

Interacting Electron-Photon System

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- 2 Single Photon System
- 3 Single Electron System
- 4 Two-body, Non-interacting System
- 5 Two-body, Interacting System

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- In Bohmian Mechanics, particles have definite positions that change with time. One wave function, defined in the configuration space of a system of particles, guides the motion of all particles by giving the guiding equation for each particle.
- We conducted our research gradually, first examining the system of a single photon, then that of a single electron, then that of the two without any interaction, and, at last, that of two interacting.

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- Kiessling & Tahvildar-Zadeh also discovered this equation in 2018. It is a Dirac-type equation, and in 1-dim. it reads:

$$-i\hbar\gamma^{\mu}\frac{\partial\Psi_{ph}}{\partial x^{\mu}} = 0,$$

where \hbar = reduced Planck's constant, $x^0 = t, x^1 = s$,

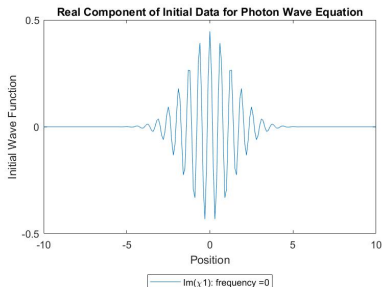
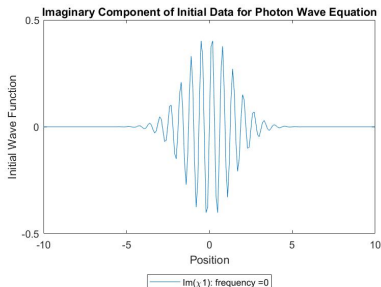
$\gamma^0 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$, $\gamma^1 = \begin{pmatrix} 0 & -1 \\ 1 & 0 \end{pmatrix}$, and repeated indices are summed over the range $\mu = 0, 1$.

Initial Wave Function

- The photon wave equation needs to be solved given an initial wave function $\Psi_{ph}(0, s) = \Psi_{ph}^0(s)$.

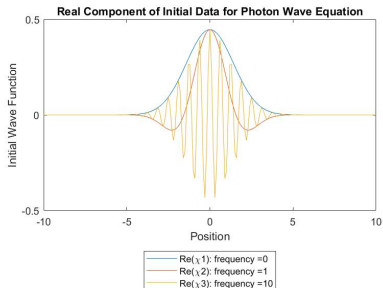
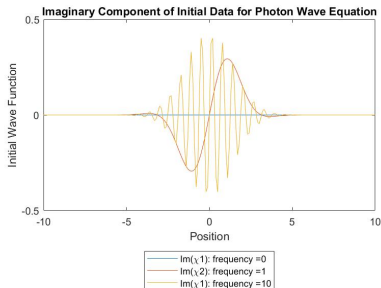
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Photon Probability Current and Velocity Field

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- Here $X = (X^0, X^1)$ is a constant vector field computed from Ψ_{ph}^0 , $\gamma(X) := \gamma_0 X^0 + \gamma_1 X^1$, and $\overline{\Psi} := \gamma^0 \Psi^{\dagger} \gamma^0$.

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- $\rho(t, s)$ is normalized, namely $\int \rho(t, s) ds = 1$.

Photon Probability Density

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http://reu.dimacs.rutgers.edu/~aas377/photon_pdf.mp4

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- Varying initial conditions gives us the following: http://reu.dimacs.rutgers.edu/~aas377/multiple_photon_pdf.mp4

The Guiding Equation

- The motion of the photon is guided by its wave function:

$$\begin{cases} \frac{dq}{dt} = v_{ph}(t, q(t)) = \frac{j^1(t, q(t))}{j^0(t, q(t))} \\ q(0) = q_0 \end{cases}$$

where $q(t)$ is the actual position of the photon at time t .

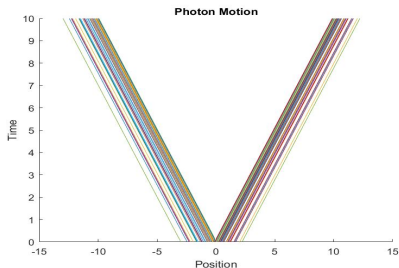
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- q_0 is the actual initial position of the photon. All we know about it is that it is randomly distributed according to the initial probability density $\rho(0, s)$.



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- Like in the case of a single photon, the wave function of a single electron also satisfies a relativistic equation. In particular, it satisfies the massive Dirac equation:

$$-i\hbar\gamma^{\mu}\partial_{\mu}\Psi_{el} + m_{el}\Psi_{el} = 0,$$

where m_{el} = the mass of electron.

Electron Probability Current and Velocity Field, and Guiding Equation

- The probability current of an electron is known:

$$j_{el}^{\mu}(time, position) = \overline{\Psi}_{el} \gamma^{\mu} \Psi_{el}$$

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- Similarly to the photon case, the guiding equation for the electron is:

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Electron Trajectories and Parameters

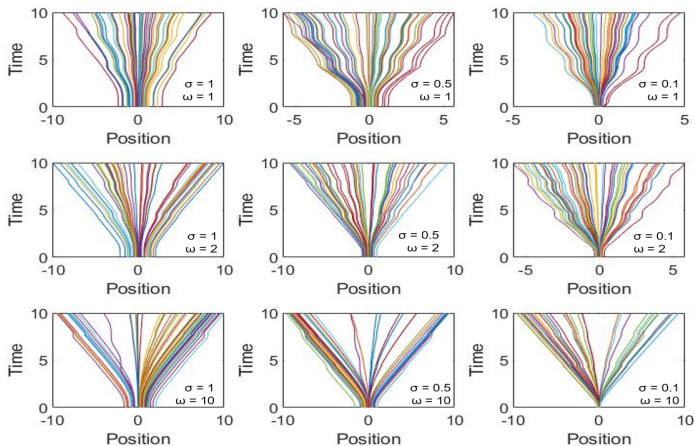
- The electron wave function Ψ_{el} , has a mass term: $\omega = \text{mass}/\hbar$, and a parameter we can change: standard deviation of the initial distribution: σ . The following graph shows the trajectory of an electron guided by the velocity field with different parameters.

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- To get a wave function that describes both a photon and an electron in a non-interacting system, we take the Tensor Product (\otimes) of the electron and the photon wave functions, giving us a four component object $\psi = (\psi_{++}, \psi_{+-}, \psi_{-+}, \psi_{--})$

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- The guiding equations for photon and electron are derived using the Hypersurface Bohm-Dirac (HBD) Theory, which allows us to describe the motion of the photon and electron in a common time

Wave Equation and Probability Current

- The tensored wave function satisfies a relativistic wave equation obtained by Tensor Product of the photon and electron wave equations

$$\begin{cases} -i\hbar\gamma^\mu\partial x_{ph}^\mu\psi = 0 \\ -i\hbar\gamma^\mu\partial x_{el}^\mu\psi + m_{el}\psi = 0 \\ \psi(0, s_{ph}, 0, s_{el}) = \psi^0(s_{ph}, s_{el}) \end{cases}$$

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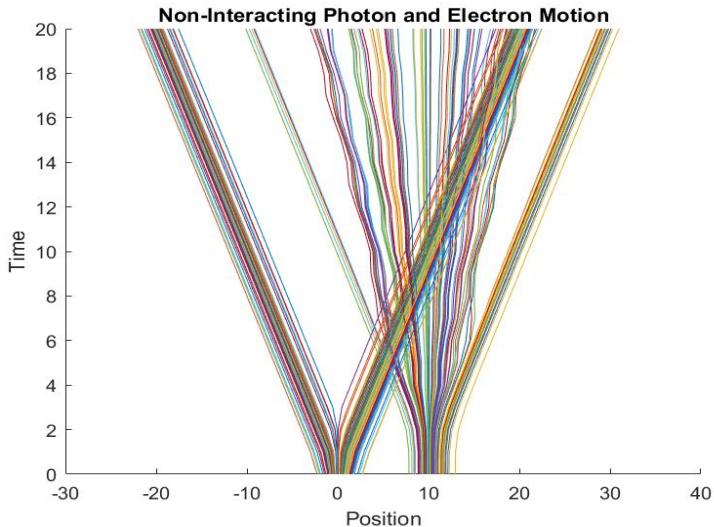
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- The probability current is the following:

$$\begin{pmatrix} j^{10} & = & |\psi_{++}|^2 + |\psi_{+-}|^2 - |\psi_{-+}|^2 - |\psi_{--}|^2 \\ j^{01} & = & |\psi_{++}|^2 - |\psi_{+-}|^2 + |\psi_{-+}|^2 - |\psi_{--}|^2 \\ j^{00} & = & |\psi_{++}|^2 + |\psi_{+-}|^2 + |\psi_{-+}|^2 + |\psi_{--}|^2 \\ j^{11} & = & |\psi_{++}|^2 - |\psi_{+-}|^2 - |\psi_{-+}|^2 + |\psi_{--}|^2 \end{pmatrix}$$

Trajectories of the Two-body, Non-interacting System

- The following graph shows the trajectories of a non-interacting system of one electron and one photon.



Chapter 4: Two-body, Interacting System

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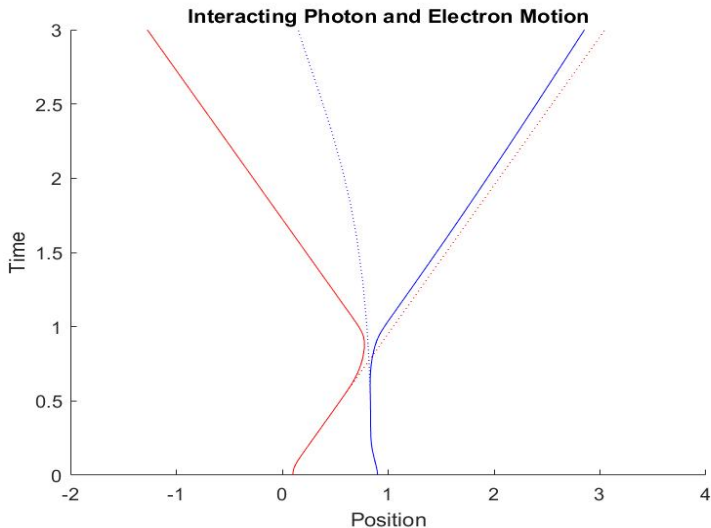
- To obtain an interacting system from a non-interacting system, it is necessary to add some conditions such that the particles do not simply go through each other.
- We do this by adding a boundary condition: we set the relative velocities of photon and electron to be 0 when the particles are at the same space and time.

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- We do this by adding a boundary condition: we set the relative velocities of photon and electron to be 0 when the particles are at the same space and time.
- Adding the boundary condition to the wave function gives us a modified probability density function: http://reu.dimacs.rutgers.edu/~aas377/interacting_pdf_2.mp4

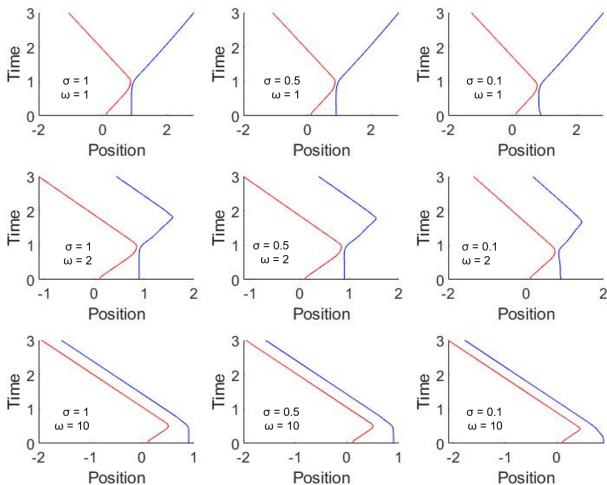
Trajectories of the Two-body, Interacting System

- Adding the boundary condition to the wave function gives us the trajectories of an interacting electron-photon system.



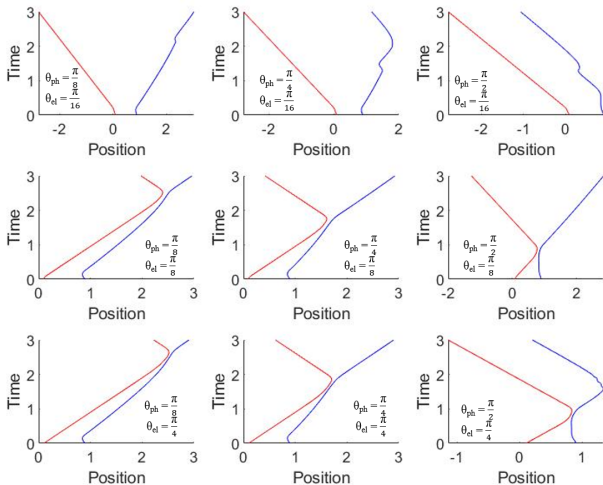
Varying Parameters in the Interacting System

- Changing the sigma and omega of the electron gives us the following changes in trajectories of the electron:



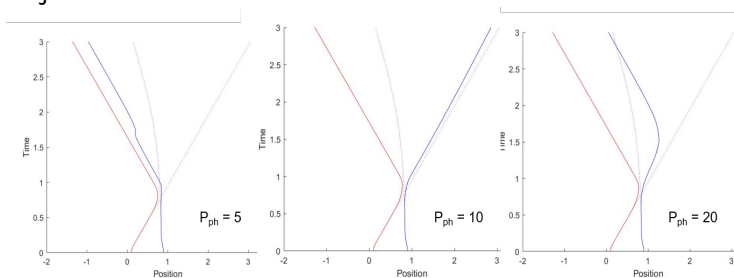
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- Changing the polarization angles of electron and photon gives us:



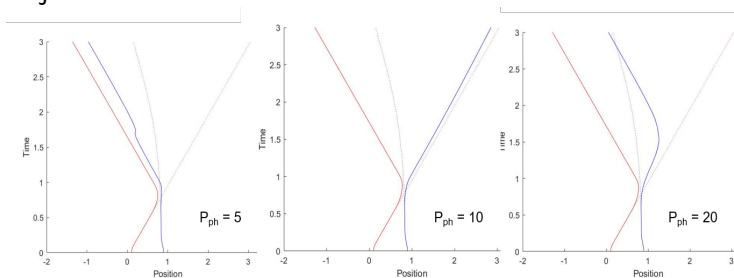
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Varying Parameters in the Interacting System

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- Momentum is related to energy, so if the photon does not have enough momentum, it cannot get the electron to bounce away.

Thank You

- This research is made possible by the Rutgers Math Department, REU-DIMACS, and the generous help from Professor Shadi Tahvildar-Zadeh