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### Enhanced Correlation of Electron-Positron Pair in Two and Three Dimensions \*

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Early time electron-positron correlation in vacuum pair-production in an external field is investigated. The entangled electron and positron wave functions are obtained analytically in the configuration and momentum spaces. It is shown that, relative to that of the one-dimensional theory, two- and three-dimensional calculations yield enhanced spatial correlation and broadened momentum spectra. In fact, at early times the electron and positron almost coincide spatially. The correlation also depends on the direction of the applied field. For the spatial correlation, the transverse correlation is stronger than the longitudinal correlation.

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Pair creation in a vacuum can be considered as a 'vacuum breakdown' by a supercritical field.<sup>[1]</sup> Since Sauter<sup>[2]</sup> and Schwinger<sup>[3]</sup> obtained the pair creation rate in a static field, many theoretical and experimental studies on this topic have been performed.<sup>[4-12]</sup> In particular, it has been shown that the pair creation rate can be improved by tailoring the applied field.<sup>[13-16]</sup> However, the pair birth process itself, such as the space and momentum correlations between the just-created electrons and positrons, is less well studied.<sup>[6,17,18]</sup>

In this Letter, using computational quantum field methods<sup>[19]</sup> we investigate the electron-positron joint probability distributions in the configuration and momentum spaces. As in Ref. [6], we are interested in the very early stage, namely at times  $t \ll 1/c^2$  after the applied potential is turned on. The one-dimensional (1D) spatial density distribution<sup>[6]</sup> is generalized to higher dimensions. The correlated spatial and momentum distributions, parallel and perpendicular (referred to as the longitudinal and transverse correlations, respectively) to the external field are obtained.

In our model, we use the Sauter potential<sup>[2]</sup>  $V(\mathbf{r}) = V_0[1 + \tanh(x/W)]/2$ , where W is the spatial extent of the corresponding electric field. The potential is abruptly turned on at t=0. The evolution of the field operator  $\hat{\Psi}(\mathbf{r},t)$  is given by the Dirac equation<sup>[20]</sup>  $i\partial_t \hat{\Psi}(\mathbf{r},t) = (c\alpha \hat{\mathbf{P}} + \beta c^2 + V) \hat{\Psi}(\mathbf{r},t)$ , where  $\alpha$  and  $\beta$  are the Dirac matrices, c is the vacuum light speed,  $\hat{\mathbf{P}}$  is the momentum operator, and V is the external potential. The field operator can be written in terms of the electron creation and annihilation operators as

$$\hat{\Psi}(\boldsymbol{r},t) = \sum_{\boldsymbol{p}} \hat{b}_{\boldsymbol{p}}(t) W_{\boldsymbol{p}}(\boldsymbol{r}) + \sum_{\boldsymbol{n}} \hat{d}_{\boldsymbol{n}}^{\dagger}(t) W_{\boldsymbol{n}}(\boldsymbol{r})$$
$$= \sum_{\boldsymbol{p}} \hat{b}_{\boldsymbol{p}} W_{\boldsymbol{p}}(\boldsymbol{r},t) + \sum_{\boldsymbol{n}} \hat{d}_{\boldsymbol{n}}^{\dagger} W_{\boldsymbol{n}}(\boldsymbol{r},t), \quad (1)$$

where  $W_{\boldsymbol{p}(\boldsymbol{n})}(\boldsymbol{r})$  is the energy eigenfunction of the field-free Dirac equation, and  $W_{\boldsymbol{p}(\boldsymbol{n})}(\boldsymbol{r},t) = \langle \boldsymbol{r}|U(t)|\boldsymbol{p}(\boldsymbol{n})\rangle$  is the solution of the time-dependent Dirac equation with the time evolution operator  $U(t) = \exp[-i(c\boldsymbol{\alpha}\hat{\boldsymbol{P}} + \beta c^2 + V)t]$ . The electron-positron wave function is given by the positive-frequency parts of the field operator and its charge-conjugated field operator  $\phi(\boldsymbol{r}_1, \boldsymbol{r}_2, t) =$  $\langle 0|\hat{\Psi}^{(+)}(\boldsymbol{r}_1, t) \otimes \hat{\Psi}_c^{(+)}(\boldsymbol{r}_2, t)|0\rangle$ , where  $\boldsymbol{r}_1$  and  $\boldsymbol{r}_2$  are the spatial coordinates of the created electron and positron, respectively. Unless otherwise stated, in this work atomic units ( $e = m_e = \hbar = 1$ ) are used. Based on the eigenstates of the free Dirac operator, we can express the pair wave function as<sup>[6]</sup>

$$\phi(\boldsymbol{r}_1, \boldsymbol{r}_2, t) = \sum_{\boldsymbol{n}} \sum_{\boldsymbol{p}} A_{\boldsymbol{p}\boldsymbol{n}}(t) W_{\boldsymbol{p}}(\boldsymbol{r}_1) \otimes CW_{\boldsymbol{n}}^*(\boldsymbol{r}_2), \quad (2)$$

where the matrix C is the charge-conjugated operator and  $A_{pn}(t) = \sum_{\mathbf{P}} \langle \mathbf{p} | U(t) | \mathbf{P} \rangle \langle \mathbf{n} | U(t) | \mathbf{P} \rangle^*$  is the expansion coefficient.

For  $t \ll 1/c^2$ , up to  $O(t^2)$  accuracy one obtains  $A_{pn} = i \langle p | V | n \rangle t = i V_{pn} t$ . Thus, the wave function  $\phi(\mathbf{r}_1, \mathbf{r}_2, t) = \phi_0(\mathbf{r}_1, \mathbf{r}_2) t$  grows linearly with

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time. One can then explore the pair creation process in this early-time regime by following the reduced time-independent wave function  $\phi_0(\mathbf{r}_1, \mathbf{r}_2) = i \sum_{\mathbf{n}} \sum_{\mathbf{p}} V_{\mathbf{p}n} W_{\mathbf{p}}(\mathbf{r}_1) \otimes CW^*_{\mathbf{n}}(\mathbf{r}_2)$ . To investigate the spatial entanglement between the electron and positron, we consider the spatial joint probability distribution  $\rho(\mathbf{r}_1, \mathbf{r}_2) = |\phi_0(\mathbf{r}_1, \mathbf{r}_2)|^2$  for the electron at  $\mathbf{r}_1$  and the positron at  $\mathbf{r}_2$ . The probability of finding the pair is given by<sup>[21]</sup>  $P(t) = t^2 \int d\mathbf{r}_1 d\mathbf{r}_2 \rho(\mathbf{r}_1, \mathbf{r}_2) = t^2 \int d\mathbf{p} d\mathbf{n} |V_{\mathbf{p}n}|^2$ . The corresponding spectrum of the spatial joint probability distribution (i.e., the momentum joint probability distribution) is  $\rho(\mathbf{p}, \mathbf{n}) = |V_{\mathbf{p}n}|^2$ , which measures the electron with momentum  $\mathbf{p}$  and the positron with momentum  $\mathbf{n}$ .

According to the Dirac theory, the wave functions of the free eigenstates are

$$W_{\boldsymbol{p}}(\boldsymbol{r}) = \frac{1}{(2\pi)^{3/2}} \sqrt{\frac{c^2}{E_{\boldsymbol{p}}}} \mu_{\boldsymbol{p}} \exp(i\boldsymbol{p}r), \qquad (3)$$

with the eigenvalue  $E_{\boldsymbol{p}}=\sqrt{c^2p_x^2+c^2p_y^2+c^2p_z^2+c^4}$  and

$$W_{\boldsymbol{n}}(\boldsymbol{r}) = \frac{1}{(2\pi)^{3/2}} \sqrt{\frac{c^2}{E_{\boldsymbol{n}}}} \nu_{\boldsymbol{n}} \exp(-i\boldsymbol{n}\boldsymbol{r}), \qquad (4)$$

with the eigenvalue  $-E_{\mathbf{n}} = -\sqrt{c^2 n_x^2 + c^2 n_y^2 + c^2 n_z^2 + c^4}$ , where  $\mu_p$  and  $\nu_n$  are the Dirac 4D spinors of the electron and positron, respectively. Because the external potential V depends only on x, the simplified expansion coefficient  $V_{pn} = \langle p | V(x) | n \rangle$  should contain the factors  $\delta(p_y + n_y)$  and  $\delta(p_z + n_z)$ , which serve to ensure momentum conservation in the transverse directions. For simplicity, for the 2D case we can choose the spinors as<sup>[22]</sup>  $\mu_p = \sqrt{(E_p + c^2)/2c^2} (1, 0, 0, cp_+/(E_p + c^2)/2c^2)$  $(c^2)$ ) and  $\nu_n = \sqrt{(E_n + c^2)/2c^2}(cn_-/(E_n + c^2), 0, 0, 1),$ where  $p_{+} = p_{x} + ip_{y}$  and  $n_{-} = n_{x} - in_{y}$ . There is no need to take all of the spinors into account because the nonzero 8 components of  $4 \times 4 = 16$  entangled wave function can be divided into two sets of 4-component form and they are conjugate with each other in the 2D system. Thus, without losing validity, we need only one set of 4 components; for example,  $\phi_0^i(\xi_x, \xi_y)$ , where i=1,2,3,4. In the limit of  $W = \infty$ , the potential approximates to  $V(x,y) = V_0(1+x/W)/2$ . The 2D reduced wave function  $\phi_0(\xi_x,\xi_y)$  (here and throughout the paper we omit i, which does not cause any confusion) is then

$$\phi_{0}(\xi_{x},\xi_{y}) = \frac{V_{0}c}{32W\pi^{2}} \int dP_{x}dP_{y} \frac{E_{P_{x}P_{y}} + 1 + P_{y}^{2} + iP_{x}P_{y}}{E_{P_{x}P_{y}}^{3}} \\ \times \begin{pmatrix} 1 \\ -\frac{P_{x} + iP_{y}}{1 + E_{P_{x}P_{y}}} \\ \frac{P_{x} + iP_{y}}{1 + E_{P_{x}P_{y}}} \\ -\left(\frac{P_{x} + iP_{y}}{1 + E_{P_{x}P_{y}}}\right)^{2} \end{pmatrix} \exp(iP_{x}\xi_{x})\exp(iP_{y}\xi_{y}), \quad (5)$$

where  $E_{P_xP_y} = \sqrt{1 + P_x^2 + P_y^2}$ ,  $\xi_x = c(x_1 - x_2)$ ,  $\xi_y = c(y_1 - y_2)$ , and  $P_x = (p_x - n_x)/2c$  and  $P_y = p_y/c$  are the relative momenta in the x and y directions measured by c. Clearly, except for a constant factor the momentum-space wave function  $\phi_0(P_x, P_y)$  can be regarded as the Fourier transformation of  $\phi_0(\xi_x, \xi_y)$ , or

$$\phi_{0}(P_{x}, P_{y}) = \frac{E_{P_{x}P_{y}} + 1 + P_{y}^{2} + iP_{x}P_{y}}{E_{P_{x}P_{y}}^{3}} \cdot \left( \begin{array}{c} 1 \\ -\frac{P_{x} + iP_{y}}{1 + E_{P_{x}P_{y}}} \\ \frac{P_{x} + iP_{y}}{1 + E_{P_{x}P_{y}}} \\ -\left(\frac{P_{x} + iP_{y}}{1 + E_{P_{x}P_{y}}}\right)^{2} \end{array} \right), \quad (6)$$

so that the momentum spectrum is

$$\rho(P_x, P_y) = |\phi_0(P_x, P_y)|^2 = 4 \frac{1 + P_y^2}{(1 + P_x^2 + P_y^2)^2}.$$
 (7)



**Fig. 1.** (Color online) (a) The spatial density distribution  $\rho(\xi_x)$  for the 1D case. (b) The corresponding momentum spectrum  $\rho(P_x)$ . The circles represent the results for  $W = \infty$ . The data are scaled to match at  $\xi_x = 0$  and  $P_x = 0$  for better graphical clarity.

To see more detail of the relation between the joint distributions in configuration and momentum spaces, we first consider a 1D system<sup>[6]</sup> by setting  $p_y = n_y = 0$ . Without loss of generality, we can assume that the positron is always detected at x = 0. In Fig. 1, we show the joint distribution in the presence of fields of different widths. As one can see, if the spatial extent of the field is wider than  $3\lambda_e$ , where  $\lambda_e$  is the electron Compton wavelength, then the joint distributions agree with the analytical result for an infinite-width field, shown as circles in Fig. 1. As the field width becomes narrower, the average distance between the particles shrinks until it vanishes.<sup>[6]</sup> As

is expected, the corresponding momentum spectrum broadens until the width becomes infinity.

Figure 2 shows the electron joint distribution obtained from the analytical wave functions Eqs. (5) and (6) for a quasi-2D system<sup>[23]</sup> with finite y momenta. We can see that, as  $p_y$  increases, the spatial density distribution in the *x*-direction becomes narrower, accompanied by broadening of the corresponding momentum spectrum, as expected since the spatial distribution width  $\Delta x$  and the momentum spectrum width  $\Delta P_x$  should satisfy the uncertainly principle,  $\Delta x \Delta P_x \sim 1$ . It can also be verified that in the 1D limit ( $p_y = 0$ ), both the spatial and momentum distributions agree with that shown in Fig. 1.



Fig. 2. (Color online) The electron-positron spatial density distribution  $\rho(\xi_x)$  in (a) and the corresponding momentum spectrum  $\rho(P_x)$  in (b) in the quasi-2D case for different given  $P_y$ . The analytical wave function Eqs. (5) and (6) are used. The data are scaled to match at  $\xi_x = 0$ and  $P_x = 0$ .

Now we consider a full 2D system. For a field of infinite width we can express the spatial joint density as  $\rho(\xi_x, \xi_y) = |\phi_0(\xi_x, \xi_y)|^2$ . The longitudinal and transverse joint densities in different directions are given by  $\rho(\xi_x) = \int d\xi_y \rho(\xi_x, \xi_y)$  and  $\rho(\xi_y) = \int d\xi_x \rho(\xi_x, \xi_y)$ . Since  $\rho(\xi_x = 0, \xi_y = 0)$ ,  $\rho(\xi_x = 0)$ , and  $\rho(\xi_y = 0)$ are all divergent (i.e., (0,0) is a singular point of the density  $\rho$ ) we cannot scale the spatial density distributions to match at  $\xi_x=0$  or/and  $\xi_y=0$ , as carried out above. On the other hand, Eq. (7) shows that, in the dual space we can still obtain the momentum spectrum, which qualitatively reflects the properties of the spatial density. The longitudinal and transverse components of the momentum spectrum are

$$\rho(P_x) = \int dP_y \rho(P_x, P_y) = \frac{2\pi}{(1+P_x^2)^{1/2}} + \frac{2\pi}{(1+P_x^2)^{3/2}},$$
  
$$\rho(P_y) = \int dp_x \rho(P_x, P_y) = \frac{2\pi}{(1+P_y^2)^{1/2}}.$$
 (8)



**Fig. 3.** (Color online) The momentum spectrum  $\rho(P_x)$  and  $\rho(P_y)$  in the 2D case for different W. The circles are the spectra in the 1D case. The data are scaled to match at  $(P_x, P_y) = 0$ .



**Fig. 4.** (Color online) The density distributions  $\rho(\xi_x)$  and  $\rho(\xi_y)$  in the 2D case for  $W = \infty$ . The data are not scaled to match at  $\xi_x$ ,  $\xi_y = 0$  because of the divergence at this point. Instead they are scaled to match at  $\xi_x, \xi_y = 0.00005$ .

In Fig. 3 we show the momentum spectrum in the 2D system for different field widths. For comparison, the corresponding distributions in the 1D system are also shown (circles). Equation (8) is used for the field width  $W = \infty$ . The distributions corresponding to the other field widths are obtained numerically. One sees that the momentum spectra in the 2D system are much wider than that in the 1D system. We can, thus, expect that the density distribution in 2D is much narrower than that in 1D; that is, in the 2D system the electron-positron momentum correlation is weakened, while the spatial correlation is enhanced. Moreover, from Fig. 3 we can see that the transverse spectrum is always wider than the longitudinal one, so that one can expect that  $\rho(\xi_y)$  is always narrower than  $\rho(\xi_x)$ , or the electron-positron spatial correlation perpendicular to the external field is stronger. In order to verify these conclusions, the longitudinal and transverse density distributions of the 2D system are obtained directly by numerically solving Eq. (5) together with the definitions of  $\rho(\xi_x)$  and  $\rho(\xi_y)$ . The results are shown in Fig. 4. Clearly, both  $\rho(\xi_x = 0)$  and  $\rho(\xi_y = 0)$ are divergent. Moreover, we can clearly see that their widths are narrower than that in the 1D case and the width of  $\rho(\xi_y)$  is narrower than that of  $\rho(\xi_x)$ .

Similarly, for the 3D system we can obtain analytically for the field of width  $W = \infty$  the joint momentum spectrum

$$\rho(\mathbf{P}) = 8 \frac{1 + P_y^2 + P_z^2}{(1 + P_x^2 + P_y^2 + P_z^2)^2},$$
(9)

where  $P_z = p_z/c$  and  $\mathbf{P} = (P_x, P_y, P_z)$  and all of the four spinors are taken into account. We can see that  $\rho(P_x)/\rho(P_x = 0) \rightarrow 1$ , so that the width of the momentum spectrum in the parallel direction approaches infinity,  $\Delta P_x \rightarrow \infty$ , and the width of the spatial distribution approaches zero,  $\Delta \xi_x \rightarrow 0$ . The same situation occurs for the perpendicular directions, namely,  $\rho(P_y)/\rho(P_y = 0) \rightarrow 1$ ,  $\rho(P_z)/\rho(P_z = 0) \rightarrow 1$ ,  $\Delta \xi_y \rightarrow 0$ , and  $\Delta \xi_z \rightarrow 0$ . Accordingly, in the 3D system the electron and positron are pair-created at practically the same location. This result differs strongly from that of the 1D case, where the average spatial extent of pair creation is of order  $\lambda_e$ .

The enhancement comes from the coupling influence of the momenta in the different directions, which can be illuminated from Eqs. (5) and (6). For example, in the quasi-2D case the transverse momentum makes the longitudinal momentum spectrum broaden with the corresponding spatial distribution approaching zero slowly as momentum increases compared to the 1D case. In other words, the coupling of entangled wave function due to different dimensions leads to the characteristic that the spatial correlation degree is greatly increased.

In summary, we have considered electron-positron correlations in the pair-creation process during the early stage. Our results suggest that the correlations in configuration and momentum spaces exhibit reciprocal duality, which is consistent with the uncertainty principle; that is,  $\Delta P_x \Delta x \approx 1$ . It is found that the correlations depend on the dimension of the computational space. Specifically, the particle pairs are created almost at the same location in the 3D system while the average distance between the particles is finite in lower-dimensional spaces for the infinite width field. We also find that the transverse spatial correlation is stronger than the longitudinal correlation in the 2D system and in the quasi-2D system, the longitudinal spatial correlation increases with the transverse momentum.

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