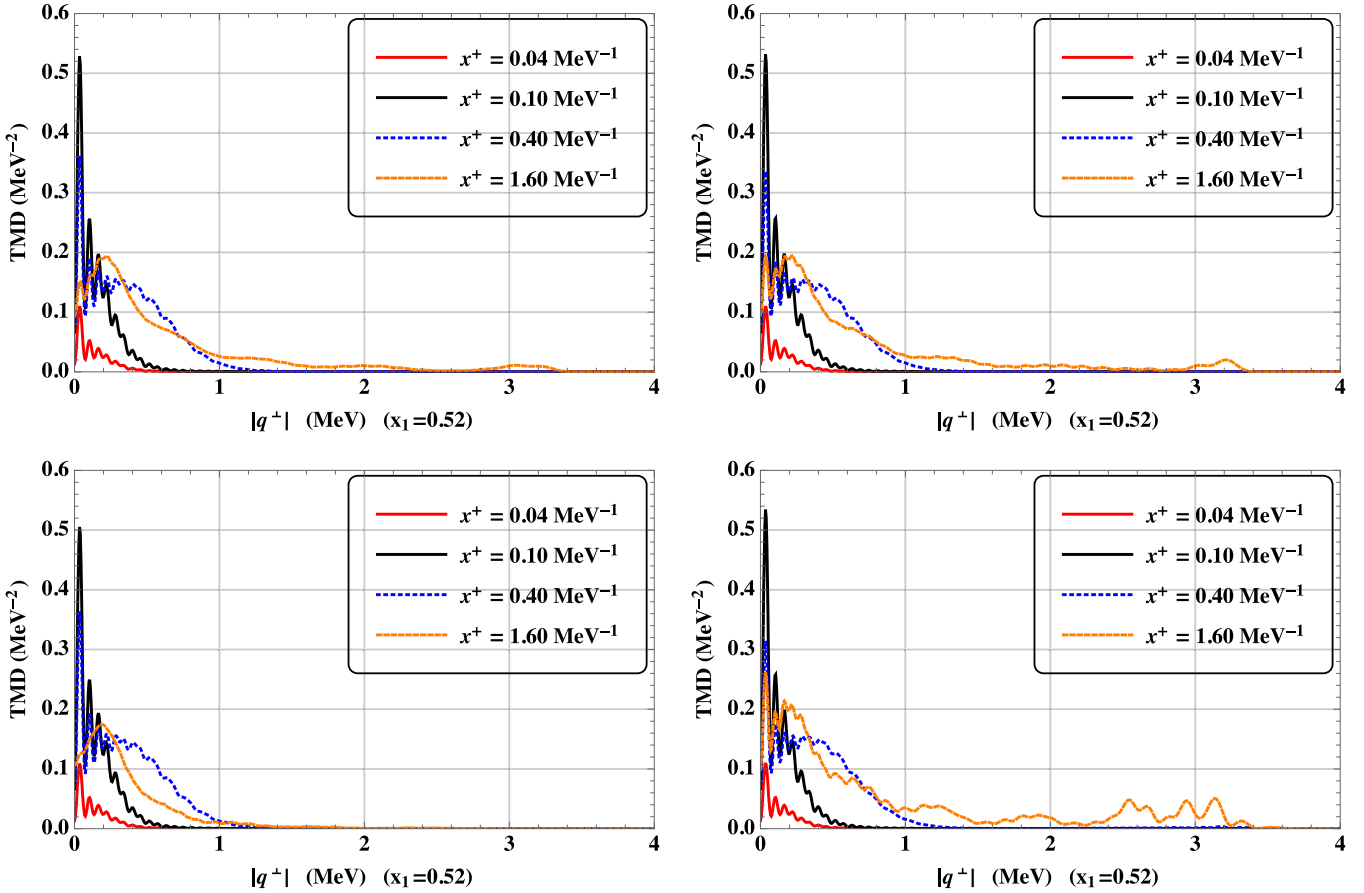


FIG. 7. Time-evolution of TMDs. In all plots, $N_{\max} = 100$, $K_{\text{total}} = 10.5$, $b_b = 0.50m_e$, $b_l = m_e$, $x_1 = 0.52$. In the left-hand column, $a_0 = -200$; in the right-hand column $a_0 = 200$. Longitudinal momentum P^+ decreases from $P^+ = 10.5$ MeV to $P^+ = 1.05$ MeV as we move from the upper to the lower row. In the left column (attractive field), as P^+ decreases, the higher momentum excitations decrease. Contrastingly, in the right column (repulsive field), the higher momentum excitations increase as P^+ decreases.

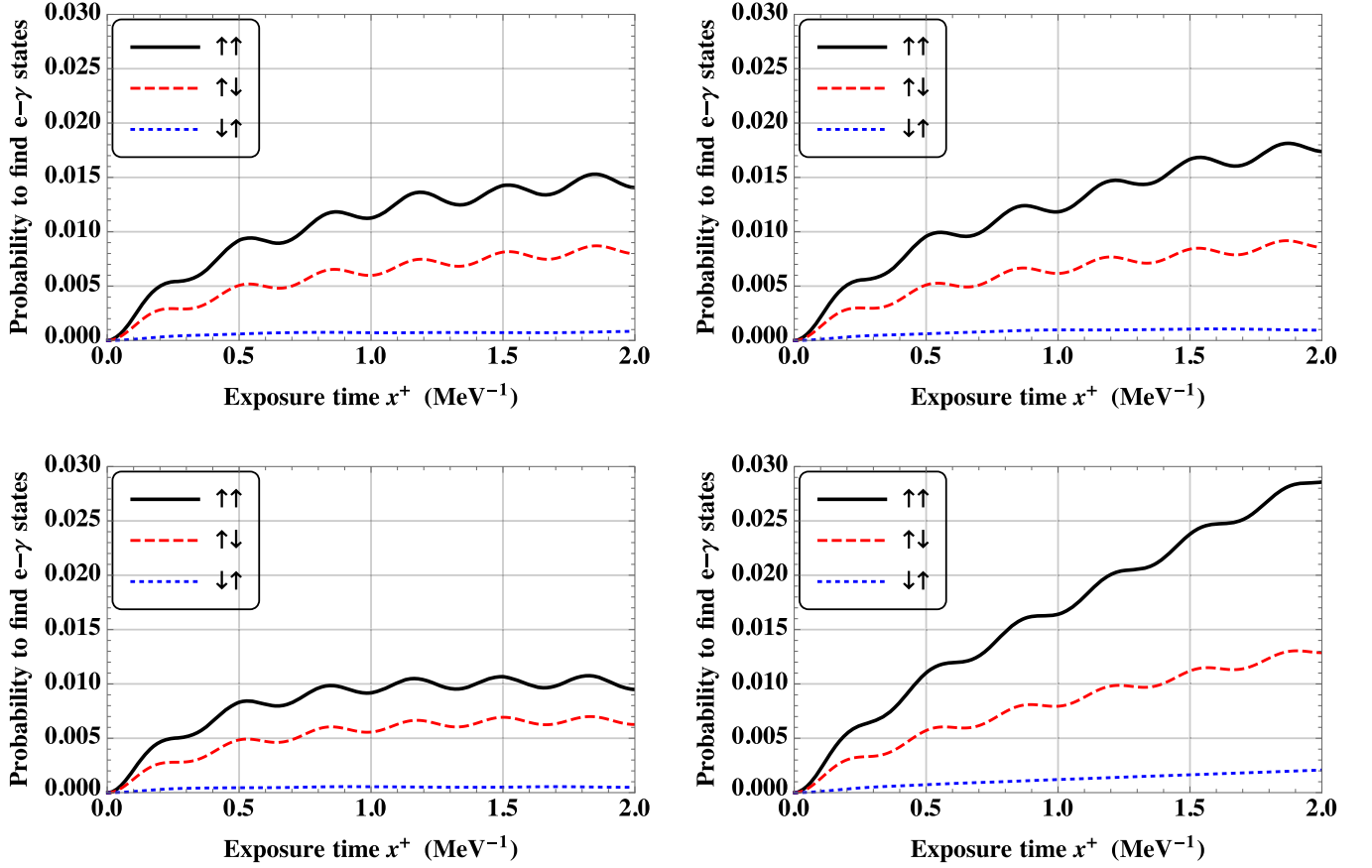


FIG. 8. Spin-resolved probability of finding electron-photon states. In all plots, $N_{\max} = 100$, $K_{\text{total}} = 10.5$, $b_l = m_e$. In the left-hand column, $a_0 = -200$; in the right-hand column $a_0 = 200$. Longitudinal momentum decreases from $P^+ = 10.5$ MeV to $P^+ = 1.05$ MeV as we move from the upper to the lower panels. At higher momentum, the probability of finding any particular spin configuration increases faster for $a_0 > 0$ than for $a_0 < 0$. At lower momentum, the curves show more differences between $a_0 > 0$ and $a_0 < 0$.

UV cutoff, proportional to $\sqrt{N_{\max}}$, and a lower IR cutoff, proportional to $1/\sqrt{N_{\max}}$ in momentum space [20]. Thus, larger N_{\max} offers a more complete description of the system. However, Fig. 5 shows that the time-evolution of the probability of photon emission only depends weakly on N_{\max} , at least initially.²

Figure 6 shows the time-evolution of a TMD in the background (16) with $a_0 = 200$. The two panels show the same calculation performed for two values of N_{\max} , namely 60 (left) and 100 (right). Now, while the background does not couple directly to the photon, it can excite the electron to higher excited states in which the relative motion between the electron and the photon is more significant; this is clear in the figure. As is shown by comparing the two panels, the overall shape of the TMD is not strongly dependent on N_{\max} , with the main difference being that for the larger N_{\max} the curve is smoother, since the resolution is better. This is,

²In our calculation, we retain only the lowest 1921 (1121) eigenstates of P_{QED}^- , which have no CM excitation, for $N_{\max} = 100$, $K_{\text{total}} = 10$ ($N_{\max} = 60$, $K_{\text{total}} = 10$); this means that we retain the states with invariant mass less than 6.67 MeV (5.02 MeV).

however, all *provided* that the momentum excitations do not reach the “boundary” implied by the UV cutoff of the basis, beyond which the results are subject to truncation artifacts. For the larger basis, the curves hit the (larger) cutoff at later

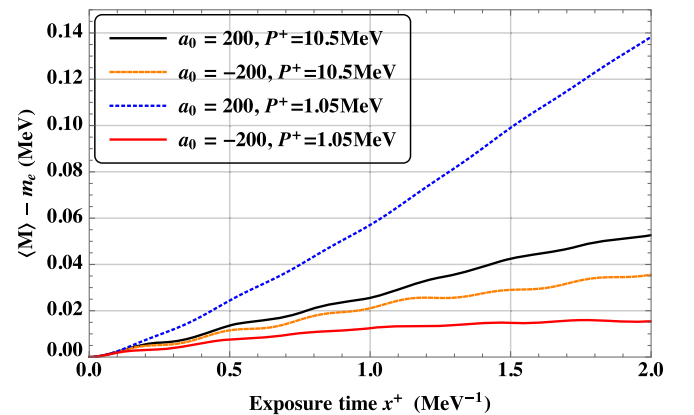


FIG. 9. Time-evolution of the invariant mass of the physical electron in background fields (the initial invariant mass $m_e = 0.511$ MeV is subtracted). As time passes, the invariant mass is excited to higher values in all cases. Other parameters: $N_{\max} = 100$, $K_{\text{total}} = 10.5$, $b_l = m_e$.

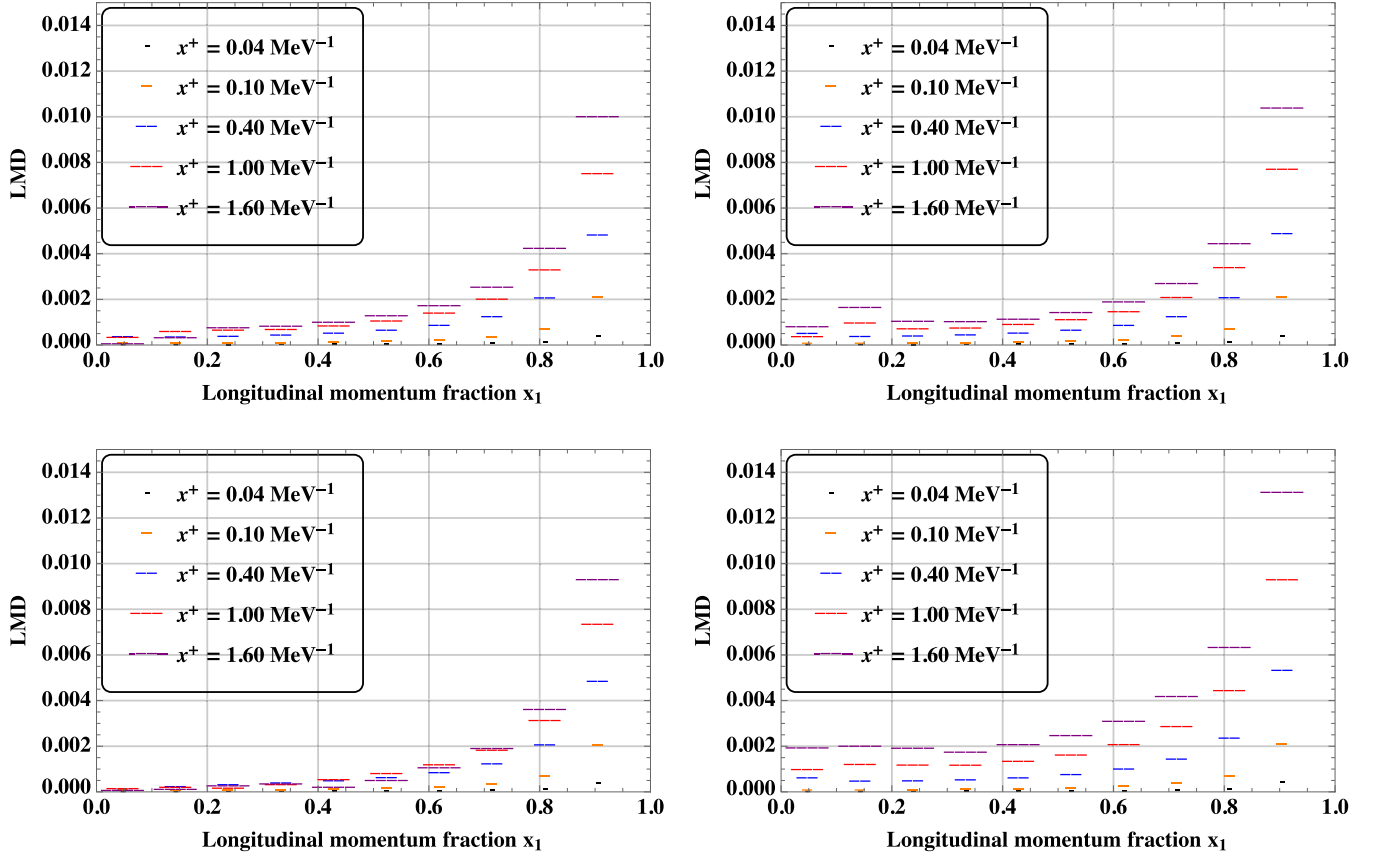


FIG. 10. Time-evolution of the LMD. In all plots, $N_{\max} = 100$, $K_{\text{total}} = 10.5$, $b_l = m_e$. In the left-hand column, $a_0 = -200$; in the right-hand column $a_0 = 200$. Longitudinal momentum P^+ decreases from $P^+ = 10.5$ MeV to $P^+ = 1.05$ MeV as we move from the upper to the lower row. In all the panels the distributions continue to increase with time at almost every electron longitudinal momentum fraction x_1 . In the left column (attractive field), as P^+ decreases, the smaller x_1 excitations decrease. Contrastingly, in the right column (repulsive field), the smaller x_1 excitations increase as P^+ decreases.

times. (We remark that while there exist “rescaling” methods [69] which improve the convergence of the wave function destroyed by the Fock-sector truncation, we need not use the method in this paper because of the good convergence of our results with N_{\max} .)

2. a_0 dependence and P^+ dependence

Following the above, we now fix $N_{\max} = 100$. In Fig. 7 we consider the dependence of the TMD on the sign of a_0 , i.e. on whether the background is attractive or repulsive, and on the total longitudinal momentum of the physical electron P^+ . We begin with $P^+ = 10.5$ MeV as shown in the upper two panels of the figure. In the left-hand panel, $a_0 = -200$ and the field is essentially attractive. In the right-hand panel, $a_0 = +200$ and the field is repulsive. In both cases, as time passes, TMDs become distributed across an increasingly broad range of momenta. TMDs are almost identical for short times, while for longer larger times there are, by time $x^+ = 1.6$ MeV $^{-1}$, slightly more higher momentum excitations for $a_0 > 0$. (In both cases the curves hit the boundary at about $x^+ = 1.60$ MeV $^{-1}$ and are then subject to artifacts.)

Figures 5 and 7, as well as Figs. 8 and 9 below, all show that the short-time evolution of the physical electron is not strongly dependent on the sign of the background, at least for $P^+ = 10.5$ MeV. We attribute this to the large light-front momentum chosen, implying high energy, for which accelerated charges emit in a narrow cone around their forward direction (instantaneous synchrotron radiation [70]); hence it makes little difference whether the background is attractive or repulsive.

In order to accentuate the differences due to the sign of a_0 , we therefore reduce the total longitudinal momentum P^+ by changing the length of the circle in Eq. (11) from³ $L = 2\pi$ MeV $^{-1}$ to $L = 20\pi$ MeV $^{-1}$. As a result, P^+ becomes equal to $0.1K_{\text{total}}$ in units of MeV. The resulting time-evolution of the electron with this smaller P^+ is shown in the lower panels of Fig. 7 (see also Figs. 8 and 9, below). In this figure $K_{\text{total}} = 10.5$ but the longitudinal momentum P^+ equals 10.5 or 1.05 MeV. As expected, there are larger

³The choice of $L = 2\pi$ MeV $^{-1}$ allows the longitudinal quantum number k to be interpreted as the longitudinal momentum in units of MeV.

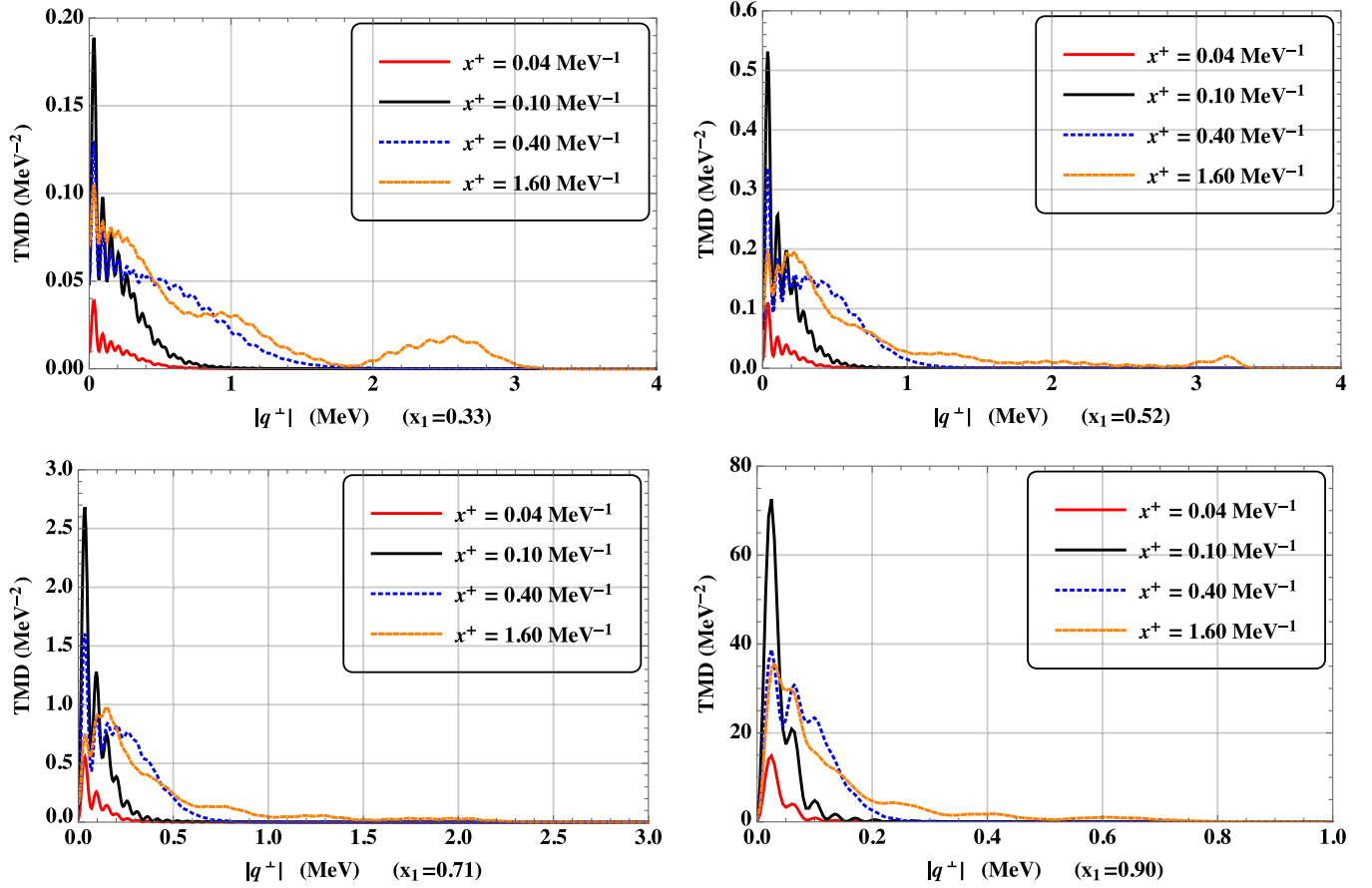


FIG. 11. Time-evolution of TMDs for longitudinal momentum fractions $x_1 = 0.33, 0.52, 0.71$, and 0.90 , as indicated in each plot; the corresponding widths of the HO basis are $b_b = 0.47m_e, 0.50m_e, 0.45m_e$, and $0.29m_e$. Other parameters: $N_{\max} = 100$, $K_{\text{total}} = 10.5$, $b_l = m_e$, $a_0 = 200$, $p^+ = 10.5$ MeV. Two obvious differences between the plots (noting the different scales on both the horizontal and vertical axes) are that the larger x_1 electrons have larger probabilities of being excited and that the distributions are excited to higher momentum more quickly for TMDs with smaller x_1 .

differences visible in the plots for opposite signs of a_0 when P^+ is smaller.

3. Time-evolution in different spin components

Figure 8 shows the spin-resolved probability of photon emission. We consider the electron/photon spin configurations $\uparrow\uparrow, \uparrow\downarrow, \downarrow\uparrow$. (The $\downarrow\downarrow$ contribution is zero and not included in the plots.) As seen in each of the four panels, the spin-flip component $\downarrow\uparrow$ is very small. The reason for this is that the background does not directly change the particle spin, and the spin-flip matrix elements from the background-free parts of the QED Hamiltonian are small [54]. Figure 8 also shows that there is a more significant dependence on the sign of the field for smaller P^+ than for larger, as was also observed above. Note that if we sum over all the spin configurations, we recover Fig. 5 as we should.

4. Invariant mass time-evolution

The invariant mass is defined as $M^2 = P^2 = (p_e + p_\gamma)^2$ and therefore its expectation value is

$$\langle M^2 \rangle = \sum_{\beta} |c_{\beta}(x^+)|^2 P_{\beta}^- P^+, \quad (27)$$

where $c_{\beta}(x^+) = \langle \beta | \Psi(x^+) \rangle$ is the wave function in the $|\beta\rangle$ basis. (Again, the transverse momentum is not included since the CM motion in the transverse directions has been eliminated.) Figure 9 shows the time-evolution of the expectation value of the invariant mass of the physical electron (the initial mass $m_e = 0.511$ MeV is subtracted) in the presence of background fields, with $a_0 = \pm 200$. Initially, there is only a physical electron, but as time increases a photon can be produced, which also contributes and increases the invariant mass. The increase in the invariant mass also means that energy is being injected into the system by the background.

While the invariant mass in Fig. 9 increases for both signs of a_0 , the increase is faster for $a_0 > 0$. This is because, recalling that the electrons are initially centered around the origin, the repulsive background drives electrons outward toward the strong parts of the field, while in an attractive background the electrons are kept near the origin, where the

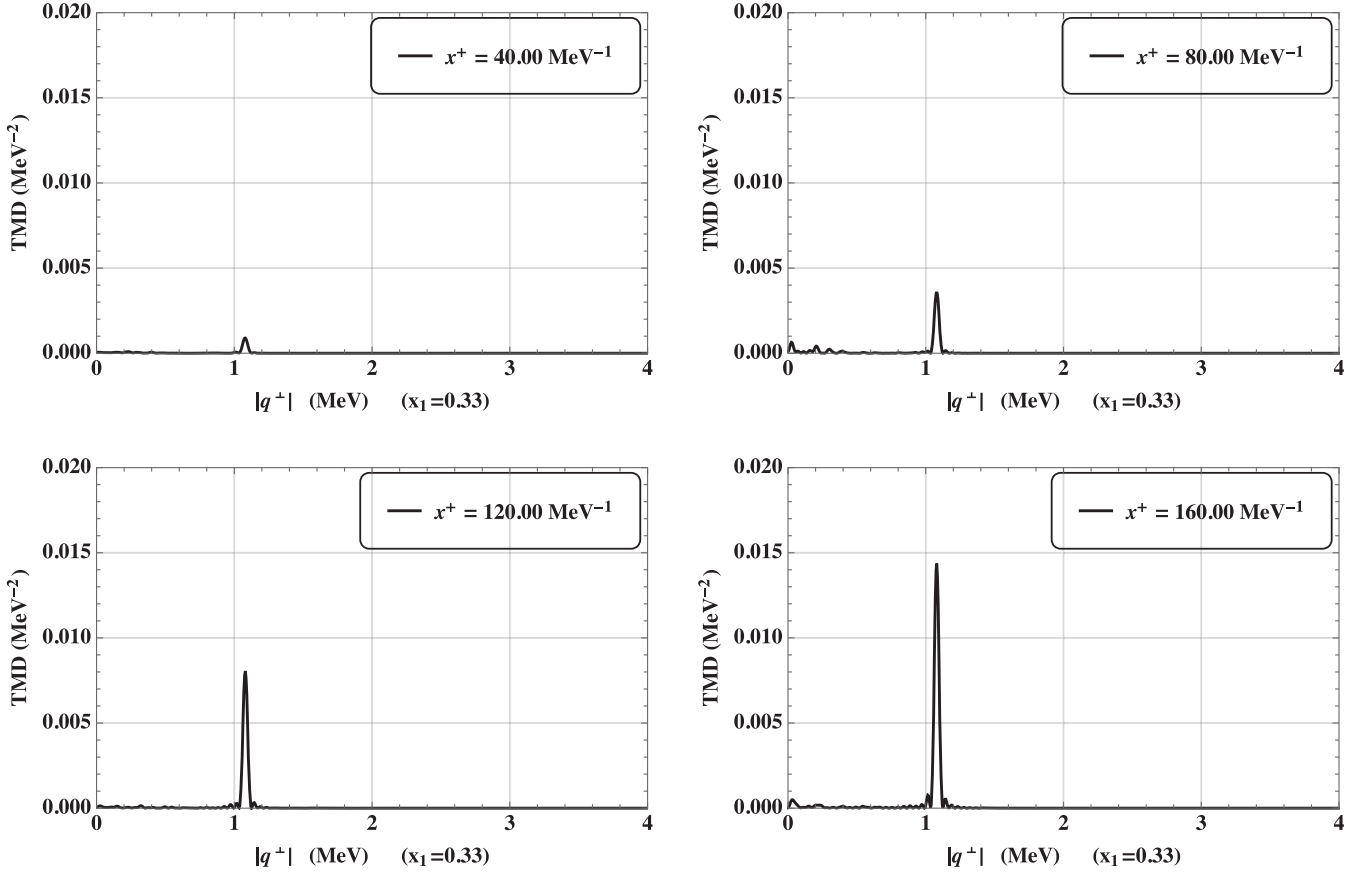


FIG. 12. Time-evolution of the TMD in a time-dependent background field ($a_0 = 1$) with a time dependence $f(x^+) = \sin(\omega x^+)$ in which $\omega = 1.92$ MeV. Four panels show TMDs in increasing light-front time x^+ . As time passes, only the excitation from the state with $\Delta P^- = 2\omega$ increases significantly. Other parameters: $N_{\max} = 160$, $K_{\text{total}} = 1.5$, $b_b = 0.47m_e$, $b_l = m_e$, $x_1 = 0.33$, $P^+ = 1.5$ MeV.

field is weak. Again, the differences between $a_0 > 0$ and $a_0 < 0$ are accentuated by decreasing P^+ .

5. Reinstating the longitudinal degrees of freedom

As discussed in Sec. III, there is no longitudinal coordinate dependence in our background field (16) and so the total longitudinal momentum of the system is conserved even in the presence of the background field. However, since the QED interaction couples the transverse and longitudinal degrees of freedom, the relative fraction of the total (conserved) longitudinal momentum each particle carries can change as time passes. In this section we study the longitudinal degrees of freedom.

In fact, we will begin with the longitudinal degrees of freedom *only*, by integrating out the transverse degrees of freedom and considering the longitudinal motion of the electron-photon state (LMD) and its dependence on the background field. Figure 10 shows the time-evolution of LMD in backgrounds with both $a_0 > 0$ and $a_0 < 0$ (left- and right-hand columns respectively). In the upper and lower rows we use total longitudinal momenta $P^+ = 10.5$ MeV and $P^+ = 1.05$ MeV respectively. The

probability at (almost) every x_1 increases with time in background fields with both $a_0 > 0$ and $a_0 < 0$. As shown in the left column ($a_0 = -200$), there are fewer small x_1 excitations as we reduce the total longitudinal momentum (from the upper to the lower panel). By contrast, in the right column ($a_0 = 200$), there are more low x_1 excitations as we reduce the total longitudinal momentum (from the upper to the lower). This is because the photons with higher transverse momentum (see Fig. 7) tend to have higher longitudinal momenta, which results in low electron momenta, i.e. small x_1 .

Next, we consider the time-evolution of TMDs in four different $k_1 = 3.5, 5.5, 7.5,$ and 9.5 , i.e. four different longitudinal momentum fractions of the electron, $x_1 \approx 0.33, 0.52, 0.71,$ and 0.90 . They are shown in Fig. 11. As time passes, basis states with higher momenta are populated. One significant difference between the four panels is that the larger x_1 electrons have larger probabilities of being excited (note the scale of the y-axes), meaning that the generated photons tend to have small longitudinal momenta. To explain this, consider the following approximation of the invariant mass $s_2 \equiv P^2$ of the outgoing electron-photon states:

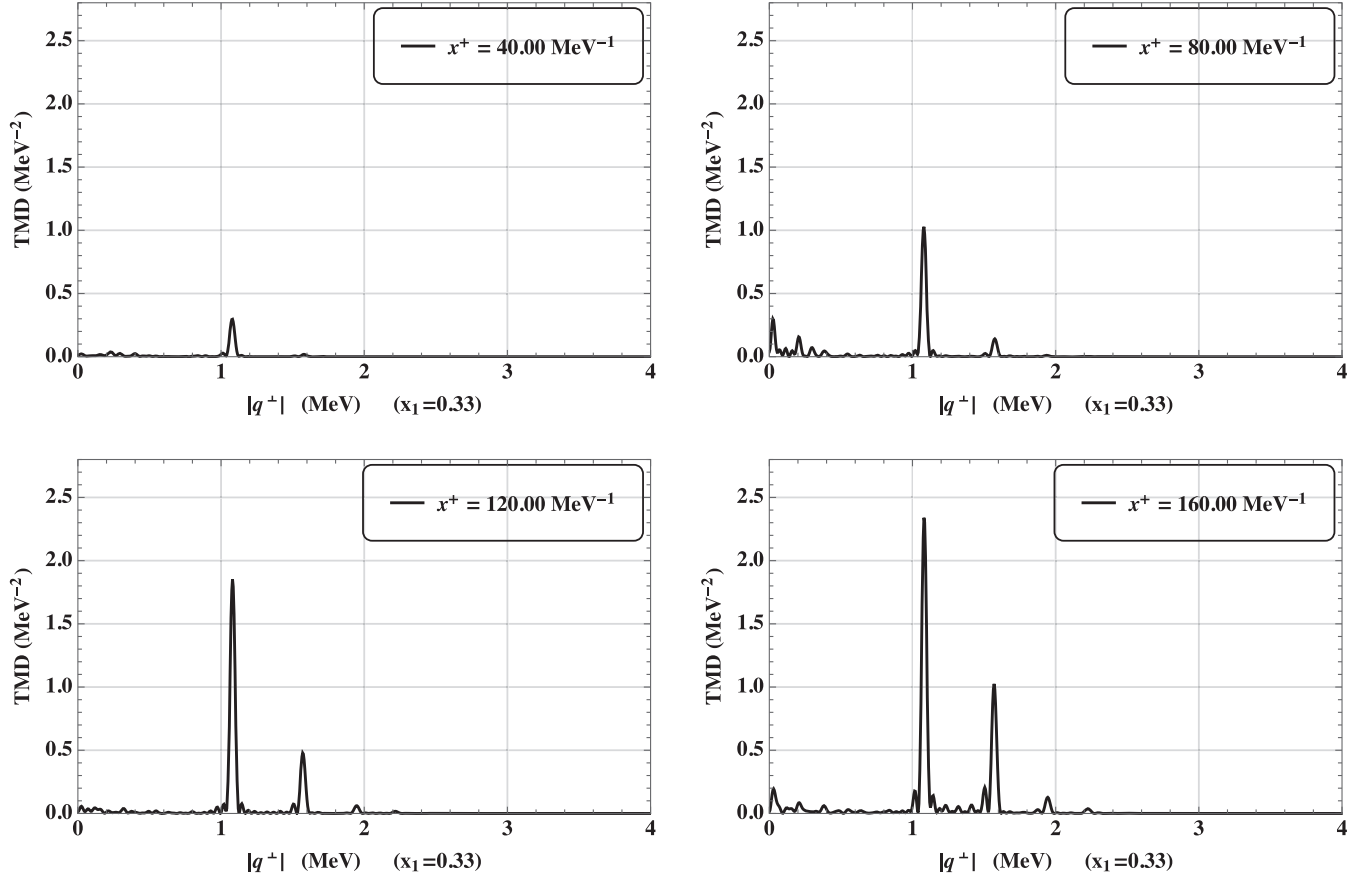


FIG. 13. Time-evolution of the TMD in a time-dependent background field ($a_0 = 20$) with a time dependence $f(x^+) = \sin(\omega x^+)$ in which $\omega = 1.92$ MeV. Four panels show TMDs in increasing light-front time x^+ . At initial stages there is only the contribution from state with $\Delta P^- = 2\omega$, as time passes, the higher excited states increase as shown in the down two panels. Other parameters: $N_{\max} = 160$, $K_{\text{total}} = 1.5$, $b_b = 0.47m_e$, $b_l = m_e$, $x_1 = 0.33$, $P^+ = 1.5$ MeV.

$$\begin{aligned}
 s_2 &= P^+ P^- - (P^\perp)^2 \approx P^+ \left[\frac{(p_1^\perp)^2 + m_e^2}{p_1^+} + \frac{(p_2^\perp)^2}{p_2^+} \right] \\
 &\quad - (p_1^\perp + p_2^\perp)^2 \\
 &= \frac{(p_1^\perp)^2 + m_e^2}{x_1} + \frac{(p_2^\perp)^2}{x_2} - (p_1^\perp + p_2^\perp)^2 \\
 &= \frac{(q^\perp)^2 + m_e^2}{x_1} + \frac{(q^\perp)^2}{x_2}. \tag{28}
 \end{aligned}$$

As Eq. (28) shows, because of the nonzero mass of the electron, small x_1 electrons carry more energy and therefore the corresponding photons, with $x_2 = 1 - x_1$, have a lower probability of being excited. Another main difference between the four panels in Fig. 11 is that the momenta increase faster (note the scale of the x-axes) in panels with smaller longitudinal momentum fractions. This is because, as in BLFQ and perturbation theory [26], the physical electron (our initial state) has broader TMDs in smaller x_1 ; in a Gaussian background field of width b_l , the width of TMDs does not change too much and therefore TMDs with smaller x_1 are still wider (move faster).

V. PHOTON EMISSION IN TIME-DEPENDENT BACKGROUND FIELDS

In previous sections the background field was independent of time; i.e. we took $f \equiv 1$ in (15). In such backgrounds, the probability that the physical electron emits a physical photon will naturally continue to rise as time evolves (as will the relative momentum between the electron and photon). In this section we add a time dependence to the background field by taking $f(x^+) = \sin(\omega x^+)$ in Eq. (15), where $\omega = 1.92$ MeV, chosen such that the energy difference between the lowest state and the 100th state (with energies labeled P_1^- and P_{100}^-) is $P_{100}^- - P_1^- = 2\omega$. We fix $K_{\text{total}} = 1.5$ and $N_{\max} = 160$, and thus the basis in the transverse plane is more complete⁴; as a result, the longitudinal momentum fraction can only take one value $x_1 = 0.5/1.5 \approx 0.33$. In this section, we fix $L = 2\pi$ MeV⁻¹ and therefore $P^+ = 1.5$ MeV.

⁴In this section we retain the lowest 318 states, which means that we retain the states with invariant mass less than 8.85 MeV.

We consider two cases, $a_0 = 1$ and $a_0 = 20$, corresponding to comparatively weak and strong backgrounds. Figure 12 shows the TMD for the case $a_0 = 1$. The TMD consists of, essentially, a single peak, the height of which increases as time passes. The peak is located at an energy difference, relative to the ground state, of $\Delta P^- = P_\beta^- - P_1^- = 2\omega$. (Recall that the excited states in $|\beta\rangle$ are scattering states, and therefore appear as peaks in the TMD.) This agrees well with Fermi's golden rule.

We now increase the background field amplitude to $a_0 = 20$, and Fig. 13 shows the resulting TMD. Initially, for small elapsed time, there is only a single peak, again at an energy corresponding to $\Delta P^- = 2\omega$, as for $a_0 = 1$. As time continues to pass, however, additional states are excited, and further peaks appear in the spectrum. These additional peaks are located at an energy difference of $\Delta P^- = 4\omega, 6\omega, 8\omega$, and so on. These can be understood as *higher harmonics* of the laser field, which are excited due to the *nonlinear* interaction of the background with the system.

VI. CONCLUSIONS AND OUTLOOK

In this paper, we have used tBLFQ to study the dynamics of an electron interacting with an intense background field, and its consequent emission of radiation (nonlinear Compton scattering). Our chosen background was a model of an axicon laser beam, which has a nontrivial spacetime structure in the directions transverse to the beam propagation. It is by including this structure that we have been able to go beyond previous tBLFQ studies in this area. Notably, the geometrical properties of the basis used in tBLFQ seem particularly well matched with, and thus suitable for the study of, the geometry of the considered backgrounds.

In our analysis of nonlinear Compton scattering we began with the acceleration of the electron in the background, within only the $|e\rangle$ Fock sector included in the calculation; the results agree with a sampling of the classical calculation. Then, in the study of the relative motion, we defined two useful observables. The first, the LMD, gives the probability of finding a scattered electron-photon state at some longitudinal momentum fraction. We saw that the LMD increases with time for every value of the longitudinal momentum fraction, implying that photons can be generated at all possible longitudinal momenta. The second observable, the TMD, gives the distribution of the relative momentum between the electron and the photon in the transverse plane. As time passes, TMDs increase with the center of the distributions moving to higher momentum; this implies that photons are being emitted from the physical electron due to the presence of the background, which naturally leads to an excitation in the relative motion between the photons and the electrons.

We have considered fields which have spacetime dependence on transverse position, and also fields which

depend additionally on (light-front) time; including also a dependence on the remaining longitudinal dependence is required to make the field fully realistic, of course. This provides no real obstacle to tBLFQ, however, other than an increase in numerical complexity, and will be considered elsewhere.

Future research directions include the further development of tBLFQ itself, and the study of other processes [71,72]. For example, here we have included only the first two Fock sectors ($|e\rangle$ and $|e\gamma\rangle$), but in order to obtain a complete picture of e.g. multiphoton emission in Compton scattering, more Fock sectors are needed. Other potential applications of tBLFQ include studying the time-evolution of states in the strong electromagnetic fields generated by heavy ion collisions and studying particle production (the Sauter-Schwinger effect) in strong backgrounds. Developments are also needed in the handling of the center-of-mass motion. In Sec. IV B, we approximately factorized the system in background fields. We further approximated the system by studying only the relative motion. In future improvements, we plan to expand our basis by including the nonzero CM excitations, and thereby we can have an accurate description of the CM motion as well as the relative motion coherently and simultaneously.

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APPENDIX A: LIGHT-FRONT FULL HAMILTONIAN

We start with the QED Lagrangian

$$\mathcal{L} = -\frac{1}{4}F_{\mu\nu}F^{\mu\nu} + \bar{\Psi}(i\gamma^\mu D_\mu - m_e)\Psi, \quad (\text{A1})$$

in which the covariant derivative $D_\mu := \partial_\mu + ie(A_\mu + \mathcal{A}_\mu)$ contains both the dynamical photon field and a background field \mathcal{A} . By taking the Legendre transformation, the full light-front Hamiltonian may be obtained as [20]

$$\begin{aligned}
P^- = & \int d^2x^\perp dx^- \frac{1}{2} \bar{\Psi} \gamma^+ \frac{m_e^2 + (i\partial^\perp)^2}{i\partial^+} \Psi + \frac{1}{2} A^j (i\partial^\perp)^2 A^j + e j^\mu A_\mu + \frac{e^2}{2} J^+ \frac{1}{(i\partial^+)^2} J^+ + \frac{e^2}{2} \bar{\Psi} \gamma^\mu A_\mu \frac{\gamma^+}{i\partial^+} \gamma^\nu A_\nu \Psi \\
& + \frac{e^2}{2} \bar{\Psi} \gamma^\mu A_\mu \frac{\gamma^+}{i\partial^+} \gamma^\nu A_\nu \Psi + \frac{e^2}{2} \bar{\Psi} \gamma^\mu A_\mu \frac{\gamma^+}{i\partial^+} \gamma^\nu A_\nu \Psi + \frac{e^2}{2} \bar{\Psi} \gamma^\mu A_\mu \frac{\gamma^+}{i\partial^+} \gamma^\nu A_\nu \Psi + e j^\mu A_\mu.
\end{aligned} \tag{A2}$$

The first line is the QED light-front Hamiltonian, P_{QED}^- . The second line contains the new terms generated by the background field and is labeled as V .

APPENDIX B: MODE EXPANSIONS OF THE FIELD OPERATORS

We adopt the following mode expansions in the plane-wave basis for fermion and gauge boson fields respectively:

$$\Psi(\mathbf{x}) = \sum_\lambda \int \frac{d^3p}{\sqrt{2}(2\pi)^3} (b(\mathbf{p}, \lambda) u(\mathbf{p}, \lambda) e^{-i\mathbf{p}\cdot\mathbf{x}} + d^\dagger(\mathbf{p}, \lambda) v(\mathbf{p}, \lambda) e^{+i\mathbf{p}\cdot\mathbf{x}}), \tag{B1}$$

$$A_\mu(\mathbf{x}) = \sum_\lambda \int \frac{d^3p}{(2\pi)^3 \sqrt{2p^+}} (a(\mathbf{p}, \lambda) \epsilon_\mu(\mathbf{p}, \lambda) e^{-i\mathbf{p}\cdot\mathbf{x}} + a^\dagger(\mathbf{p}, \lambda) \epsilon_\mu^*(\mathbf{p}, \lambda) e^{+i\mathbf{p}\cdot\mathbf{x}}), \tag{B2}$$

where $\mathbf{p} \cdot \mathbf{x} := \frac{1}{2} p^+ x^- - p^\perp \cdot x^\perp$ is the 3-product for the spatial components of p^μ and x^μ . The creation and annihilation operators obey the (anti)commutation relations

$$[a(\mathbf{p}, \lambda), a^\dagger(\mathbf{p}', \lambda')] = \{b(\mathbf{p}, \lambda), b^\dagger(\mathbf{p}', \lambda')\} = \{d(\mathbf{p}, \lambda), d^\dagger(\mathbf{p}', \lambda')\} = (2\pi)^3 \delta(p^+ - p'^+) \delta^{(2)}(p^\perp - p'^\perp) \delta_\lambda^{\lambda'}. \tag{B3}$$

The single-particle states are

$$|\mathbf{p}, \lambda\rangle_e \equiv b^\dagger(\mathbf{p}, \lambda)|0\rangle \quad \text{and} \quad |\mathbf{p}, \lambda\rangle_\gamma \equiv a^\dagger(\mathbf{p}, \lambda)|0\rangle. \tag{B4}$$

The Dirac spinors are

$$\begin{aligned}
u(\mathbf{p}, \uparrow) &= \begin{pmatrix} 1 \\ 0 \\ \frac{im_e}{p^+} \\ \frac{(ip^\perp - p^2)}{p^+} \end{pmatrix}, & u(\mathbf{p}, \downarrow) &= \begin{pmatrix} 0 \\ 1 \\ \frac{(-ip^\perp - p^2)}{p^+} \\ \frac{im_e}{p^+} \end{pmatrix}, \\
v(\mathbf{p}, \uparrow) &= \begin{pmatrix} 0 \\ 1 \\ \frac{(-ip^\perp - p^2)}{p^+} \\ \frac{-im_e}{p^+} \end{pmatrix}, & v(\mathbf{p}, \downarrow) &= \begin{pmatrix} 1 \\ 0 \\ \frac{-im_e}{p^+} \\ \frac{(ip^\perp - p^2)}{p^+} \end{pmatrix}.
\end{aligned} \tag{B5}$$

The photon polarization vectors are

$$\epsilon^\mu(\mathbf{k}, \lambda) = \left(0, \frac{2\epsilon^\perp(\lambda) \cdot k^\perp}{k^+}, \epsilon^\perp(\lambda) \right), \tag{B6}$$

where $\epsilon^\perp(\uparrow) = \frac{1}{\sqrt{2}}(1, i)$ and $\epsilon^\perp(\downarrow) = \frac{1}{\sqrt{2}}(1, -i)$.

In a space with a compactified longitudinal direction x^- of length $2L$, as in BLFQ, the field operators become

$$\Psi(\mathbf{x}) = \sum_{k,\lambda} \int \frac{d^2p^\perp}{(2\pi)^2 \sqrt{2L}} (b(\mathbf{p}, \lambda) u(\mathbf{p}, \lambda) e^{-i\mathbf{p}\cdot\mathbf{x}} + d^\dagger(\mathbf{p}, \lambda) v(\mathbf{p}, \lambda) e^{+i\mathbf{p}\cdot\mathbf{x}}), \tag{B7}$$

and

$$A_\mu(\mathbf{x}) = \sum_{k,\lambda} \int \frac{d^2p^\perp}{(2\pi)^2 \sqrt{2Lp^+}} (a(\mathbf{p}, \lambda) \epsilon_\mu(\mathbf{p}, \lambda) e^{-i\mathbf{p}\cdot\mathbf{x}} + a^\dagger(\mathbf{p}, \lambda) \epsilon_\mu^*(\mathbf{p}, \lambda) e^{+i\mathbf{p}\cdot\mathbf{x}}), \tag{B8}$$

where $p^+ = \frac{2\pi}{L} k$ with the dimensionless quantity $k = 1, 2, 3, \dots$ for bosons and $k = \frac{1}{2}, \frac{3}{2}, \frac{5}{2}, \dots$ for fermions. Meanwhile, the Dirac delta functions of longitudinal momentum in (B3) should be replaced by the Kronecker delta functions

$$\begin{aligned}
[a(\mathbf{p}, \lambda), a^\dagger(\mathbf{p}', \lambda')] &= \{b(\mathbf{p}, \lambda), b^\dagger(\mathbf{p}', \lambda')\} \\
&= \{d(\mathbf{p}, \lambda), d^\dagger(\mathbf{p}', \lambda')\} \\
&= (2\pi)^2 \delta_{\mathbf{p}^+}^{\mathbf{p}'^+} \delta^{(2)}(p^\perp - p'^\perp) \delta_\lambda^{\lambda'}. \quad (\text{B9})
\end{aligned}$$

APPENDIX C: BLFQ HARMONIC OSCILLATOR BASIS

In the transverse plane, the BLFQ basis adopts the 2D-HO states, which are the eigenstates of the 2D-HO Hamiltonian

$$H_{\text{HO}} = \frac{(p^\perp)^2}{2M} + \frac{1}{2}M\Omega^2(x^\perp)^2, \quad (\text{C1})$$

where M and Ω are the mass and the frequency of the oscillator respectively. In coordinate space, the eigenfunction can be factorized into a radial and an angular function

$$\Phi_{nm}^b(\rho, \phi) = (-1)^n i^{|m|} f_{nm}^b(\rho) \chi_m(\phi). \quad (\text{C2})$$

The corresponding eigenvalue is $E_{n,m} = (2n + |m| + 1)\Omega$, in which n and m are integers; $n \geq 0$ characterizes the radial excitation and m describes the angular momentum of the oscillator; (ρ, ϕ) are polar coordinates in the transverse plane with $x^1 = \rho \cos \phi$ and $x^2 = \rho \sin \phi$. The explicit form of the radial function is

$$f_{nm}^b(\rho) = b\sqrt{2} \sqrt{\frac{n!}{(n+|m|)!}} e^{-b^2\rho^2/2} (b\rho)^{|m|} L_n^{|m|}(b^2\rho^2), \quad (\text{C3})$$

where $L_n^{|m|}(b^2\rho^2)$ is the generalized Laguerre polynomial and $b := \sqrt{M\Omega}$ is the 2D-HO scale parameter. The angular function is

$$\chi_m(\phi) = \frac{1}{\sqrt{2\pi}} e^{im\phi}. \quad (\text{C4})$$

The 2D-HO eigenfunction (C2) is orthonormalized such that

$$\int_0^\infty \int_0^{2\pi} d\rho d\phi \Phi_{nm}^{b*}(\rho, \phi) \Phi_{n'm'}^b(\rho, \phi) = \delta_n^{n'} \delta_m^{m'}.$$

$$\begin{aligned}
\langle p'^\perp, p'^+, \lambda' | V | p^\perp, p^+, \lambda \rangle &= \int dx^- d^2x^\perp \frac{1}{2} e\mathcal{A}^-(x) \langle p'^\perp, p'^+, \lambda' | \bar{\Psi}(\mathbf{x}) \gamma^+ \Psi(\mathbf{x}) | p^\perp, p^+, \lambda \rangle \\
&= \int dx^- d^2x^\perp \frac{m_e a_0}{2b_l} \sum_{ss' ll'} \int \frac{d^2q'^\perp d^2q^\perp}{2L(2\pi)^4} \Phi_{00}^{b_l}(x^\perp) e^{i(\mathbf{q}' - \mathbf{q}) \cdot x} \\
&\quad \times \bar{u}(\mathbf{q}', s') \gamma^+ u(\mathbf{q}, s) \langle p'^\perp, p'^+, \lambda' | b^\dagger(\mathbf{q}', s') b(\mathbf{q}, s) | p^\perp, p^+, \lambda \rangle \\
&= \frac{m_e a_0}{b_l} \delta_k^k \delta_\lambda^{\lambda'} \tilde{\Phi}_{00}^{b_l}(p'^\perp - p^\perp). \quad (\text{D2})
\end{aligned}$$

In order to transform the interaction Hamiltonian to the BLFQ basis which adopts HO eigenstates in the transverse plane, we convolute (D2) with the HO wave functions

The momentum-space eigenfunction can be obtained by a Fourier transform

$$\begin{aligned}
\tilde{\Phi}_{nm}^b(p^\perp) &= \int d^2x^\perp e^{-i\tilde{x}^\perp \cdot \tilde{p}^\perp} \Phi_{nm}^b(x^\perp) \\
&= (2\pi) \tilde{f}_{nm}^b(p) \tilde{\chi}_m(\phi), \quad (\text{C5})
\end{aligned}$$

in which

$$\tilde{f}_{nm}^b(p) = \frac{\sqrt{2}}{b} \sqrt{\frac{n!}{(n+|m|)!}} e^{-p^2/(2b^2)} \left(\frac{p}{b}\right)^{|m|} L_n^{|m|}\left(\frac{p^2}{b^2}\right), \quad (\text{C6})$$

and

$$\tilde{\chi}_m(\phi) = \frac{1}{\sqrt{2\pi}} e^{im\phi}. \quad (\text{C7})$$

APPENDIX D: MATRIX ELEMENTS OF THE INTERACTION HAMILTONIAN OF THE BACKGROUND FIELD

As in Eq. (16), only the longitudinal component of the background field \mathcal{A} is nonzero. From Eq. (A2), the only surviving term in the Hamiltonian involving the background is the last term

$$V = \int dx^- d^2x^\perp e j_- \mathcal{A}^-, \quad (\text{D1})$$

where $j_- = \bar{\Psi} \gamma_- \Psi = \frac{1}{2} \bar{\Psi} \gamma^+ \Psi$ is the fermion current. Note that the background field is treated as a classical field; therefore the photons are spectators and the nontrivial part of the interaction comes from the fermion fields. For now, we neglect the photon part and keep in mind that the matrix elements of the photon part contribute Kronecker delta functions conserving all the 4 quantum numbers for the spectator photons. In the momentum representation, the matrix elements of the interaction Hamiltonian can be straightforwardly obtained by using the commutation relations (B9)

$$\begin{aligned}\langle \bar{\alpha}'_e | V | \bar{\alpha}_e \rangle &= \int \frac{d^2 p'^{\perp} d^2 p^{\perp}}{(2\pi)^4} (\tilde{\Phi}_{n'm'}^{\sqrt{x}b}(p'^{\perp}))^* \langle p'^{\perp}, p'^+, \lambda' | V | p^{\perp}, p^+, \lambda \rangle \tilde{\Phi}_{nm}^{\sqrt{x}b}(p^{\perp}) \\ &= \frac{m_e a_0}{b_l} \delta_k^{\lambda'} \delta_{\lambda}^{\lambda'} \int \frac{d^2 p'^{\perp} d^2 p^{\perp}}{(2\pi)^4} \tilde{\Phi}_{00}^{b_l}(p'^{\perp} - p^{\perp}) (\tilde{\Phi}_{n'm'}^{\sqrt{x}b}(p'^{\perp}))^* \tilde{\Phi}_{nm}^{\sqrt{x}b}(p^{\perp}),\end{aligned}\quad (\text{D3})$$

in which the x (longitudinal momentum fraction) dependent HO basis is adopted, for exact factorization in a finite truncated BLFQ basis.

We now recover the photon part: for the matrix elements in the $|\epsilon\gamma\rangle$ sector, we augment (D3) with the Kronecker delta functions conserving photon quantum numbers and the matrix elements become

$$\langle \bar{\alpha}'_{e\gamma} | V | \bar{\alpha}_{e\gamma} \rangle = \langle \bar{\alpha}'_e | V | \bar{\alpha}_e \rangle \delta_{\alpha'_\gamma}^{\alpha_\gamma}. \quad (\text{D4})$$

Because there is no cross term between the $|e\rangle$ and $|\epsilon\gamma\rangle$ sectors, the full interaction Hamiltonian matrix is the direct sum of the interaction Hamiltonian matrix in these two sectors. Finally, the interaction Hamiltonian matrix in BLFQ basis $|\alpha\rangle$ can be transformed to the tBLFQ basis $|\beta\rangle$, using the wave function $\langle \alpha | \beta \rangle$ obtained by diagonalizing P_{QED}^- [54]

$$\langle \beta' | V | \beta \rangle = \sum_{\alpha\alpha'} \langle \beta' | \alpha' \rangle \langle \alpha' | V | \alpha \rangle \langle \alpha | \beta \rangle. \quad (\text{D5})$$

APPENDIX E: PERTURBATIVE CALCULATION OF PHOTON EMISSION IN A WEAK BACKGROUND FIELD

In order to validate our tBLFQ results we will compare them with an established method of calculation, namely perturbation theory. Since perturbation theory is valid for weak coupling, we focus here on the case of a weak background field with the dimensionless parameter $a_0 = 0.1$ in Eq. (16). We present below a comparison between tBLFQ and perturbation theory calculations of the TMDs (26) introduced in the text, for Compton scattering.

The natural objects to study in perturbation theory are scattering amplitudes between asymptotic free states, rather than the finite time-evolution of interacting states as in tBLFQ. In order to aid the comparison of tBLFQ with perturbation theory, we therefore switch off our background Eq. (16) outside of light-front time $0 < x^+ < \tau$:

$$e\mathcal{A}^\mu(x, \tau) = \delta_-^\mu \frac{m_e a_0}{b_l} \Phi_{00}^{b_l}(x^+) [\theta(x^+) - \theta(x^+ - \tau)]. \quad (\text{E1})$$

In light-front perturbation theory, up to order e^2 , there are four Feynman diagrams which contribute to Compton scattering in a classical background field, see Fig. 14. In our normalization, the T -matrix element is

$$\begin{aligned}\langle p', s'; k', \lambda | iT(\tau) | p, s \rangle &= \frac{-ie^2 \epsilon_\mu^*(\mathbf{k}', \lambda)}{2\sqrt{2k'^+}} \frac{1}{2} \tilde{\mathcal{A}}_\nu(p' + k' - p, \tau) \bar{u}(\mathbf{p}', s') \left[\frac{\gamma^\mu (\not{p}' + \not{k}' + m) \gamma^\nu}{(p' + k')^2 - m^2} + \frac{\gamma^\nu (\not{p}' - \not{k}' + m) \gamma^\mu}{(p - k')^2 - m^2} \right] u(\mathbf{p}, s) \\ &= (2\pi) \delta(p^+ - p'^+ - k'^+) i\mathcal{M}(p \rightarrow p', k'),\end{aligned}\quad (\text{E2})$$

in which $\tilde{\mathcal{A}}(k, \tau)$ is the Fourier transform of $\mathcal{A}(x, \tau)$

$$\tilde{\mathcal{A}}(k, \tau) = \int dx^+ dx^- d^2 x^\perp \mathcal{A}(x, \tau) e^{ikx}. \quad (\text{E3})$$

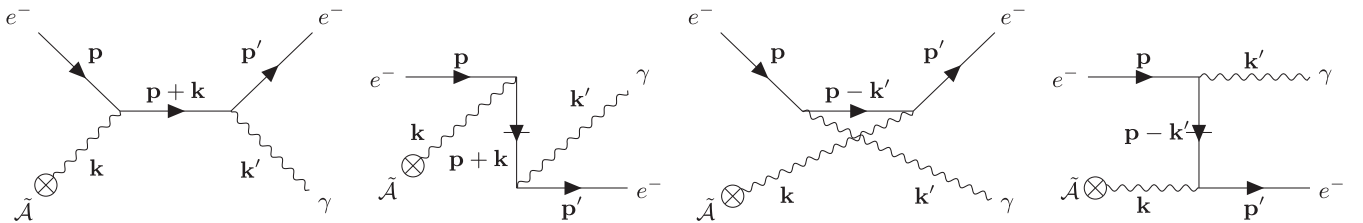


FIG. 14. Compton scattering in light-front perturbation theory. Time flows from left to right. The second and fourth terms are instantaneous-fermion contributions. The sum of the diagrams gives the same matrix element as in standard covariant perturbation theory.

around momentum p^+ , and a momentum eigenstate in the transverse plane

$$|\phi, p^\perp\rangle = \int \frac{dq^+}{2\pi} \phi(q^+) |q^+, p^\perp\rangle. \quad (\text{E4})$$

The wave packet is normalized such that

$$\langle \phi, p^\perp | \phi, p'^\perp \rangle = (2\pi)^2 \delta^{(2)}(p^\perp - p'^\perp) \quad \text{and} \quad \int \frac{dq^+}{2\pi} |\phi(q^+)|^2 = 1. \quad (\text{E5})$$

The differential probability for the initial state to scatter to a final electron-photon state with momenta in the ranges $d^3 p'$, $d^3 k'$, respectively, is

$$\begin{aligned} \mathcal{P}(e^- \rightarrow e^- \gamma) &= \frac{d^3 p'}{(2\pi)^3} \frac{d^3 k'}{(2\pi)^3} \frac{1}{(2\pi)^2} \int \frac{dq^+}{2\pi} \int \frac{dq'^+}{2\pi} \phi(q^+) \phi^*(q'^+) \mathcal{M}(q^+; p^\perp \rightarrow p', k') \\ &\quad \times \mathcal{M}^*(q'^+; p^\perp \rightarrow p', k') (2\pi) \delta(q^+ - k'^+ - p'^+) (2\pi) \delta(q'^+ - k'^+ - p'^+) \\ &= \frac{d^3 p'}{(2\pi)^3} \frac{d^3 k'}{(2\pi)^3} \frac{1}{(2\pi)^2} \int \frac{dq^+}{2\pi} |\phi(q^+)|^2 |\mathcal{M}(q^+; p^\perp \rightarrow p', k')|^2 (2\pi) \delta(q^+ - k'^+ - p'^+) \\ &\approx \frac{d^3 p' d^3 k'}{(2\pi)^8} |\mathcal{M}(p \rightarrow p', k')|^2 (2\pi) \delta(p^+ - p'^+ - k'^+). \end{aligned} \quad (\text{E6})$$

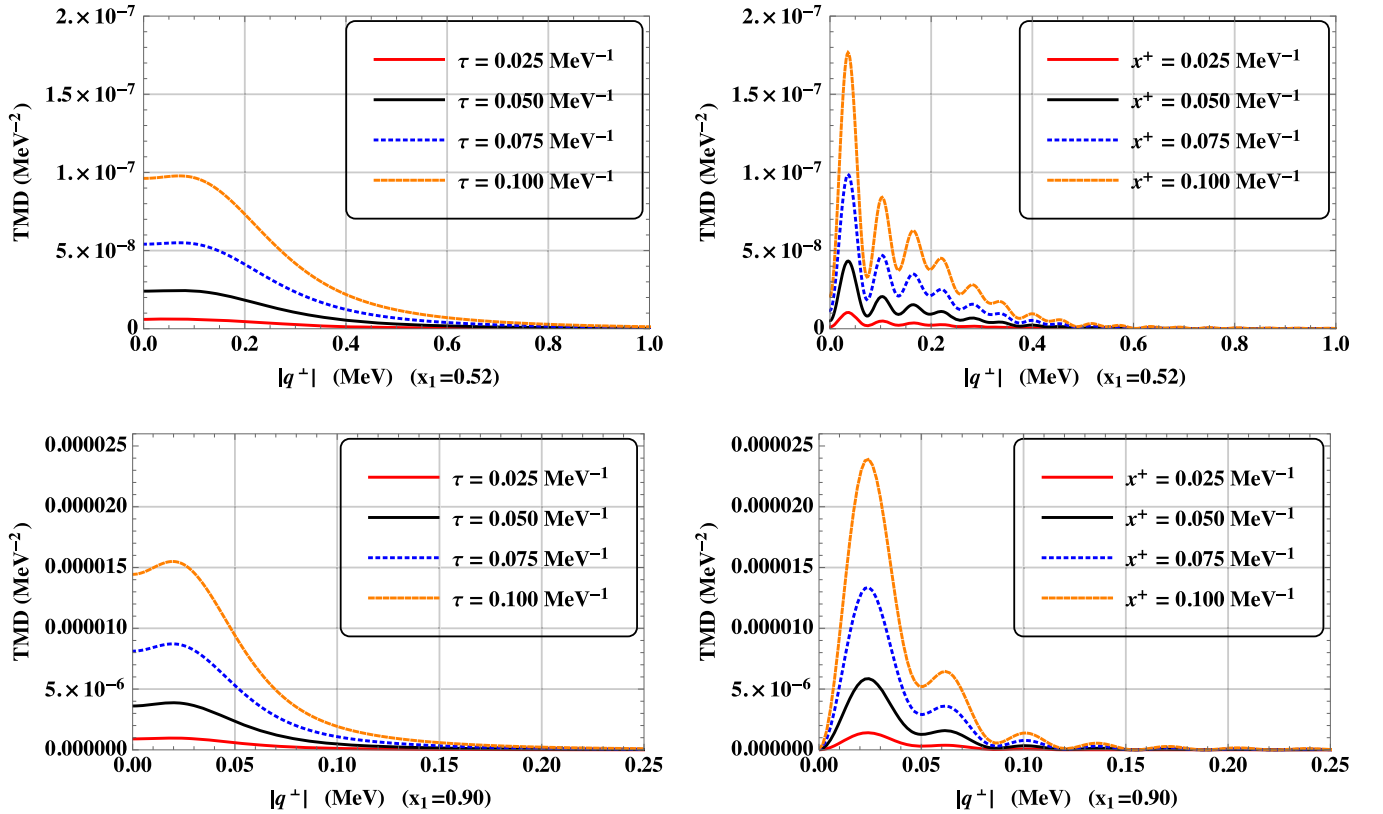


FIG. 15. TMDs in perturbation theory and tBLFQ. In all plots, $a_0 = 0.1$, $b_l = m_e$, $P^+ = 10.5$ MeV; in perturbation theory, the momentum of the incident electron is $\mathbf{p} \equiv (p^+, p^\perp) = (10.5, 0, 0)$ MeV. In the upper row, $x_1 = 0.52$ and in the lower row, $x_1 = 0.90$. The left-hand column shows TMDs obtained in perturbation theory, with 4 different values of parameter τ , as indicated in the figure; the right-hand column shows the TMD time-evolution obtained in tBLFQ, in basis with $N_{\max} = 100$ and $K_{\text{total}} = 10.5$. TMDs increase with x^+ and τ in all the cases and show similar shapes in two approaches (noting the different scales on both the horizontal and vertical axes) when $x^+ = \tau$ if we neglect the oscillations introduced by the HO basis. Overall the tBLFQ results are slightly smaller than the perturbative results, which is discussed in the main text.

As in tBLFQ, we wish to separate the CM and the relative motion. To do so we rewrite the single-particle momenta p'^{\perp} and k'^{\perp} in terms of the CM momentum $P^{\perp} = p'^{\perp} + k'^{\perp}$ and the relative momentum $q^{\perp} = x_2 p'^{\perp} - x_1 k'^{\perp}$. For consistency and simplicity, we make the same assumption as in Sec. IV B, that the amplitude (E2) can be approximately factorized. The relative probability can be obtained by integrating out the CM momentum P^{\perp} in Eq. (E6) and the TMD can be obtained by summing over spins and helicity as in Eq. (26)

$$\text{TMD}(q^{\perp}, x_1; \tau) = \frac{1}{2} \sum_{s, s', \lambda} \int \frac{p^+ dk'^+ d^2 P^{\perp}}{(2\pi)^6} |\mathcal{M}(p \rightarrow p', k')|^2 \times (2\pi) \delta(p'^+ + k'^+ - p^+). \quad (\text{E7})$$

The calculation of the spin and polarization sums in Eq. (E7) proceeds as in textbook derivations so we do not repeat it. Figure 15 compares TMDs obtained from

perturbation theory, Eq. (E7), and from tBLFQ. Four different field durations, τ , are shown for the perturbative results, compared against tBLFQ results evolved from states evolved through corresponding times x^+ . As shown in Fig. 15, TMDs obtained in two approaches increase with x^+ (and τ), indicating photons are being emitted and excited, and show broadly similar shapes and amplitudes when $x^+ = \tau$. The tBLFQ results show additional oscillations induced by the HO basis. Overall, the tBLFQ results are slightly smaller than the perturbative results. This discrepancy may be attributed to the fact that the CM motion of the initial state in our tBLFQ calculation is a Gaussian wave packet in the transverse plane (21), rather than a momentum eigenstate as in the perturbative calculation. Another main difference in Fig. 15 is that, as $q^{\perp} \rightarrow 0$, TMDs in two approaches show different limits. In this region, which is however below the IR cutoff implied by the HO basis ($\sqrt{x_1 x_2} b / \sqrt{N_{\max}} \approx 0.015 \text{ MeV}$ when $N_{\max} = 100$, $x_1 = 0.9$), the tBLFQ results are not reliable.

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