Interacting Electron-Photon System

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Outline

Introduction

- 2 Single Photon System
- 3 Single Electron System
- 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 on the configuration space of a system of particles, guides the motion of all particles through their respective guiding equations.
- First we examined the system of a single photon, then that of a single electron, then that of the two without any interaction, and, at last, the system of the photon and electron 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:

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

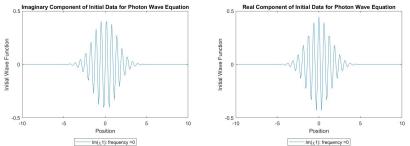
where $\hbar =$ reduced Planck's constant, $x^0 = t, x^1 = s$, $\gamma^0 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \ \gamma^1 = \begin{pmatrix} 0 & -1 \\ 1 & 0 \end{pmatrix}$, and repeated indices are summed over the range $\mu = 0, 1$.

Initial Wave Function

 The photon wave equation needs to be solved given an initial wave function Ψ_{ph}(0, s) = Ψ⁰_{ph}(s).

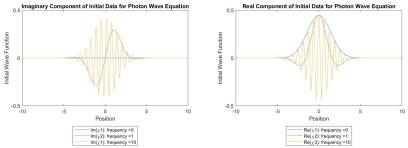
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• Here $X = (X^0, X^1)$ is a constant vector field computed from Ψ^0_{ph} , $\gamma(X) := \gamma_0 X^0 + \gamma_1 X^1$, and $\overline{\Psi} := \gamma^0 \Psi^{\dagger} \gamma^0$.

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- The current is conserved: $\partial_{\mu}j^{\mu}_{ph} = 0$, future directed $(j^0 \ge 0)$, and timelike $(j^0 \ge |j^1|)$.
- The probability density of detecting the photon at event (t, s) is $\rho(t, s) = j_{ph}^0(t, s)$.

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Photon Probability Density

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- The photon probability density looks like this: http://reu.dimacs.rutgers.edu/~aas377/photon_pdf.mp4
- 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.

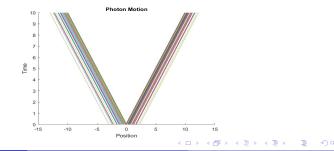
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• q_0 is the actual initial position of the photon. All we know about it is that it is randomly distributed according to the initial probability density $\rho(0, s)$.



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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:

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

where $\overline{\Psi}:=\Psi^\dagger\gamma^0$ is the Dirac adjoint for rank-one bispinors.

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

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Electron Probability Density

• http://reu.dimacs.rutgers.edu/~aas377/electron_pdf.mp4

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Electron Trajectories and Parameters

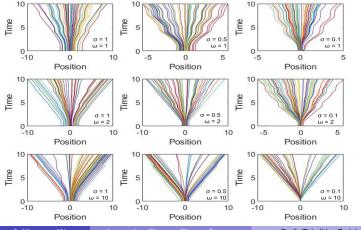
The electron wave function Ψ_{el}, has a mass term: ω =mass/ħ, and a parameter we can change: standard deviation of the initial distribution: σ.The following graph shows the trajectory of an electron guided by the velocity field with different parameters.

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- The wave function, ψ , is a function of four variables, namely the time and position of each particle.
- To get a wave function that describes both a photon and an electron in a non-interacting system, we take the Tensor Product (⊗) of the electron and the photon wave functions, giving us a four component object ψ = (ψ₊₊, ψ_{+−}, ψ_{-−})

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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

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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

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

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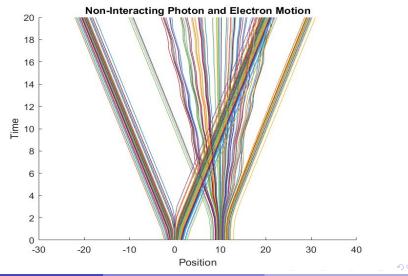
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$$\begin{pmatrix} j^{00} &= |\psi_{++}|^2 + |\psi_{+-}|^2 + |\psi_{-+}|^2 + |\psi_{--}|^2 \\ j^{10} &= |\psi_{++}|^2 + |\psi_{+-}|^2 - |\psi_{-+}|^2 - |\psi_{--}|^2 \\ j^{01} &= |\psi_{++}|^2 - |\psi_{+-}|^2 + |\psi_{-+}|^2 - |\psi_{--}|^2 \\ j^{11}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

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

• 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.

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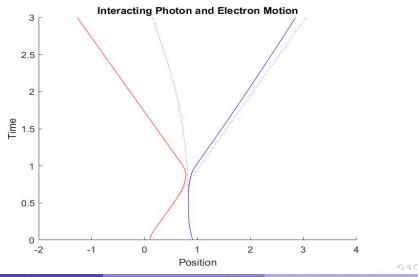
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- We do this by adding a boundary condition: we set the relative velocities of photon and electron to be 0 when the particles are at the same space and time.
- Adding the boundary condition to the wave function gives us a modified probability density 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.

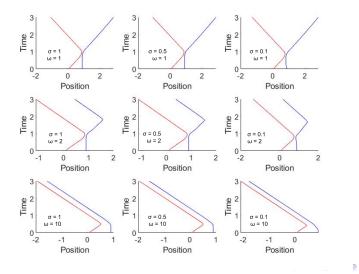


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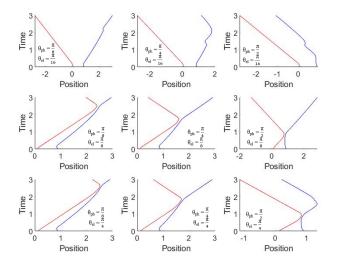
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• Changing the sigma and omega of the electron gives us the following changes in trajectories of the electron:

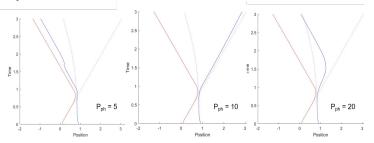


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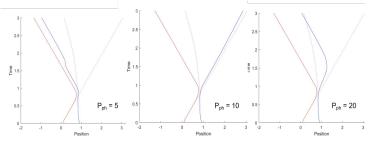
• Changing the polarization angles of electron and photon gives us:



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

Future Steps

• We need to continue performing numerical experiments in order to clarify the roles of different parameters

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- We need to continue performing numerical experiments in order to clarify the roles of different parameters
- In order to verify our framework, we need to find a way to calculate momentum after the photon-electron collision and compare our findings to the results of Compton Scattering in one space dimension

Thank You

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

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- For more information, visit us at CoRE 417!