# Investigation of two-photon emission in strong field QED using channeling in a crystal

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(Received 16 May 2019; published 12 August 2019)

We investigate the second order process of two photons being emitted by a high-energy electron dressed in the strong background electric field found between the planes in a crystal. The strong crystalline field combined with ultrarelativistic electrons is one of very few cases where the Schwinger field can be experimentally achieved in the electron's rest frame. The radiation being emitted, the so-called channeling radiation, is a well studied phenomenon. However, only the first order diagram corresponding to emission of a single photon has been studied so far. We elaborate on how the two-photon emission process should be understood in terms of a two-step versus a one-step process, i.e., if one can consider one photon being emitted after the other, or if there is also a contribution where the two photons are emitted "simultaneously." From the calculated full probability we see that the two-step contribution is simply the product of probabilities for single photon emission while the additional one-step terms are, mainly, interferences due to several possible intermediate virtual states. These terms can contribute significantly when the crystal is thin. Therefore, in addition, we see how one can, for a thick crystal, calculate multiple photon emissions quickly by neglecting the one-step terms, which represents a solution of the problem of quantum radiation reaction in a crystal beyond the usually applied constant field approximation. We explicitly calculate an example of 180 GeV electrons in a thin silicon crystal and argue why it is, for experimental reasons, more feasible to see the one-step contribution in a crystal experiment than in a laser experiment.

DOI: 10.1103/PhysRevD.100.036002

### I. INTRODUCTION

Strong field QED is the study of physical processes that take place in a strong background field, and nonlinear effects of quantum nature arise when the size of the Lorentz invariant parameter

$$\chi = e \sqrt{(F_{\mu\nu}p^{\nu})^2}/m^3 \tag{1}$$

is on the order of unity, which is the ratio of the electromagnetic field experienced in the electron's rest frame compared to the Schwinger field strength  $E_{\rm Sch} = 1.32 \times 10^{18}$  V/m. Here *e* is the elementary charge, *m* the electron mass,  $F_{\mu\nu}$  the electromagnetic field tensor of the background field, and  $p^{\nu}$  the electron 4-momentum. We use natural units such that  $\hbar = c = 1$ ,  $\alpha = e^2$ . Lindhard was one of the first to realize that when high-energy charged particles are aimed close to the direction along an axis or plane in a crystal, the charged particle can

become transversely trapped [1]. Later it was studied how this motion leads to radiation emission called channeling radiation, especially relevant for electrons and positrons. This is well-studied both experimentally [2–10] and theoretically [5,11–15]. Crystal channeling represents one of the only phenomena where the Schwinger field can be experimentally achieved in the electron's rest frame [6,16–18], with the only other example being the famous E-144 SLAC experiment on nonlinear Compton scattering [19] using relativistic electrons colliding with a laser beam. Crystals with ultrarelativistic electrons or positrons therefore present a unique possibility to study physics in such strong fields. However, a calculation from first principles of



FIG. 1. The Feynman diagrams corresponding to the process under study. The double fermion lines correspond to positron solutions of the Dirac equation in the background field of the interplanar crystal potential.

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emission of more than one photon has not been carried out for crystal channeling (see Fig. 1). The recent studies of two-photon emission in the collision of relativistic electrons with a laser pulse [20-23] show that the emission of two photons is not exactly the product of probabilities for each emission; however, under certain conditions it is an acceptable approximation. The experimental verification of such results are, however, complicated in the case of the laser pulse colliding with an electron bunch because any two (or more) emitted photons cannot be known to be emitted by the same electron. In crystal experiments as in, e.g., [18], it is standard that each incoming particle is recorded as a separate event, and therefore the measured outgoing photons are sure to stem from the single incoming particle. Therefore, in this paper, we will calculate the emission of two photons during electron channeling in a crystal, which could potentially be studied experimentally in an experiment similar to the one seen in [18]; however, with a modified setup to allow for the detection of an additional photon. For the theory of channeling radiation, in particular, the development of the semiclassical operator method by Baier et al. [24] stands out and has been extensively applied to the phenomenon of channeling [25]. This method allowed the inclusion of quantum effects such as the electron spin and the photon recoil, which are important when  $\chi$  is no longer small, while needing only the classical trajectory of the electron/positron in the external field. The authors of this method, seeking analytical results, in most applications to channeling, applied the approximation of the local constant field, which greatly simplifies calculations. The constant field approximation means that while a particle moves in an external field, which is not constant, one applies the result of the constant field formula locally, i.e., in a small time step. Effectively this means neglecting that the radiation emitted before or after can interfere with this radiation. This is valid only for certain parameters of fields and particle energies. However, the semiclassical operator method can be used to calculate the radiation emission under general circumstances without much effort, also when the constant field approximation is no longer valid [26,27], which with modern computing power makes it one of the most powerful methods to calculate the radiation emitted by ultrarelativistic electrons in a general field configuration. There are caveats, however, which are twofold. First, the notion of a classical trajectory should make sense. Or, in other words, the quantum numbers associated with the motion should be large, a subject recently studied in [28,29]. Second, the derivation starts out from the first-order diagram of a dressed electron emitting a single photon. Therefore the emission rate of two, or more, photons cannot be predicted by this method without approximations. The emission of a single photon yields a rate, an emission probability per unit time, and as such one can construct the probability for emitting several photons by applying this rate for each consecutive

emission. In this way, the probability to emit, e.g., two photons would be proportional to time, or thickness of the crystal, squared, and so on. We will call this process the "cascade" process. Herein lies an approximation, where interference between different emissions is neglected. We show that the two-photon emission probability contains the cascade along with one-step terms that scale linearly with the crystal thickness. Therefore, for sufficiently thin crystals, these one-step terms will become important. This phenomenon is also discussed in pair production of electron/positron pairs from high-energy photons in a strong field where one also distinguishes between the two-step and the one-step, or "trident," process. This has been investigated in crystals in [17] and has received renewed interest with the prospect of studying such phenomena in high-intensity laser fields [30-35]. In this paper we make quantitative calculations of the angularly integrated probability, differential in photon energies, of the emission of two photons by an electron in the planar Doyle-Turner potential [25,36–38]. We do this by finding numerical solutions of the Dirac equation by solving the problem in a basis of plane waves, which is possible due to the periodicity of the transverse potential in a crystal, as shown in [28]. Whether the cascade terms are enough to properly describe the radiation emission is a highly relevant question as it closely relates to the phenomenon of quantum radiation reaction, the emission of multiple photons when  $\gamma$  is large [39], recently studied using channeling radiation as well as in laser experiments [18,40,41]. In the crystal experiment it was seen that even for energies as high as 180 GeV positrons, where it could be expected that the constant field approximation would be acceptable, it was shown that discrepancies arise due to this, and therefore a more general theory was called for. The current theory of quantum radiation reaction in lasers relies on the local constant field approximation [25,39,42-46], and it is unknown if one can calculate the emission of many photons in a way that avoids calculating all the corresponding higher order diagrams, when going beyond the constant field approximation. This question will be addressed in the case of a crystal, in the current paper.

We use the Feynman slash notation such that  $\not = a_{\mu}\gamma^{\mu}$ , where  $\gamma^{\mu}$  are the Dirac gamma matrices and  $a^{\mu}$  an arbitrary four-vector. We adopt the metric tensor  $\eta^{\mu\nu} = \text{diag}(+1, -1, -1, -1)$ .

### **II. FORMALISM**

In QED the transition amplitude from a given initial state  $|i\rangle$  to a final state  $|f\rangle$  is given by

$$S_{fi} = \langle f | U(\infty, -\infty) | i \rangle, \tag{2}$$

where U is the time evolution operator, often written as  $U(\infty, -\infty) = \mathcal{T} \exp(-i \int_{-\infty}^{\infty} V(t) dt)$  where  $\mathcal{T}$  is the timeordering operator and  $V(t) = \int e \bar{\Psi} \mathcal{A} \Psi d^3 x$  is the quantized interaction. We then write our quantized fields as

$$\Psi = \sum_{s=1}^{2} \int \frac{d^3 p}{(2\pi)^3} [b_p^s \psi_{p,s}^-(x) + c_p^{s\dagger} \psi_{p,s}^+(x)], \quad (3)$$

$$A = \int \frac{d^3k}{(2\pi)^3} \sqrt{\frac{4\pi}{2\omega}} \sum_{r=1}^{2} [e_r a_k^r e^{-ikx} + e_r^* a_k^{r\dagger} e^{ikx}], \quad (4)$$

where  $\psi_{p,s}^{-}(x)$  and  $\psi_{p,s}^{+}(x)$  are an orthonormal and complete set of electron and positron solutions, respectively, in the background field.  $\int \frac{d^3p}{(2\pi)^3}$  denotes a summation over all states, and p the relevant quantum numbers that we will find later. The b, c, and a operators are the annihilation operators of the electron, positron, and photon fields, respectively, obeying the relations that the only nonzero (anti)commutators are  $\{b_p^r, b_q^{s\dagger}\} = \{c_p^r, c_q^{s\dagger}\} = [a_p^r, a_q^{s\dagger}] = (2\pi)^3 \delta^{rs} \delta^{(3)}(\mathbf{p} - \mathbf{q})$ , where the  $\{\}$  brackets denote the anticommutator and [] the commutator.

In [28,29] we discussed the Dirac equation with the potential found in the crystal, but we will here repeat the results we need in order to calculate the emission of two photons. It was found in [28] that the electron solution can be written as follows:

$$\psi^{-}(x) = \frac{1}{\sqrt{2\varepsilon}} e^{i(p_x x + p_z z - \varepsilon t)} U(y), \tag{5}$$

the positron solutions can then be written as (see the Appendix A)

$$\psi^{+}(x) = \frac{1}{\sqrt{2\varepsilon}} e^{-i(p_x x + p_z z - \varepsilon t)} V(y), \tag{6}$$

and U and V are given by

$$U(y) = \sqrt{\varepsilon + m} \left(\frac{s^{-}}{\frac{\sigma \tilde{p}}{\varepsilon + m}} s^{-}\right) I^{-}(y),$$
$$V(y) = \sqrt{\varepsilon + m} \left(\frac{\frac{\sigma \tilde{p}}{\varepsilon + m}}{s^{+}} s^{+}\right) I^{+}(y),$$
(7)

where  $\tilde{\boldsymbol{p}} = (p_x + q\varphi(y), \operatorname{sign}(q)i\frac{d}{dy}, p_z), \quad \varphi(y)$  is the electrostatic potential,  $q = \pm e$  is the charge, the superscript on I(y) refers to the charge sign, and *s* is a two component vector describing the spin, which we can choose as either  $(1 \ 0)^T$  or  $(0 \ 1)^T$ , corresponding to spin-up and spin-down, respectively, for the electron, and opposite for the positron. From the choice of the form of the spinors *U* and *V*, it is also clear that  $\varepsilon$  positive should be used (see Appendix A). I(y) is the solution to the equation

$$\left[-\frac{1}{2\varepsilon}\frac{d^2}{dy^2} + q\varphi(y)\right]I(y) = \frac{\varepsilon^2 - p_x^2 - p_z^2 - m^2}{2\varepsilon}I(y).$$
 (8)

For  $\varphi(y)$  we will use the Doyle-Turner model [25,36–38], chosen as symmetric around 0. In a crystal this potential  $\varphi(y)$  is periodic with the period of the interplanar distance,

which we will denote as  $d_p$ . Because of this, the solution (for the electron) can be written as a Bloch wave such that

$$I^{-}(y) = e^{ik_{B}^{-}y}u_{k_{B}}^{-}(y), \qquad (9)$$

where  $u_{k_B^-}^-(y)$  is also periodic with period  $d_p$  and  $k_B^-$  is the Bloch momentum, which can be taken to be in the interval  $0 \le k_B^- < k_0, k_0 = \frac{2\pi}{d_p}$ . It then follows from Bloch's theorem that these solutions form an orthogonal and complete set of solutions of Eq. (8). Inserting  $I^-(y)$  of Eq. (9) into Eq. (8) gives us the equation governing  $u_{k_p}^-(y)$ ,

$$\left(-\frac{1}{2\varepsilon}\left[\frac{d^2}{dy^2} + 2ik_B\frac{d}{dy} - k_B^2\right] + q\varphi(y)\right)u_{k_B}^-(y)$$
$$= \frac{\varepsilon^2 - p_x^2 - p_z^2 - m^2}{2\varepsilon}u_{k_B}^-(y). \tag{10}$$

The periodicity of  $u_{k_B}(y)$  means it can be written as a Fourier series,

$$u_{k_B}^{-}(y) = \sum_{j} c_j e^{ijk_0 y}.$$
 (11)

To ensure normalization we should have  $\sum_j |c_j|^2 = 1$  (see Appendix B of Ref. [28]). It is now clear that this is an eigenvalue problem for each  $k_B$  where the quantized eigenvalue is

$$E_n = \frac{\varepsilon^2 - p_x^2 - p_z^2 - m^2}{2\varepsilon},$$
(12)

where *n* is the quantum number corresponding to the value of this energy in ascending order and where 0 is the ground state. This equation leads to a quantization of, e.g.,  $p_x$ . The coefficients  $c_j$  are found by solving the matrix eigenvalue problem obtained by inserting Eq. (11) into Eq. (10), multiplying with  $\frac{1}{d_p}e^{-ilk_0y}$ , and integrating over *y* from 0 to  $d_p$  to exploit orthogonality

$$\sum_{j} \frac{1}{2\varepsilon} [jk_0 + k_B]^2 \delta_{j,l} c_j$$
$$+ \sum_{j} c_j \frac{1}{d_p} \int q\varphi(y) e^{i(j-l)k_0 y} dy$$
$$= \sum_{j} \frac{\varepsilon^2 - p_x^2 - p_z^2 - m^2}{2\varepsilon} \delta_{j,l} c_j.$$
(13)

This was done with the electron function  $I^{-}(y)$  in mind, but the positron coefficients can be obtained just by changing q. With these things taken into consideration, we now see that we can write the U(y) and V(y) functions in terms of the coefficients  $c_j$  such that

$$U(y) = \sum_{j} c_{j}^{-} \boldsymbol{S}_{j}^{-} e^{i(jk_{0}+k_{B})y}, \qquad (14)$$

$$V(y) = \sum_{j} c_{j}^{+} S_{j}^{+} e^{-i(jk_{0} + k_{B})y}, \qquad (15)$$

where

$$\mathbf{S}_{j}^{-} = \sqrt{\varepsilon + m} \begin{pmatrix} \mathbf{s}^{-} \\ \frac{\sigma \cdot \mathbf{p}_{j}}{\varepsilon + m} \mathbf{s}^{-} \end{pmatrix}, \tag{16}$$

$$S_{j}^{+} = \sqrt{\varepsilon + m} \left( \frac{\frac{\sigma p_{j}}{\varepsilon + m} S^{+}}{S^{+}} \right), \qquad (17)$$

where  $p_j = (p_x + E_n - \frac{(jk_0+k_B)^2}{2\varepsilon}, jk_0 + k_B, p_z)$ . For the calculation of radiation emission from electrons we will need the quantity  $\bar{S}_f \notin^* S_i^-$ , where we have put labels for the initial state *i* and final state *f*; however, these still each depend on the index *j*. This quantity can then be written as

$$\bar{S}_{f} \not \in S_{i}^{-} = -s_{f}^{T} [\boldsymbol{\epsilon} \cdot \boldsymbol{A} + i\boldsymbol{B} \cdot \boldsymbol{\sigma}] s_{i}, \qquad (18)$$

where

$$\boldsymbol{A} = \sqrt{\frac{\varepsilon_f + m}{\varepsilon_i + m}} \boldsymbol{p}_i + \sqrt{\frac{\varepsilon_i + m}{\varepsilon_f + m}} \boldsymbol{p}_f, \qquad (19)$$

$$\boldsymbol{B} = \boldsymbol{\epsilon}^* \times \left( \sqrt{\frac{\varepsilon_f + m}{\varepsilon_i + m}} \boldsymbol{p}_i - \sqrt{\frac{\varepsilon_i + m}{\varepsilon_f + m}} \boldsymbol{p}_f \right).$$
(20)

Now since we have an orthonormal complete set of solutions, we can write the propagator in terms of these states as [47]

$$G(x_2, x_1) = \int \frac{d^3 p}{(2\pi)^3} \sum_{n,s} + \theta(t_2 - t_1) e^{-i\varepsilon(t_2 - t_1)} \psi_{p,n,s}^-(\mathbf{x}_2) \bar{\psi}_{p,n,s}^-(\mathbf{x}_1) \\ - \theta(t_1 - t_2) e^{i\varepsilon(t_2 - t_1)} \psi_{p,n,s}^+(\mathbf{x}_2) \bar{\psi}_{p,n,s}^+(\mathbf{x}_1).$$
(21)

This expression can be simplified due to the simple expression for the wave functions in all coordinates but the *y* coordinate. However, we will not carry this out, as it is easier to see how the cascade part of the radiation emission arises by starting from the above expression.

### **III. SINGLE PHOTON EMISSION AND CASCADE**

We will now briefly mention some results obtained in [28] on the single photon emission probability which is relevant to build the expected cascade contribution. We found that the rate of emission is given by

$$dW_{i\to f}^{(1)} = \frac{1}{(2\pi)^2} |\mathcal{M}_{i\to f}|^2 \delta(\varepsilon_f + \omega - \varepsilon_i) d^3k, \quad (22)$$

where we defined

$$\mathcal{M}_{i\to f} = e \sqrt{\frac{4\pi}{2\omega}} \frac{1}{2\sqrt{\varepsilon_f \varepsilon_i}} \sum_j c^*_{n_B+j,f} c_{j,i} \bar{\mathbf{S}}^-_{n_B+j,f} \boldsymbol{\ell}^* \mathbf{S}^-_{j,i}, \quad (23)$$

where  $n_B$  is the integer such that  $0 \le k_{B,f} < k_0$ , where  $k_{B,f} = k_{B,i} - k_y - n_B k_0$ ,  $S_{j,i}$  corresponds to the initial state, and  $c_{j,i}$  is a coefficient with index *j* corresponding to the initial state *i*. See the Appendix of [28] for the details on why  $\mathcal{M}$  reduces to a single sum over *j*. As shown in [28] there are large terms in  $\varepsilon_f + \omega - \varepsilon_i$  that cancel, leaving behind the relevant small terms, because the relevant transverse energies  $E_n$ , comparable to the potential depth, are much smaller than the whole particle energy, i.e., eV versus GeV. We could rewrite the content of the delta function as

$$f(\theta) = \varepsilon_f + \omega - \varepsilon_i$$
  

$$\simeq E_{n_f} - E_{n_i} + \frac{m^2}{2\varepsilon_f} - \frac{m^2}{2\varepsilon_i} + \frac{\omega\theta^2}{2} \left(1 + \frac{\omega\sin^2\varphi}{\varepsilon_f}\right). \quad (24)$$

Now we may use that  $\delta(\varepsilon_f + \omega - \varepsilon_i) = \frac{1}{|f'(\theta_0)|} \delta(\theta - \theta_0)$ where  $\theta_0$  is the positive solution to  $f(\theta) = 0$ . From the formula for single photon emission, Eq. (22), we can construct the cascade contribution to two-photon emission. We wish to know the probability of finding a photon in the momentum interval  $d^3k_1$  around  $k_1$  while also finding a photon within another interval  $d^3k_2$  around  $k_2$ . This can happen in two ways, either the particle emits  $k_1$  while transitioning from the initial state, and then subsequently  $k_2$ or vice versa. We are, however, interested in the angular integrated spectrum, that is  $dP_{i \to f}^{(cascade)}/d\omega_1 d\omega_2$ , and therefore an additional factor of  $\frac{1}{2}$  must be added due to counting the same point in phase space twice [48]; and so we obtain

$$\frac{dP_{i \to f}^{(\text{cascade})}}{d\omega_1 d\omega_2} = \frac{T^2}{2} \sum_{v} \frac{1}{2} \left[ \frac{dW_{i \to v}^{(1)}}{d\omega} (\omega_1) \frac{dW_{v \to f}^{(1)}}{d\omega} (\omega_2) + \frac{dW_{i \to v}^{(1)}}{d\omega} (\omega_2) \frac{dW_{v \to f}^{(1)}}{d\omega} (\omega_1) \right].$$
(25)

#### **IV. TWO-PHOTON EMISSION**

Expanding the time evolution operator to second order, allowing for two-photon emission we have that the *S*-matrix element is

$$S_{i \to f}^{(2)} = -\langle f | \frac{1}{2} \iint_{-\infty}^{\infty} \mathcal{T} \hat{V}(t_2) \hat{V}(t_1) dt_1 dt_2 | i \rangle.$$
(26)

When specifying the final state as  $\langle p_f, k_1, k_2 |$ , an electron, two photons, and the initial state as just an electron,

 $|p_i, 0, 0\rangle$ ,  $S_{fi}^{(2)}$  can be rewritten in terms of the wave functions and the propagator. In [47] this is done for the Compton scattering matrix element, which is the same diagram as here, except that an incoming photon is instead outgoing. The matrix element is therefore

$$S_{i \to f}^{(2)} = -ie^{2} \sqrt{\frac{4\pi}{2\omega_{1}}} \sqrt{\frac{4\pi}{2\omega_{2}}} \iint d^{4}x_{2} d^{4}x_{1} \\ \times \bar{\psi}_{f}^{-}(x_{2}) \phi_{2}^{*} e^{ik_{2}x_{2}} G(x_{2}, x_{1}) \phi_{1}^{*} e^{ik_{1}x_{1}} \psi_{i}^{-}(x_{1}) \\ + (\epsilon_{1}, k_{1}) \leftrightarrow (\epsilon_{2}, k_{2}).$$
(27)

Now we define

$$\begin{split} M_{i\to v}^{-} &= e \sqrt{\frac{4\pi}{2\omega_{1}}} \int d^{3}x \bar{\psi}_{v}^{-}(x) \phi_{1}^{*} e^{-ik_{1}\cdot x} \psi_{i}^{-}(x) \\ &= (2\pi)^{3} \delta(p_{x,i} - p_{x,v} - k_{x}) \delta(p_{z,i} - k_{z,1} - p_{z,v}) \\ &\times \delta(k_{B,i} - k_{y,1} - k_{B,v} - n_{B,1}k_{0}) e \sqrt{\frac{4\pi}{2\omega_{1}}} \frac{1}{2\sqrt{\varepsilon_{i}\varepsilon_{v}}} \\ &\times \sum_{j} c_{n_{B}+j,v}^{*} c_{j,i} \bar{S}_{n_{B}+j,v} \phi_{1}^{*} S_{j,i}^{-} \\ &= (2\pi)^{3} \delta(p_{x,i} - p_{x,v} - k_{x}) \delta(p_{z,i} - k_{z,1} - p_{z,v}) \\ &\times \delta(k_{B,i} - k_{y,1} - k_{B,v} - n_{B,1}k_{0}) \times \mathcal{M}_{i\to v}^{-}(k_{1},\epsilon_{1}), \end{split}$$

$$(28)$$

where  $\mathcal{M}$  is defined as in Eq. (23) where v is used to denote the virtual state from the propagator and is shorthand for the dependence on  $p_{x,v}$ ,  $k_{B,v}$ ,  $p_{z,v}$ ,  $n_v$ , and  $s_v$ . The superscript – on  $M^-$  and  $\mathcal{M}^-$  denotes that the virtual state is the electron state  $\psi_v^-$ , and  $M^+$ ,  $\mathcal{M}^+$  is the same but with the positron virtual state. The matrix element may then be written as

$$S_{fi}^{(2)} = i \iint dt_2 dt_1 \sum_{n_v, s_v} \int \frac{d^3 p_v}{(2\pi)^3} - \theta(t_2 - t_1) e^{i(\omega_1 + \varepsilon_v - \varepsilon_i)t_1} e^{i(\omega_2 + \varepsilon_f - \varepsilon_v)t_2} M_{i \to v}^- M_{v \to f}^- + \theta(t_1 - t_2) e^{i(\omega_1 - \varepsilon_i - \varepsilon_v)t_1} e^{i(\omega_2 + \varepsilon_f + \varepsilon_v)t_2} M_{i \to v}^+ M_{v \to f}^+.$$
(29)

Therefore the term in the second line is seen as the electron first emits a photon with momentum  $k_1$  at  $t_1$  and then propagates to a later time  $t_2$  and emits a second photon with momentum  $k_2$ . The term in the third line is then the electron emitting the photon with momentum  $k_1$  at a time  $t_1$  turning the electron into a positron going into the past and emitting the photon with momentum  $k_2$  at the earlier time  $t_2$ . This last term is heavily suppressed in our case, which we can see as follows. Denote  $a = \varepsilon_v + \omega_1 - \varepsilon_i$  and  $b = \varepsilon_f + \omega_2 - \varepsilon_v$ , then we may use that

$$\theta(t_1 - t_2) = \frac{i}{2\pi} \int_{-\infty}^{\infty} \frac{1}{\varepsilon_V + i\epsilon} e^{i(t_2 - t_1)\varepsilon_V} d\varepsilon_V, \quad (30)$$

where  $\epsilon$  is a small real number for which one in the end should take the limit  $\epsilon \to 0$  and therefore we have

$$\iint_{-\infty}^{\infty} dt_2 dt_1 \theta(t_1 - t_2) e^{iat_2} e^{ibt_1}$$
  
=  $2\pi i \delta(a + b) \frac{1}{-a + i\epsilon}$   
=  $2\pi i \delta(\epsilon_f + \omega_1 + \omega_2 - \epsilon_i) \frac{1}{\epsilon_i - \epsilon_v - \omega_1 + i\epsilon}.$  (31)

We have also that  $-\theta(t_2 - t_1) = \frac{i}{2\pi} \int \frac{1}{\varepsilon_V - i\varepsilon} e^{i(t_2 - t_1)\varepsilon_V} d\varepsilon_V$ , and therefore we have the term from the third line of Eq. (29) that carries the factor of

$$2\pi i\delta(\varepsilon_f + \omega_1 + \omega_2 - \varepsilon_i)\frac{1}{\varepsilon_i + \varepsilon_v - \omega_1 - i\varepsilon}.$$
 (32)

Therefore this term will always be very far off-shell, as the virtual particle on-shell condition can never be met as it corresponds to the spontaneous production of an electron, positron, and photon from the crystal field, where the produced positron is subsequently annihilated with the incoming electron to emit another photon. Having carried out the integrations over time we obtain that

$$S_{fi}^{(2)} = -\sum_{n_v, s_v} \int \frac{d^3 p_v}{(2\pi)^3} 2\pi \delta(\varepsilon_f + \omega_1 + \omega_2 - \varepsilon_i) \\ \times \left( \frac{M_{i \to v}^- M_{v \to f}^-}{\varepsilon_i - \varepsilon_{v_1} - \omega_1 - i\varepsilon} + \frac{M_{i \to v}^+ M_{v \to f}^+}{\varepsilon_i + \varepsilon_{v_1} - \omega_1 + i\varepsilon} + (\varepsilon_1, k_1) \leftrightarrow (\varepsilon_2, k_2) \right).$$
(33)

Now we may integrate over  $p_v$  to obtain

$$S_{fi}^{(2)} = -\sum_{n_v, s_v} \left( \frac{\mathcal{M}_{i \to v_1}^- \mathcal{M}_{v_1 \to f}^-}{\varepsilon_i - \varepsilon_{v_1} - \omega_1 - i\epsilon} + \frac{\mathcal{M}_{i \to v_1}^+ \mathcal{M}_{v_1 \to f}^+}{\varepsilon_i + \varepsilon_{v_1} - \omega_1 + i\epsilon} \right) \\ \times (2\pi)^4 \delta(\varepsilon_f + \omega_1 + \omega_2 - \varepsilon_i) \\ \times \delta(p_{x,i} - k_{x,1} - k_{x,2} - p_{x,f}) \\ \times \delta(p_{z,i} - k_{z,1} - k_{z,2} - p_{z,f}) \\ \times \delta(k_{B,i} - k_{y,1} - k_{y,2} - k_{B,f} - (n_{B,1} + n_{B,2})k_0) \\ + (\varepsilon_1, k_1) \leftrightarrow (\varepsilon_2, k_2),$$
(34)

and then  $v_1$  denotes the virtual state with momentum given by  $p_{x,v_1} = p_{x,i} - k_{x,1}$ ,  $p_{z,v_1} = p_{z,i} - k_{z,1}$  and  $k_{B,v_1} = k_{B,i}^- - k_{y,1} - n_{B,1}^- k_0$  and  $-k_{B,v_1}^+ = k_{B,i}^- - k_{y,1} - n_{B,1}^+ k_0$ , i.e., that photon with label 1 is emitted at the vertex connected with the initial particle. From the amplitude we get the transition probability according to

$$dP^{(2)} = \frac{1}{2} \int \sum_{n_f, s_f} |S_{fi}^{(2)}|^2 \frac{dp_{x,f} dk_{B,f} dp_{z,f}}{(2\pi)^3} \frac{d^3 k_1}{(2\pi)^3} \frac{d^3 k_2}{(2\pi)^3}$$
$$= \sum_{n_f, s_f} \left| \sum_{n_v, s_v} \frac{\mathcal{M}_{i \to v_1}^- \mathcal{M}_{v_1 \to f}^-}{\varepsilon_i - \varepsilon_{v_1} - \omega_1 - i\varepsilon} + \frac{\mathcal{M}_{i \to v_1}^+ \mathcal{M}_{v_1 \to f}^+}{\varepsilon_i + \varepsilon_{v_1} - \omega_1 + i\varepsilon} + (\varepsilon_1, k_1) \leftrightarrow (\varepsilon_2, k_2) \right|^2$$
$$\times \frac{T}{(2\pi)^5} \delta(\varepsilon_f + \omega_1 + \omega_2 - \varepsilon_i) d^3 k_1 d^3 k_2, \qquad (35)$$

where we have added a factor of 1/2 in front due to identical particles in the final state that we in the end want to integrate over all angles and would therefore, again, be counting double [48]. From this full result, it is seen that the result can diverge when  $\epsilon \to 0$  because  $\varepsilon_i - \varepsilon_v - \omega_1 = 0$  is possible. The nature of the divergence is, however, different for some of the terms, namely the ones that are the norm square of each term underneath the sum,  $|\mathcal{M}_{i\to v_1}^- \mathcal{M}_{v_1\to f}^-/(\varepsilon_i - \varepsilon_{v_1} - \omega_1 - i\varepsilon)|^2$ , where the limit of  $\epsilon \to 0$  will yield an infinite result, even after integration over one of the angles  $\theta_1$  or  $\theta_2$ . On the other hand, while the remaining terms, of the interference type, still diverge, they can be integrated over  $\theta_1$  or  $\theta_2$  to yield a convergent result. To learn the meaning of this divergence due to the denominator (see also [49]), we may write

$$\left|\frac{1}{b-i\epsilon}\right|^2 = \frac{1}{b^2 + \epsilon^2},\tag{36}$$

and note that

$$\lim_{\epsilon \to 0} \frac{\epsilon}{b^2 + \epsilon^2} = \pi \delta(b). \tag{37}$$

If we evaluate the integrals of  $\mathcal{M}_{i \to v_1}^- \mathcal{M}_{v_1 \to f}^-$  with the factor  $\delta(a+b)\delta(b)$ , we get well defined results, as this just amounts to the product of two first order emissions. It is therefore useful to write

$$\left|\frac{1}{b-i\epsilon}\right|^2 = \frac{1}{\epsilon} \frac{\epsilon}{b^2 + \epsilon^2},\tag{38}$$

where then the factor  $\epsilon/(b^2 + \epsilon^2)$  acts as a delta function for small enough  $\epsilon$ , yielding a finite value when we perform the integrals in Eq. (35), and then it is clear that this is divergent as  $\epsilon \to 0$  due to the factor of  $1/\epsilon$ . However, this should be understood in terms of an additional factor of Tfor this term. To see this, consider the origin of this expression from Eq. (31), but consider instead that we had a finite time, and integrate over a and b,

$$\int \left| \int_{0}^{T} \theta(t_{1} - t_{2}) e^{iat_{2}} e^{ibt_{1}} dt_{1} dt_{2} \right|^{2} dadb$$
$$= (2\pi)^{2} \int_{0}^{T} \theta(t_{1} - t_{2}) dt_{1} dt_{2} = (2\pi)^{2} \frac{T^{2}}{2}, \quad (39)$$

and we also have that

$$\int \left| 2\pi i \delta(a+b) \frac{1}{b-i\epsilon} \right|^2 dadb = 2\pi T \frac{\pi}{\epsilon}.$$
 (40)

So we see that we must replace  $|\frac{1}{b-i\epsilon}|^2 \rightarrow \pi T \delta(b)$ , and therefore these terms turn out to give us the cascade contribution. To see how the probability from Eq. (35) splits up into this cascade along with additional terms, we will denote the quantity underneath the norm square as  $R^- = \sum_{n_v} J_1^-(n_v) + J_2^-(n_v)$  corresponding to the terms with the virtual electron and similarly  $R^+ = \sum_{n_v} J_1^+(n_v) + J_2^+(n_v)$ , where

$$J_1^-(n_v) = \sum_{s_v} \frac{\mathcal{M}_{i \to v_1}^- \mathcal{M}_{v_1 \to f}^-}{\varepsilon_i - \varepsilon_{v_1} - \omega_1 - i\varepsilon},$$
(41)

$$J_1^+(n_v) = \sum_{s_v} \frac{\mathcal{M}_{i \to v_1}^+ \mathcal{M}_{v_1 \to f}^+}{\varepsilon_i + \varepsilon_{v_1} - \omega_1 + i\epsilon},$$
(42)

and  $J_2$  is  $J_1$  with  $(\epsilon_1, k_1) \leftrightarrow (\epsilon_2, k_2)$ ; we then define  $R = R^+ + R^-$ . The quantity we want is then  $|R|^2 = |R^+|^2 + |R^-|^2 + 2\text{Re}[R^+(R^-)^*]$ . In the  $R^+$  term, it is never possible for the denominator to become 0, and therefore it can be directly calculated (see Appendix B). For

$$|R^{-}|^{2} = \left(\sum_{n_{v}} J_{1}^{-}(n_{v}) + J_{2}^{-}(n_{v})\right) \times \left(\sum_{n_{v}'} J_{1}^{-}(n_{v}') + J_{2}^{-}(n_{v}')\right)^{*}, \quad (43)$$

the product of the terms with the same subscript and where  $n_v = n'_v$  are the cascade process that are the only problematic terms and so need special attention as described above. Therefore it is useful to employ that

$$|R^{-}|^{2} = \sum_{n_{v}} [J_{1}^{-}(n_{v})(R^{-} - J_{1}^{-}(n_{v}))^{*} + J_{2}^{-}(n_{v})(R^{-} - J_{2}^{-}(n_{v}))^{*} + |J_{1}^{-}(n_{v})|^{2} + |J_{2}^{-}(n_{v})|^{2}], \qquad (44)$$

and so the terms in the first two lines are convergent contributions to the one-step process and the terms in the last line are the cascade terms, except that the spin sum is still underneath the norm square. In Appendix C we show that the interference due to spin will be 0 when the photon polarization can be taken as real and that either the sum over initial or final spins (we will do both) is carried out. And so we can write the differential probability of emission, with a given initial state, as

$$dP^{(2)} = \frac{1}{2} \frac{T}{(2\pi)^5} \delta(\varepsilon_f + \omega_1 + \omega_2 - \varepsilon_i) d^3 k_1 d^3 k_2$$
  

$$\times \sum_{n_f} \left( |R^+|^2 + 2\operatorname{Re}[R^+(R^-)^*] + \sum_{n_v} \left\{ \sum_{s_v} [|\mathcal{M}^-_{i \to v_1} \mathcal{M}^-_{v_1 \to f}|^2 \pi T \delta(b)] + J^-_1(n_v) (R^- - J^-_1(n_v))^* + (\varepsilon_1, k_1) \leftrightarrow (\varepsilon_2, k_2) \right\} \right).$$
(45)

For the sake of clarity, we can write this as

$$dP^{(2)} = dP^{(\text{cascade})} + dP^{(1-\text{step})}, \tag{46}$$

where the cascade contribution proportional to  $T^2$  is given by

$$dP^{(\text{cascade})} = \frac{1}{2} \frac{T^2}{(2\pi)^5} \delta(\epsilon_f + \omega_1 + \omega_2 - \epsilon_i) d^3 k_1 d^3 k_2$$
$$\times \sum_{n_f} \sum_{n_v} \left\{ \sum_{s_v} [|\mathcal{M}_{i \to v_1}^- \mathcal{M}_{v_1 \to f}^-|^2 \pi \delta(b)] + (\epsilon_1, k_1) \leftrightarrow (\epsilon_2, k_2) \right\},$$
(47)

which is equivalent to Eq. (25) and where the one-step contribution, proportional to *T*, is given by

$$dP^{(1-\text{step})} = \frac{1}{2} \frac{T}{(2\pi)^5} \delta(\epsilon_f + \omega_1 + \omega_2 - \epsilon_i) d^3 k_1 d^3 k_2$$
  

$$\times \sum_{n_f} \left( |R^+|^2 + 2\text{Re}[R^+(R^-)^*] + \sum_{n_v} \{J_1^-(n_v)(R^- - J_1^-(n_v))^* + (\epsilon_1, k_1) \leftrightarrow (\epsilon_2, k_2)\} \right).$$
(48)

### **V. CHOICE OF REGULARIZATION**

Above we chose a certain way to regularize the divergence, by recognizing that the divergent terms correspond to the cascade terms. In taking the time limit from  $\pm \infty$ , some information about the duration of interaction was lost, which we put back in, in a way that is correct when *T* is large enough, i.e., larger than the photon formation length roughly estimated by  $l_f = 2\gamma^2(1 - \omega/\varepsilon)/\omega$  [25], which in our case is

roughly  $\gamma/m \sim 0.8 \ \mu\text{m}$ , because  $\omega$  is on the order of  $\varepsilon$ . Another way often found in literature [50–54] is to say that the virtual state is unstable and therefore replace the energy of the virtual particle according to  $\varepsilon_v \rightarrow \varepsilon_v - i\Gamma_v/2$  where  $\Gamma_v = \sum_f W_{v \rightarrow f}$  is the total decay width of the virtual state from all processes. This is equivalent to adding the effect of the linewidth in atomic Raman scattering [55]. Effectively this corresponds to replacing the  $\varepsilon$  in the denominator with  $-\Gamma_v/2$ , which lifts the divergence. However, one can see that with this substitution [see Eq. (38)], one would obtain that

$$\left|\frac{1}{b+i\Gamma_v/2}\right|^2 = \frac{2\pi}{\Gamma_v}f(b),\tag{49}$$

where f(b) is a function peaked around b = 0 that obeys  $\int f(b)db = 1$  and therefore resembles the delta function  $\delta(b)$  but with a nonzero width  $\Gamma_v$ . If we then again calculate the cascade part according to this, we would obtain

$$dP^{(\text{cascade})*} = \frac{1}{2} \frac{1}{(2\pi)^5} T d^3 k_1 d^3 k_2$$
  
 
$$\times \sum_{n_f} \sum_{n_v, s_v} |\mathcal{M}_{i \to v_1}^- \mathcal{M}_{v_1 \to f}^-|^2 \delta(a+b) \frac{2\pi}{\Gamma_{v_1}} f(b)$$
  
 
$$+ (\epsilon_1, k_1) \leftrightarrow (\epsilon_2, k_2).$$
(50)

If we assume that the dominant contribution to the decay width is due to radiation emission, we have that the total width is

$$\Gamma_{v} = \sum_{f} \int \frac{1}{(2\pi)^{2}} |\mathcal{M}_{v \to f}|^{2} \delta(\varepsilon_{f} + \omega - \varepsilon_{v}) d^{3}k.$$
 (51)

Therefore if we approximate  $f(b) \simeq \delta(b)$ , integrate over  $d^3k_2$ , and sum over  $n_f$ , we will obtain a factor of the total rate  $\Gamma_{v_1}$ , which cancels out, and so we have that

$$dP^{(\text{cascade})*} = \frac{1}{2} \frac{1}{(2\pi)^2} T d^3 k_1 \sum_{n_v, s_v} |\mathcal{M}_{i \to v_1}^-|^2 \delta(a)$$
  
+  $(\epsilon_1, k_1) \leftrightarrow (\epsilon_2, k_2),$  (52)

which is just the single photon emission probability. Therefore this approach leads to the prediction that it is just as likely to emit two photons as it is one. This is not a meaningful result and the reason is that the integration over time has been carried out over all times; i.e., it is assumed that  $T \gg 1/\Gamma_v$ , which means it is guaranteed that the virtual state decays. However, in that case not only two-photon emission is likely but also a larger number of photons, which we do not take into account. For Raman scattering the approach is reasonable when  $T \gg 1/\Gamma_v$  such that it is guaranteed that an excited state will decay before the observation is made. However, if the interaction time is very short,  $T \ll 1/\Gamma_v$ , it is also expected that Raman



FIG. 2. The differential emission probability of two photons with energy  $\omega_1$  and  $\omega_2$  integrated over angles for the cascade contribution of Eq. (47) divided by  $T^2$  and multiplied by  $\omega_1 \omega_2$  for the case mentioned in the text.

scattering should have a dependence as  $T^2$ , as each subprocess, excitation, and decay is characterized by a rate, and the probability is therefore the product of  $(W^{\text{excite}}T) \times$  $(W^{\text{decay}}T)$ . The substitution  $\varepsilon_v \rightarrow \varepsilon_v - i\Gamma_v/2$  therefore corresponds to the replacement  $W^{\text{decay}}T \rightarrow 1$  and then combines the processes corresponding to the first order diagrams of excitation first, and subsequently decay, with the second order diagram, which allows for off-resonant excitation and decay. We are interested in the case when  $T < 1/\Gamma_v$  such that two-photon emission is unlikely compared to onephoton emission, and therefore a higher number of photon emissions can be neglected. In this case one can also think of the previously obtained result for the cascade contribution, as the contribution of the finite crystal length to the linewidth, which corresponds to setting  $\Gamma_v/2 = 1/T$ , which will be the dominant contribution to the linewidth when  $T \ll 1/\Gamma_v$ .

# **VI. DISCUSSION OF RESULTS**

In the figures in this paper we show the calculations made for a 180 GeV electron in the Doyle-Turner potential [25,36-38] for the (110) planes in silicon and for the state n = 25. This is a quite low lying state which for electrons will have a high radiation power [28]. Electrons were chosen for this reason as it is not as numerically heavy when the quantum numbers are relatively small, as opposed to the positron case, which would require large quantum numbers to obtain an appreciable value of the quantum nonlinearity parameter  $\chi$ , which means that quantum effects such as spin and recoil are important in the emission process. To compare with an experiment one should average over the distribution of the initial states, which depends on the particle beam angular mean and divergence. In Eq. (45) the integrals over  $\varphi$  and  $\theta$ are carried out numerically over the intervals  $0 < \varphi < 2\pi$ and  $0 < \theta < \frac{1.5}{v} \times (1 + \xi)$ , and therefore include nearly all emitted radiation. From the result of Eq. (45) we see that the part scaling with  $T^2$  is the cascade, obtained by simple multiplication of probabilities, and will dominate unless the crystal is very thin, due to the remaining terms being proportional to T. Therefore, if one made a Monte Carlo approach using the single photon emission rate using the quantum numbers of the current state, instead of using the constant field approximation with the current value of the field, one would obtain the dominant (cascade) contribution, which will be accurate also when the constant field approximation is no longer valid. In Fig. 2 we show the result from the cascade process. In Fig. 3 we show the one-step terms,





FIG. 3. The differential emission probability of two photons with energy  $\omega_1$  and  $\omega_2$  integrated over angles for the one-step contribution of Eq. (48) divided by *T* and multiplied by  $\omega_1\omega_2$ , for the case mentioned in the text and as in Fig. 2.

FIG. 4. The ratio of differential emission probabilities of two photons with energy  $\omega_1$  and  $\omega_2$  integrated over angles for the onestep contribution of Eq. (48) to the cascade contribution of Eq. (47) when  $T = 20 \ \mu$ m, for the case mentioned in the text and as in Fig. 2. This ratio therefore scales as 1/T.

and finally in Fig. 4 we show the ratio of these one-step terms to the cascade terms for  $T = 20 \ \mu m$ . From this figure we see that the one-step terms can become significant compared to the cascade terms for short crystals. This ratio scales as 1/T. Therefore one needs a thin crystal for the one-step contribution to be significant, so thin that the probability to emit more than one photon becomes small. One may rightfully ask based on these figures, if one picks a very small value of T, could the total probability seemingly become negative? However, the results shown are valid only when  $T \gg l_f \sim$ 0.8  $\mu$ m as estimated earlier. For the 180 GeV case calculated here, the probability to emit a photon with energy above 1 GeV from a 20  $\mu$ m crystal is roughly 7%, and therefore the probability corresponding to the cascade for two-photon emission above this photon energy is 0.25%, and as can be seen in Fig. 4 the spectrum in the region where the radiation is most abundant, the ratio is around  $\pm 20\%$ . This number serves as an upper limit to the size of the effect, because under experimental conditions one would obtain the average from a population of many different levels with different quantum numbers n, and this averaging would likely reduce the size of the effect. If we assume the size of the effect to be this upper limit, one would need enough events such that one would have enough statistics to see an effect of such a size from only 0.25% of the events. If this setup was realized by adding a calorimeter to a setup as the one used in [18], we can estimate the number of particles required to see this. Making a histogram of 20 bins in each direction of  $\omega_1$  and  $\omega_2$  and assuming 100 counts on average in each bin, one would need roughly  $3.2 \times 10^8$  electrons; assuming an electron rate of  $10^4$  / min, this translates into roughly 22 days of measuring time. This would therefore be a challenging experiment, and having in mind that there would likely also be systematic uncertainties, the realistic outcome of such an experiment would be to put a constraint on the size of such one-step terms, rather than their direct observation.

### VII. CONCLUSION

In conclusion, we have shown how to accurately calculate the two-photon emission rate for a high-energy electron (or positron) channeled in a crystal. This calculation shows that the full probability contains what is known as the cascade, which could have been obtained by multiplying probabilities of single photon emissions, as well as additional interference terms, called the one-step contribution. The one-step contribution scales only linearly with the crystal length, and therefore one needs a thin crystal to see the effect of these terms. We have calculated the size of all contributions to the emission probability for 180 GeV electrons in silicon and found that with a long measuring time, the one-step contribution could possibly be seen. Since these effects are, however, small, we also see how to solve the problem of the quantum radiation reaction under general circumstances, in a crystal, by using the single photon emission rate in consecutive emissions, corresponding to the particle's current state.

## ACKNOWLEDGMENTS

The author gratefully acknowledges useful discussions with Antonino Di Piazza and Karen Z. Hatsagortsyan. This work was partially supported by a research grant (VKR023371) from VILLUM FONDEN and later by the Alexander von Humboldt-Stiftung. In addition, the author acknowledges the support of NVIDIA Corporation with the donation of the Titan V GPU used for this research.

## **APPENDIX A: POSITRON WAVE FUNCTION**

The general (unnormalized) solution to the Dirac equation with potential energy  $V(\mathbf{r}) = -e\varphi(\mathbf{r})$  can be written as

$$\psi(\mathbf{r},t) = e^{-i\varepsilon t} \begin{pmatrix} \phi(\mathbf{r}) \\ \chi(\mathbf{r}) \end{pmatrix}.$$
 (A1)

The Dirac equation then becomes

$$(\varepsilon + e\varphi - m)\phi(\mathbf{r}) = \boldsymbol{\sigma} \cdot \hat{\boldsymbol{p}}\chi(\mathbf{r}),$$
 (A2)

$$(\varepsilon + e\varphi + m)\chi(\mathbf{r}) = \boldsymbol{\sigma} \cdot \hat{\boldsymbol{p}}\phi(\mathbf{r}).$$
 (A3)

The electron solution is then (see Ref. [28])

$$\psi^{-}(\mathbf{r},t) = e^{-i\varepsilon t} \begin{pmatrix} \phi(\mathbf{r}) \\ \frac{\sigma \cdot \hat{p}}{\varepsilon + e\varphi(\mathbf{r}) + m} \phi(\mathbf{r}) \end{pmatrix}.$$
 (A4)

We then obtained an equation for  $\phi(\mathbf{r})$  by isolating  $\chi(\mathbf{r})$  in Eq. (A3) and inserting it in Eq. (A2). This solution has the property that it is well defined when  $\varepsilon = m$ . Another solution can be found by isolating  $\phi(\mathbf{r})$  in Eq. (A2) and inserting it in (A3). However, this solution is not well defined when  $\varepsilon = m$ , and therefore one must use the negative energy solution  $\varepsilon = -\sqrt{p^2 + m^2} = -\varepsilon_p$ . Therefore we have

$$\psi(\mathbf{r},t) = e^{i\varepsilon_p t} \begin{pmatrix} -\frac{\sigma \cdot \hat{\mathbf{p}}}{\varepsilon_p - e\varphi(\mathbf{r}) + m} \chi(\mathbf{r}) \\ \chi(\mathbf{r}) \end{pmatrix}, \quad (A5)$$

where  $\varepsilon_p$  is the positive energy of the positron. The equation for  $\chi$  can now be obtained by using

$$(\varepsilon + e\varphi + m)\chi(\mathbf{r}) = \boldsymbol{\sigma} \cdot \hat{\boldsymbol{p}} \frac{1}{\varepsilon + e\varphi - m} \boldsymbol{\sigma} \cdot \hat{\boldsymbol{p}}\chi(\mathbf{r}),$$
 (A6)

which is equivalent to

$$(\varepsilon_p - e\varphi - m)\chi(\mathbf{r}) = \boldsymbol{\sigma} \cdot \hat{\boldsymbol{p}} \frac{1}{\varepsilon_p - e\varphi + m} \boldsymbol{\sigma} \cdot \hat{\boldsymbol{p}}\chi(\mathbf{r}).$$
 (A7)

This is the same equation as the one we obtained for  $\phi(\mathbf{r})$ , except with the sign of *e* changed such that, after making the same approximations as we did in [28,29],

$$[\hat{p}^2 + 2\varepsilon_p e\varphi(\mathbf{r}) - (\varepsilon_p^2 - m^2)]\chi(\mathbf{r}) = 0.$$
 (A8)

We therefore make the ansatz in line with the usual approach (the sign on the momenta is changed):

$$\chi(\mathbf{r}) = \mathbf{s}I^+(\mathbf{y})e^{-i(p_x \mathbf{x} + p_z z)},\tag{A9}$$

$$I^{+}(y) = e^{-ik_{B}y} \sum_{j} c_{j} e^{-ijk_{0}y}.$$
 (A10)

Then

$$\psi(\mathbf{r},t) = e^{i(-p_x x - p_z z + \varepsilon_p t)} \begin{pmatrix} \frac{\sigma \cdot p}{\varepsilon_p + m} \mathbf{s} I^+(y) \\ \mathbf{s} I^+(y) \end{pmatrix}, \qquad (A11)$$

with  $\boldsymbol{p} = (p_x + q\varphi(\boldsymbol{r}), i\frac{d}{dy}, p_z)$ . Inserting  $I^+(y)$ , this becomes

$$\psi(\mathbf{r},t) = e^{i(-p_x x - p_z z + \varepsilon_p t)} \sum_j c_j e^{-i(k_B + jk_0)y} \begin{pmatrix} \frac{\sigma p}{\varepsilon_p + m} \mathbf{s} \\ \mathbf{s} \end{pmatrix}, \quad (A12)$$

with 
$$\mathbf{p} = (p_x + E_n - \frac{(jk_0 + k_B)^2}{2\varepsilon}, jk_0 + k_B, p_z).$$

# **APPENDIX B: POSITRON MATRIX ELEMENT**

Even though we consider the radiation from electrons, the propagator contains terms from the positron  $\psi_{p,n,s}^+(\mathbf{x}_2)\bar{\psi}_{p,n,s}^+(\mathbf{x}_1)$ . Therefore we will need to calculate

$$\mathcal{M}_{i\to v}^{+} = e \sqrt{\frac{4\pi}{2\omega}} \frac{1}{2\sqrt{\varepsilon_v \varepsilon_i}} \sum_{j,l} c_{l,v}^* c_{j,i} \bar{\mathbf{S}}_{l,v}^+ \boldsymbol{\mathscr{I}}^* \mathbf{S}_{j,i}^- \quad (B1)$$

and  $\mathcal{M}^+_{v \to f}$  so we need

$$\begin{aligned} -\frac{1}{\sqrt{\varepsilon_{i}+m}\sqrt{\varepsilon_{v}+m}}\bar{S}_{v}^{T}q^{s}S_{i}^{-} \\ &= \left[\left(S_{v}^{T}\frac{\sigma p_{v}}{\varepsilon_{v}+m}}S_{v}^{T}\right)\left(\begin{array}{cc}0&\sigma\cdot\epsilon^{*}\\\sigma\cdot\epsilon^{*}&0\end{array}\right)\left(\frac{s_{i}}{\frac{\sigma p_{i}}{\varepsilon_{i}+m}}s_{i}\right)\right] \\ &= \left[\left(S_{v}^{T}\frac{\sigma p_{v}}{\varepsilon_{v}+m}}S_{v}^{T}\right)\left(\begin{array}{cc}\sigma\cdot\epsilon^{*}\frac{\sigma p_{i}}{\varepsilon_{i}+m}s_{i}\\\sigma\cdot\epsilon^{*}s_{i}\end{array}\right)\right] \\ &= s_{v}^{T}\left[\frac{\sigma\cdot p_{v}}{\varepsilon_{v}+m}\sigma\cdot\epsilon^{*}\frac{\sigma\cdot p_{i}}{\varepsilon_{i}+m}+\sigma\cdot\epsilon^{*}\right]s_{i} \\ &= s_{v}^{T}\left[\frac{1}{(\varepsilon_{i}+m)(\varepsilon_{v}+m)}(\sigma\cdot p_{v})(\sigma\cdot\epsilon^{*})(\sigma\cdot p_{i})+\sigma\cdot\epsilon^{*}\right]s_{i} \\ &= s_{v}^{T}\left[\frac{1}{(\varepsilon_{i}+m)(\varepsilon_{v}+m)}(\sigma\cdot p_{v})(\epsilon^{*}\cdot p_{i}+i\sigma\cdot[\epsilon^{*}\times p_{i}])+\sigma\cdot\epsilon^{*}\right]s_{i} \\ &= s_{v}^{T}\left[\frac{1}{(\varepsilon_{i}+m)(\varepsilon_{v}+m)}\left\{(\sigma\cdot p_{v})(\epsilon^{*}\cdot p_{i})+i(\sigma\cdot p_{v})\sigma\cdot[\epsilon^{*}\times p_{i}]\right\}+\sigma\cdot\epsilon^{*}\right]s_{i} \\ &= s_{v}^{T}\left[\frac{1}{(\varepsilon_{i}+m)(\varepsilon_{v}+m)}\left\{(\sigma\cdot p_{v})(\epsilon^{*}\cdot p_{i})+ip_{v}\cdot(\epsilon^{*}\times p_{i})+i\sigma\cdot(p_{v}\times(\epsilon^{*}\times p_{i}))\right]\right\}+\sigma\cdot\epsilon^{*}\right]s_{i} \\ &= s_{v}^{T}\left[\frac{1}{(\varepsilon_{i}+m)(\varepsilon_{v}+m)}\left\{(\sigma\cdot p_{v})(\epsilon^{*}\cdot p_{i})+ip_{v}\cdot(\epsilon^{*}\times p_{i})-\sigma\cdot(p_{v}\times(\epsilon^{*}\times p_{i}))\right\}+\sigma\cdot\epsilon^{*}\right]s_{i} \\ &= s_{v}^{T}\left[\frac{1}{(\varepsilon_{i}+m}(\varepsilon_{v}+m)}\left\{(\sigma\cdot p_{v})(\epsilon^{*}\cdot p_{i})+ip_{v}\cdot(\epsilon^{*}\times p_{i})-\sigma\cdot(p_{v}\times(\epsilon^{*}\times p_{i}))\right\}+\sigma\cdot\epsilon^{*}\right]s_{i} \\ &= s_{v}^{T}\left[\frac{1}{(\varepsilon_{i}+m}(\varepsilon_{v}+m)}\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{(\varepsilon_{v}+m)\left\{$$

Then

$$C = \frac{\boldsymbol{p}_{v} \cdot (\boldsymbol{\epsilon}^{*} \times \boldsymbol{p}_{i})}{(\boldsymbol{\varepsilon}_{i} + m)(\boldsymbol{\varepsilon}_{v} + m)} = \frac{\boldsymbol{\epsilon}^{*} \cdot (\boldsymbol{p}_{i} \times \boldsymbol{p}_{v})}{(\boldsymbol{\varepsilon}_{i} + m)(\boldsymbol{\varepsilon}_{v} + m)},$$
(B3)

$$\boldsymbol{D} = \frac{(\boldsymbol{\epsilon}^* \cdot \boldsymbol{p}_i)\boldsymbol{p}_v - \boldsymbol{p}_v \times (\boldsymbol{\epsilon}^* \times \boldsymbol{p}_i)}{(\boldsymbol{\epsilon}_i + m)(\boldsymbol{\epsilon}_v + m)} + \boldsymbol{\epsilon}^*.$$
(B4)

Here we may use that  $p_v \times (\epsilon^* \times p_i) = \epsilon^* (p_i \cdot p_v) - p_i (p_v \cdot \epsilon^*)$ , and so

$$\boldsymbol{D} = \frac{(\boldsymbol{\epsilon}^* \cdot \boldsymbol{p}_i)\boldsymbol{p}_v + \boldsymbol{p}_i(\boldsymbol{p}_v \cdot \boldsymbol{\epsilon}^*)}{(\boldsymbol{\varepsilon}_i + m)(\boldsymbol{\varepsilon}_v + m)} + \boldsymbol{\epsilon}^* \left(1 - \frac{\boldsymbol{p}_i \cdot \boldsymbol{p}_v}{(\boldsymbol{\varepsilon}_i + m)(\boldsymbol{\varepsilon}_v + m)}\right).$$
(B5)

Now consider the other part for  $\mathcal{M}_{v \to f}^+(\mathbf{k}, \epsilon)$ ,

This is the same as before except with  $i \rightarrow v$  and  $v \rightarrow f$ . And now we want the quantity

$$\sum_{j,l} c_{l,v}^* c_{j,i} \bar{\mathbf{S}}_{l,v}^+ \boldsymbol{\xi}^* \mathbf{S}_{j,i}^{-} e^{i(k_{B,v} + k_{B,i} - k_y)y} e^{i(j+l)k_0 y}$$

$$= 2\pi \delta(k_{B,v} + k_{B,i} - k_y - n_{B,1}^+ k_0) \sum_j c_{-(n_{B,1}^+ + j),v}^* c_{j,i} \bar{\mathbf{S}}_{-(n_{B,1}^+ + j),v}^+ \boldsymbol{\xi}^* \mathbf{S}_{j,i}^{-}, \tag{B7}$$

where now  $n_{B,1}^+$  is chosen such that  $k_{B,v} = k_y - k_{B,i} + n_{B,1}^+ k_0$  is in the first brillouin zone. Note that  $-k_{B,v} = k_{B,i} - k_y - n_{B,1}^+ k_0$  for which we already have the solution, called  $k_{B,v}^- = k_{B,i} - k_y - n_{B,1}^- k_0$ , and therefore

$$k_{B,v} = -k_{\overline{B},v} + k_0 = -k_{B,i} + k_y + (n_{\overline{B},1} + 1)k_0; \quad (B8)$$

therefore  $n_{B,1}^+ = n_{B,1}^- + 1$ . For the  $\mathcal{M}_{v \to f}^+$  term one obtains that  $k_{B,f} = -k_{B,v} - k_y - n_{B,2}^+ k_0$ , and for this term one has that  $n_{B,2}^+ = n_{B,2}^- - 1$ , in terms of the  $n_{B,2}^-$  value for the corresponding electron term in the propagator. The *l* index is given by  $l = n_{B,2}^+ - j$ .

### **APPENDIX C: SPIN INTERFERENCE**

We need to consider  $|\sum_{s_v} \mathcal{M}_2 \mathcal{M}_1|^2$ ; in particular we would like to show that  $\operatorname{Re}([\mathcal{M}_{2,\uparrow} \mathcal{M}_{1,\uparrow}][\mathcal{M}_{2,\downarrow} \mathcal{M}_{1,\downarrow}]^{\dagger})$  is 0, where the arrows denote the spin state of the virtual particle. This we may rearrange and consider therefore the product  $\mathcal{M}_{2,\downarrow}^{\dagger} \mathcal{M}_{2,\uparrow}$ . Now we may use that  $\mathcal{M}$  can be written as

$$\mathcal{M}_{2} = e \sqrt{\frac{4\pi}{2\omega}} \frac{1}{2\sqrt{\varepsilon_{f}\varepsilon_{i}}} \sum_{j} c^{*}_{n_{B,2}+j,f} c_{j,v} \bar{\mathbf{S}}_{n_{B,2}+j,f} \boldsymbol{\mathscr{E}}^{*} \mathbf{S}_{j,v}$$
$$= -e \sqrt{\frac{4\pi}{2\omega}} \frac{1}{2\sqrt{\varepsilon_{f}\varepsilon_{i}}} \sum_{j} c^{*}_{n_{B,2}+j,f} c_{j,v} \mathbf{s}^{\dagger}_{f}$$
$$\times [\boldsymbol{\varepsilon}^{*} \cdot \boldsymbol{A}_{n_{B,2}+j,j} + i\boldsymbol{B}_{n_{B,2}+j,j} \cdot \boldsymbol{\sigma}] \boldsymbol{s}_{v}. \tag{C1}$$

Now for simplicity we define

$$\tilde{A} = -e\sqrt{\frac{4\pi}{2\omega}} \frac{1}{2\sqrt{\varepsilon_f \varepsilon_i}} \sum_j c^*_{n_{B,2}+j,f} c_{j,v} A_{n_{B,2}+j,j}, \quad (C2)$$

$$\tilde{\boldsymbol{B}} = -e\sqrt{\frac{4\pi}{2\omega}} \frac{1}{2\sqrt{\varepsilon_f \varepsilon_i}} \sum_j c^*_{n_{B,2}+j,f} c_{j,v} \boldsymbol{B}_{n_{B,2}+j,j}, \quad (C3)$$

and then we have that

$$\mathcal{M}_2 = \mathbf{s}_f^{\dagger} [\boldsymbol{\epsilon}^* \cdot \tilde{\boldsymbol{A}} + i \tilde{\boldsymbol{B}} \cdot \boldsymbol{\sigma}] s_v.$$
(C4)

Therefore

$$\mathcal{M}_{2,\downarrow}^{\dagger}\mathcal{M}_{2,\uparrow} = \mathbf{s}_{\downarrow}^{\dagger} [\boldsymbol{\epsilon}^{*} \cdot \tilde{\boldsymbol{A}} + i\tilde{\boldsymbol{B}} \cdot \boldsymbol{\sigma}]^{\dagger} \boldsymbol{s}_{f} \boldsymbol{s}_{f}^{\dagger} [\boldsymbol{\epsilon}^{*} \cdot \tilde{\boldsymbol{A}} + i\tilde{\boldsymbol{B}} \cdot \boldsymbol{\sigma}] \boldsymbol{s}_{\uparrow}.$$
(C5)

We assume that  $\epsilon^* = \epsilon$ , which is possible if we choose linear polarization as our basis, and we will perform the summation of final spins. Therefore  $s_f s_f^{\dagger}$  is the identity

$$\mathcal{M}_{2,\downarrow}^{\dagger}\mathcal{M}_{2,\uparrow} = \mathbf{s}_{\downarrow}^{\dagger} [\boldsymbol{\epsilon} \cdot \tilde{\boldsymbol{A}} - i\tilde{\boldsymbol{B}} \cdot \boldsymbol{\sigma}] [\boldsymbol{\epsilon} \cdot \tilde{\boldsymbol{A}} + i\tilde{\boldsymbol{B}} \cdot \boldsymbol{\sigma}] \mathbf{s}_{\uparrow}$$
$$= \mathbf{s}_{\downarrow}^{\dagger} [(\boldsymbol{\epsilon} \cdot \tilde{\boldsymbol{A}})^2 + \tilde{\boldsymbol{B}}^2] \mathbf{s}_{\uparrow}$$
$$= 0, \qquad (C6)$$

where we used that  $\tilde{\boldsymbol{B}}$  is a real vector. For the other term,  $\mathcal{M}_{1,\uparrow}\mathcal{M}_{1,\downarrow}^{\dagger}$ , the same can be done, and here the argument hinges upon summation over initial spins. Therefore, if a summation is carried out over either initial or final spins, the spin interference terms will be 0.

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