Derivation of the Pauli Exclusion Principle and Meaning Simplification of the Dirac Theory of the Hydrogen Atom

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Abstract: In generally, the Pauli Exclusion Principle follows from the spectroscopy whereas its origin is not good understood. To understand fully this principle, most important is origin of quantization of the azimuthal quantum number i.e. the angular momentum quantum number. Here, on the base of the theory of ellipse and starting from very simple physical condition, I quantized the azimuthal quantum number. The presented model leads directly to the eigenvalue of the square of angular momentum and to the additional potential energy that appears in the equation for the modified wave function. I formulated the very simple semiclassical analog to the Dirac and Sommerfeld theories of the hydrogen atom. The constancy of the base of the natural logarithm for the quantum fields is the reason that the three theories are equivalent.

1. Introduction

The Pauli Exclusion Principle says that no two identical half-integer-spin fermions may occupy the same quantum state simultaneously. For example, no two electrons in an atom can have the same four quantum numbers. They are the principal quantum number $n$ that denotes the number of the de Broglie-wave lengths $\lambda$ in a quantum state, the azimuthal quantum number $l$ (i.e. the angular momentum quantum number), the magnetic quantum number $m$ and the spin $s$.

On the base of the spectrums of atoms, placed in magnetic field as well, follows that the quantum numbers take the values:

$n = 1, 2, 3, \ldots$
$l = 0, 1, 2, \ldots n - 1$
$m = -l, \ldots +l$
$s = \pm 1/2.$

The three first quantum numbers $n$, $l$, and $m$ are the integer numbers and define a state in which can be maximum two electrons with opposite spins.

The magnetic quantum number $m$ determines the projection of the azimuthal quantum number $l$ on the arbitrary chosen axis. This axis can overlap with a diameter of the circle $l = 0$.

To understand fully the Pauli Exclusion Principle we must answer following questions concerning the azimuthal quantum number $l$:

1. What is physical meaning of this quantum number?
2. Why the $l$ numbers are the natural numbers only?
3. Why the zero is the lower limit?
4. Why the \( n - 1 \) is the upper limit?

To answer these questions we must apply the theory of ellipse, especially the formula for its circumference \( C \) and eccentricity \( e \). When we use the complete elliptic integral of the second kind and the Carlson symmetric form [1], we obtain for circumference \( C \) of an ellipse following formula

\[
C = 2\pi a \left[ 1 - \frac{1}{2} e^2 - \frac{1}{3} \left( 1 - \frac{1}{2} \right) e^4 - \left( 1 - \frac{1}{2} \right) \frac{1}{4} e^6 - \ldots \right],
\]

where \( a \) is the major radius and \( e \) is the eccentricity defined as follows

\[
e = \frac{\sqrt{a^2 - b^2}}{a},
\]

where \( b \) is the minor radius.

Here, I formulated as well the very simple semiclassical analog to the Dirac and Sommerfeld theories of the hydrogen atom and proved that these three theories are equivalent.

2. Calculations
2.1 Angular momentum quantum number

In the figure, the circumference of the ellipse \( C_{\text{de-Broglie}} \) is \( C_{\text{de-Broglie}} = n\lambda = 2\pi n\lambda \), where the \( n \) is the principal quantum number whereas the \( \lambda \) is the reduced de Broglie-wave length.

Assume that there are allowed only ellipses that circumference is the arithmetic mean of the circumferences of two circles that radii are equal to the major and minor radii of the ellipse.

Similarly as for the circumference of the ellipse, the circumferences of the circles must be equal to a natural number multiplied by the de Broglie-wave length. This leads to following definitions

\[
a = j\lambda \text{ and } b = k\lambda.
\]

Notice that \( j = k = 0 \) has no sense.

Then, we can rewrite formula (2) as follows

\[
e = \frac{\sqrt{j^2 - k^2}}{j}.
\]

It is the natural assumption that the allowed circumferences of the ellipse should be the arithmetic mean of the sum of the circumferences of the two circles. It leads to following conclusion

\[(j + k)/2 = n.\]

Define some number \( l \) as follows

\[(j - k)/2 = l.\]
Formulae (5) and (6) lead to following relations

\[ j = n + l, \]  
\[ k = n - l. \]  
(7)  
(8)

Since the \( j, k \) and \( n \) are the integers so the number \( l \) must be an integer as well.

On the base of formulae (7) and (8) we can rewrite formula (4) as follows

\[ e = 2\sqrt{nl}/(n + l). \]  
(9)

We can see that due to the square root, this formula has no real sense for \( l < 0 \). Since the \( l \) cannot be negative then from formulae (5) and (6) follows that \( l < n \).

On the base of formulae (3) and (7), we can rewrite formula (1) as follows

\[ C_K = 2\pi(n + l)\lambda[1 - (1/2)^2e^2/1 - (1-3/(2-4))^2e^4/3 - (1-3-5/(2-4-6))^2e^6/5 - \ldots]. \]  
(10)

Notice that for \( n = l \) is \( e = 1 \) and then \( C_{\text{de-Broglie}} > C_K \) i.e. \( l \) cannot be equal to \( n \). For \( l = 0 \) is \( C_{\text{de-Broglie}} = C_K \) and because \( l \) cannot be negative then the \( l = 0 \) is the lower limit for \( l \).

Some recapitulation is as follows. We proved that the azimuthal quantum number \( l \)

1) is associated with transitions between the states \( j \) and \( k \),
2) is the integer,
3) cannot be negative and the lower limit is zero,
4) the \( n - 1 \) is the upper limit.

Some abbreviation of it is as follows

\( l = 0, 1, 2, \ldots n - 1. \)

The Quantum Physics is timeless because a quantum particle disappears in one region of a field or spacetime and appears in another, and so on. There are no trajectories of individual quantum particles. Quantum Physics is about the statistical shapes and their allowed orientations. Such procedure simplifies considerably the Quantum Physics.

2.2 Eigenvalue of the square of angular momentum

An ellipse/electron-state we can resolve into two circles that radii are defined by the semi-
axes of the ellipse. The two circles in a pair are entangled due to the exchanges of the binary
systems of the closed strings (the entanglons [2]) the Einstein-spacetime components, from
which are built up all the Principle-of-Equivalence particles, consist of [2]. Spin of the
entanglons is 1 [2] and they are responsible for the infinitesimal transformations that lead to
the commutators [3]. Calculate a change in the azimuthal quantum number \( l \) when the smaller
circle or one of identical two circles emits one entanglon (since in this paper is \( j \geq k \) so there is
the transition \( k \to k - 1 \) whereas the second circle in the pair almost simultaneously absorbs
the emitted entanglon (there is the transition \( j \to j + 1 \)). Such transition causes that ratio of the
major radius to the minor radius of the ellipse (or circle) increases. From formula (5) follows
that such emission-absorption does not change the principal quantum number \( n \) whereas from
formula (6) follows that there is following transition for the azimuthal quantum number \( l \):
\( l \to l + 1 \). The geometric mean is \( \sqrt{(l(l + 1))} \) and this expression multiplied by \( \hbar \) is the mean
angular momentum \( L \) for the described transition. This leads to conclusion that eigenvalue of
the square of angular momentum \( L^2 \) is \( l(l + 1)\hbar^2 \).

The eigenvalue of the square of angular momentum leads to the additional potential energy
\( E_A \) (it follows from the radial transitions i.e. from the changes in shape of the ellipses) equal to

\[ E_A = L^2/(2m^2) = l(l + 1)\hbar^2/(2m^2). \]  
(11)

The energy \( E_A \) appears in the equation for the modified wave function.

The theory of baryons [2] shows that inside the baryons are only the \( l = 0 \) states (i.e. there
are only the circles) so the quantum mechanics describing baryons is much simpler than for
atoms.
2.3 The very simple semiclassical analog to the Dirac and Sommerfeld theories of the hydrogen atom

Since the fermions consist, at least for period of spinning, of the stable/classical structures/bare-fermions and of the quantum fields so the semiclassical theories are simplest, most fruitful and contain least parameters. And such method is not a mathematical trick – just in such a way behaves Nature.

Gravity is associated with the inverse square law. It is because gravitational fields are the gradients produced in the modified Higgs field by masses [2]. There are the divergently moving classical tachyons so there appears the inverse square law

\[ F \sim \frac{1}{r^2}. \] (12)

Today, in the modified Higgs field cannot be created any virtual pairs as it is in the Einstein spacetime [2]. The produced structures in the Einstein spacetime and associated with them virtual pairs cause that field \( \psi \) of the virtual pairs changes according following function (see Paragraphs “Mathematical Constants” and “Fractal Field” here [2])

\[ \psi = \psi_0 e^{-r}. \] (13)

where \( e = 2.7182818… \) is the base of the natural logarithm. In reality, this formula is more complicated for \( r \to 0 \) because there appears torus/charge/spin and central condensate [2]. It causes that the quantum physics is the incomplete theory. By the way, notice that, for example, we define mean-lifetime in such a way that after this time there do not decay \( 1/e \) particles. Amplitude \( \psi_0 \) should be in proportion to coupling constant characteristic for a field i.e. \( \psi_0 \sim \alpha \). On the other hand, a physical meaning has following expression

\[ \psi_0 \psi_0 \sim \alpha^2. \] (14)

Formulae (12) and (13) and associated with them the sets of applied methods differ very much so unification of gravity and quantum physics within one of the two sets is impossible.

Due to the gluons \( \to \) photons transitions on edges of the strong fields, there leaks the internal structure of nucleons but, of course, there change the coupling constants. On the other hand, the mass of nucleons dominates in atoms. The coupling constant for the strong interactions of pions is \( \alpha_s = 1 \) [2], whereas for weak interactions of baryons is \( \alpha_W = 0.0187229 \) [2]. Define following expression

\[ X = e^{-1} \alpha = e^{-1}(\alpha_S + \alpha_W) = 0.37476… \] (15)

Can we quantize the value \( X \) i.e. can we write an expression that leads to \( X \)? Notice that

\[ Y = m(1 - 1/2 - 1/8)/m_{\text{electron}} = 0.37479…, \] (16)

where \( m \) is the reduced mass of electron i.e. \( m/m_{\text{electron}} = m_{\text{proton}}/(m_{\text{proton}} + m_{\text{electron}}) = 0.9994557. \)

It leads to following expression

\[ Y = m(1 - \Sigma_{d=1,2} 1/(2^{2d-1}))/m_{\text{electron}} = 0.37479… \] (17)

In the Quantum Theory of Fields (QTF) we apply Lagrangians (energy) in which as well appear products of two generators and each generator is proportional to coupling constant i.e. there are components for which energy is proportional to \( \alpha^2 \). This remark and formula (14) suggests that we should expand energy into series type \( (\alpha/n)^{2d} \), where \( d = 1, 2 \), whereas \( n = 1, 2, 3, 4,… \). For the hydrogen atom we obtain

\[ E = m c^2 \left[ 1 - \left(\alpha_{\text{em}}/n\right)^2/2 - \left(\alpha_{\text{em}}/n\right)^4/8 \right], \] (18)

where \( m c^2 = 0.5107208 \) MeV is the reduced mass of electron, \( n \) is the principal quantum number whereas \( \alpha_{\text{em}} = 1/137.036 \) is the fine-structure constant.
The first component \( E_R = mc^2 \) is the internal energy concerning the reduced mass of electron.

The second component

\[
E_{B,n} = -mc^2\left(\frac{\alpha_{em}}{n}\right)^2/2
\]

(19)
is equal to the energies of the Bohr orbits in the hydrogen atom and \( E_{B,n=1} = -13.598 \text{ eV} \).

The third component is the fine structure energy

\[
E_{FS,n} = -mc^2\left(\frac{\alpha_{em}}{n}\right)^4/8.
\]

(20)

This component depends on classical and quantum structure of electron so we must write it in such a way to interpret it correctly. Write the factor 1/8 as follows

\[
1/8 = \left(\frac{1}{2} - \frac{3}{4}\right)/2.
\]

(21)

The 3/4 represents the classical mass of electron [2] (see Chapter “Foundations of Quantum Physics”; the factor 3/4 follows from the internal structure of the bare electron – there is the torus/electric charge) whereas the unity represents the quantum mass of electron. We can see that we separated these two very different masses. We know that maximum azimuthal quantum number \( l \) is \( l_{\text{max}} = n - 1 \) so \( n/(l_{\text{max}} + 1) = 1 \). This means that we can rewrite formula (21) as follows

\[
1/8 = \left[\frac{n}{l_{\text{max}} + 1}\right] - \frac{3}{4}/2.
\]

(22)

The \( n \) and \( l_{\text{max}} + 1 \) define the lengths of the de Broglie waves but the additional potential energy \( E_A = l(l + 1)\hbar^2/(2mr^2) \) suggests that for defined \( n \) there can appear spontaneously as well the other standing waves defined by \( l + 1 \). For smaller \( l \) waves are shorter so corresponding absolute energy is greater. Since in formula (20) is the sign “–” so the levels defined by smaller and smaller \( l \) are closer and closer to the ground state \( n = 1 \). Finally, we can rewrite formula (20) as follows

\[
E_{FS,n} = -mc^2\left(\frac{\alpha_{em}}{n}\right)^4\left[\frac{n}{l_{\text{max}} + 1}\right] - \frac{3}{4}/2.
\]

(23)

The ground state is shifted by \( E_{FS,n=1} = -mc^2\left(\frac{\alpha_{em}}{1}\right)^4\left[\frac{1}{2} - \frac{3}{4}\right]/2 = -1.81 \cdot 10^{-4} \text{ eV} \).

Calculate the energy distance between the states \( l = 0, 1 \) for \( n = 2 \). The general solution for the energy distance between the extreme levels for defined \( n \) is as follows

\[
\Delta E_{FS,n} = -mc^2\left(\frac{\alpha_{em}}{n}\right)^4(n - 1)/2.
\]

(24)

From it we obtain \( \Delta E_{FS,n=2} = -mc^2\left(\frac{\alpha_{em}}{2}\right)^4(2 - 1)/2 = -mc^2\left(\frac{\alpha_{em}}{2}\right)^4/32 = -4.53 \cdot 10^{-5} \text{ eV} \).

Why we obtained results the same as in the Sommerfeld theory [4]? Why we obtained results the same as in the Dirac theory [5] neglecting the relativistic effects, the spin-orbit interactions, and so on? It follows from the fact that for the quantum fields is \( X \approx Y \). It follows from the constancy of the base of the natural logarithm for the quantum fields. It is due to the applied methods – just the standing waves defined by the quantum numbers cannot by changed by any phenomena. Just the quantum numbers define the total picture and must be conserved. The three theories are equivalent because the numbers \( n_0 \) in the Sommerfeld theory, \( j + 1/2 \) in the Dirac theory and \( l + 1 \) in presented here theory, are the integers and change from 1 to \( n \). But only presented here theory of hydrogen atom proves equivalence of the three theories and describes in all respects physical origin of the final equation.

The Lamb-Retherford shift is associated with the internal structure of proton and I calculated it here [2]. Since the all levels inside baryons are the \( l = 0 \) states so the Lamb-Retherford shift, due to the resonance, concerns only the \( l = 0 \) states in atoms. This shift is the energy distance between the \( l = 0 \) and \( l = 1 \) states for the same \( n \) and \( j \). This shift decreases binding energy in the \( l = 0 \) state.
Notice that absolute value of the second component (the Bohr theory of atoms) for coupling constant for the strong interactions of pions ($\alpha = 1$) is $E_{B,n=1,\alpha=1} = 0.25536$ MeV and it is in approximation the mass of the torus/electric-charge inside the bare electron whereas for the coupling constant for the strong interactions of the nucleons at very low energy ($\alpha = 14.4$ [2]) we obtain $E_{B,n=1,\alpha=14.4} = 52.95$ MeV and it is in approximation the mass of the torus/electric-charge of muon [2].

3. Summary

In generally, the Pauli Exclusion Principle follows from the spectroscopy whereas its origin is not good understood. To understand fully this principle, most important is origin of quantization of the azimuthal quantum number i.e. the angular momentum quantum number. Here, on the base of the theory of ellipse and starting from very simple physical condition, I quantized the azimuthal quantum number. The presented model leads directly to the eigenvalue of the square of angular momentum and to the additional potential energy that appears in the equation for the modified wave function.

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I formulated as well the very simple semiclassical analog to the Sommerfeld and Dirac theories of the hydrogen atom. The Sommerfeld, Dirac and presented here theories of hydrogen atom are equivalent due to the constancy of the base of the natural logarithm for the quantum fields. The standing waves defined by the quantum numbers cannot by changed by any phenomena. Just the quantum numbers define the total picture and must be conserved. The three theories are equivalent because the numbers $n_\theta$ in the Sommerfeld theory, $j + 1/2$ in the Dirac theory and $l + 1$ in presented here theory, change from 1 to $n$. But only presented here theory of hydrogen atom proves equivalence of the three theories and describes in all respects physical origin of the final equation.

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