

The Hydrogen Atom in Wave Mechanics

Michael Brodskiy

Professor: Q. Yan

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Contents

1	A One-Dimensional Atom	3
2	Angular Momentum in the Hydrogen Atom	3
3	The Hydrogen Atom Wave Functions	4

1 A One-Dimensional Atom

- Analyzing a Hydrogen Atom in Quantum Mechanics

- The potential energy, derived from the Coulomb force, is:

$$U(r) = -\frac{e^2}{4\pi\epsilon_0 r}$$

- This may be converted to be in terms of x , and plugged into the Schrödinger equation:

$$-\frac{\hbar^2}{2m} \frac{d^2\psi(x)}{dx^2} - \frac{e^2}{4\pi\epsilon_0 x} \psi(x) = E\psi(x)$$

- To keep this finite, the following two conditions need to be met:

$$\begin{cases} x \rightarrow 0, & \psi(x) = 0 \\ x \rightarrow \infty, & \psi(x) = 0 \end{cases}$$

- Using $\psi(x) = Axe^{-bx}$ with the Schrödinger equation, solving would obtain:

$$b = \frac{me^2}{4\pi\epsilon_0 \hbar^2} = \frac{1}{\alpha_o} \Rightarrow \text{Bohr Radius}$$

- Using the normalization condition, and $\psi(x) = Axe^{-bx} = Axe^{\frac{x}{\alpha_o}}$, we get:

$$A = \frac{2}{\alpha_o^{\frac{3}{2}}}$$

- Conclusions:

- * For the ground state, there is an uncertainty in the location of e^-
- * The most probable region to find e^- is near $x = \alpha_o$ (consistent with Bohr model)
- * But, it's possible to find e^- anywhere (which is very difference from Bohr)

2 Angular Momentum in the Hydrogen Atom

- The angular momentum in a planetary system is constant, and vector $\vec{L}(L_x, L_y, L_z)$ has three components
 - * l is the angular momentum quantum number; it determines the length of the vector
 - * m_l is the magnetic number; it determines one of the components of the vector

* We have:

$$|\vec{L}| = \sqrt{l(l+1)}\hbar, \quad l = 0, 1, 2, \dots$$

$$\vec{L}_z = m_l \hbar, \quad m_l = 0, \pm 1, \pm 2, \dots \pm l$$

* We know l and m_l for \vec{L}_z , but what is the direction?

* Applying the uncertainty principle, we obtain:

$$\Delta \vec{L}_z \Delta \phi \geq \hbar$$

3 The Hydrogen Atom Wave Functions

- Moving from a unidimensional case to a tridimensional case, we try to apply the Schrödinger equation (in Cartesian coordinates):

$$-\frac{\hbar^2}{2m} \left(\frac{\partial^2 \psi}{\partial x^2} + \frac{\partial^2 \psi}{\partial y^2} + \frac{\partial^2 \psi}{\partial z^2} \right) + U(x, y, z) \psi(x, y, z) = E \psi(x, y, z)$$

- The potential energy is:

$$U(x, y, z) = -\frac{e^2}{4\pi\epsilon_0} \frac{1}{\sqrt{x^2 + y^2 + z^2}}$$

- The wave function may be defined as:

$$\psi(x, y, z) = X(x)Y(y)Z(z)$$

- Converting to polar coordinates to make calculations simpler, the Schrödinger equation becomes:

$$-\frac{\hbar^2}{2m} \left(\frac{\partial^2 \psi}{\partial r^2} + \frac{2}{r} \frac{\partial \psi}{\partial r} + \frac{1}{r^2 \sin(\theta)} \frac{\partial}{\partial \theta} \left(\sin(\theta) \frac{\partial \psi}{\partial \theta} \right) + \frac{1}{r^2 \sin^2(\theta)} \frac{\partial^2 \psi}{\partial \phi^2} \right) + U(r) \psi(r, \theta, \phi) = E \psi(r, \theta, \phi)$$

- The wave function then becomes

$$\psi(r, \theta, \phi) = R(r)\Theta(\theta)\Phi(\phi)$$

– Where R is the radial function, Θ is the polar function, and Φ is the azimuthal function

– This breaks ψ into 3 equations, making the wave function easier to solve for

- The defined quantum numbers are as follows:

Number	Name	Values
n	Principle	$1, 2, 3, \dots$
l	Angular Momentum	$0, 1, 2, \dots, n - 1$
m_l	Magnetic	$0, \pm 1, \pm 2, \dots, \pm l$

- When integrating:

$$|\psi|^2 dV = |R|^2 |\Theta|^2 |\Phi|^2 r^2 \sin(\theta) dr d\phi$$

- Intrinsic angular momentum (electron spin)

– $\vec{L} = \vec{r} \times \vec{p}$ is the orbital angular momentum

– \vec{s} : the electron spin

- The idea of spin was proposed in 1925

– In 1928, Paul Dirac developed the relativistic quantum theory, and determined a 4th quantum number as electron spin

– $\vec{s} = \frac{1}{2} \iff l = 0, 1, 2, \dots$

– The length of \vec{s} becomes $\sqrt{\vec{s}(\vec{s} + 1)}\hbar$

– Now the electron state may be represented by “spectroscopic notation”

$$(n, l, m_l, m_s)$$