

**Study of Schrödinger Particles under the Influence of
a Kratzer Molecular Potential in a Global Monopole
Spacetime Background Plus a Wu-Yang Magnetic
Monopole**

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ABSTRACT

Global monopoles are topological defects predicted by the Grand Unified Theory throughout the transition phase of the primordial universe. They are formed as a byproduct of spontaneous global symmetry breaking. Their behavior resembles that of the elementary particles with their energies concentrated close to the core of the monopole. They do not have gravitational interactions but they are known to modify spacetime geometry. It is an interesting topic to study Schrödinger particles in a Kratzer Molecular potential in the Global Monopole Spacetime background and under the influence of a Wu-Yang Magnetic Monopole. Some modifications in the spectroscopic structure of such molecular interactions have been observed and reported. This study presents the use of the Schrödinger equation to precisely describe the motion of a particle in the presence of Kratzer potential and a Wu-Yang magnetic monopole. The radial equation was derived and solved to yield the eigenfunctions and eigenvalues for the problem.

Keywords: PDM Schrödinger Equation, Point-like Global Monopole, Topological Defect, Kratzer Potential, Wu-Yang Magnetic Monopole.

ÖZ

Global monopoller, Büyük Birlesik Teori tarafından erken evrenin geis fazında öngörülen topolojik kusurlardir. Bu topolojik kusurlar, spontan global simetri kılmasının bir sonucu olarak ortaya ikarlar. Davranslari elementer paraciklarinkine benzer; enerjileri monopolün gekirdegine yakin yogunlasir. Gravitasyonel etkileşim göstermese de. uzay-zaman geometrisini degistirdikleri bilinir. Global Monopole uzay-zaman arka planında ve Wu-Yang Manyetik Monopolden etkilenen Kratzer moleküler potansiyel içinde Schrödinger paraciklarnin incelenmesi ilging bir konudur Bu moleküler etkileşimlerin spektroskopik yapısında bazı ayarlamalar gözlemlenmiş ve belgelenmiştir. Bu alıřma, Global Monopole uzay-zamanında, Kratzer potansiyeli ve Wu-Yang manyetik monopolden etkilenen kosullarda, göreli olmayan bir kuantum paraciginin hareketini kesin olarak tanımlamak için Schrödinger denklemini kullanir. Karsilk ge-len radyal denklem, problemin özfonksiyonların ve özdeğerlerini belirlemek için türetilmiş ve özölmüştür.

Anahtar Kelimeler: PDM Shrödinger Denklemi, Nokta Benzeri Küresel Tek Kutup, Topolojik Kusur, Kratzer Moleküles Potansiyel, Wu-Yang Manyetik Tek Kutup.

DEDICATION

...To my father, without whom I would not be the same person I am now. For his boundless love, endless patience and unwavering support have been my anchor during tough times. To my family for their support, to my friends for their love. To everyone who has taught me, believed in me and given me opportunities. I dedicate this work with gratitude to each one of you.

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Chapter 1

INTRODUCTION

A classical particle of constant mass m moving on the x -axis can be described by its location $\mathbf{x}(t)$ and velocity $\mathbf{v}(t) = \frac{d\mathbf{x}}{dt}$. The remaining dynamical and/or kinematical features of this system only depend on the variables x and v , such as linear momentum $\mathbf{p} = m\mathbf{v}$, kinetic energy $T = \frac{1}{2}mv^2$, and total energy $E = T + V$, Where V represents the potential energy. Hence, the net force exerted on such a particle is given by $\mathbf{F} = \frac{d\mathbf{p}}{dt} = m\mathbf{a}$. In non-relativistic quantum mechanics, on the other hand, the quantum particle is described by a wave function $\psi(x, t)$ that includes all information about the particle and is solved through the Schrödinger equation [1, 2]:

$$-\frac{\hbar^2}{2m} \frac{\partial^2 \psi}{\partial x^2} + V\psi = i\hbar \frac{\partial \psi}{\partial t}. \quad (1.1)$$

Here $i = \sqrt{-1}$ and \hbar is Planck's constant.

However, when the mass becomes effectively position-dependent, the dynamics of both classical and quantum particles dramatically change. In classical mechanics, for example, Mustafa [3] has stated that the net force under PDM conditions is no longer represented by the rate of change of linear momentum \mathbf{p} with respect to time (i.e., $\mathbf{F} \neq \frac{d\mathbf{p}}{dt}$). The net force is instead given as:

$$\mathbf{F} = \sqrt{m(x)} \frac{d}{dt} \left(\sqrt{m(x)} \dot{\mathbf{x}} \right) = m(x) \ddot{\mathbf{x}} + \frac{1}{2} m'(x) \dot{\mathbf{x}}^2. \quad (1.2)$$

Hence, Newton's second law in the standard one-dimensional format takes the form

$$m(x) \ddot{x} = F - \frac{1}{2} m'(x) \dot{x}^2; \quad m'(x) = \frac{dm(x)}{dx}. \quad (1.3)$$

The quadratic velocity term works as an internal force introduced by the PDM-particle itself. It has a pronounced effect on the dynamics of the PDM system. Moreover, since $\dot{x}^2 \geq 0$, the system accelerates or decelerates for $m'(x)$ negative or positive, respectively. A particle's mass being a coordinate-dependent affects the net force \mathbf{F}_{net} , which is the sum of the external force \mathbf{F} and an internal force term $(-m'(x)\dot{x}^2/2)$.

In quantum mechanics, however, the one-dimensional kinetic energy operator is replaced by the so-called von-Roos PDM kinetic energy operator [5] :

$$\hat{T} = -\frac{\hbar^2}{4}[m(x)^f \partial_x m(x)^g \partial_x m(x)^h + m(x)^h \partial_x m(x)^g \partial_x m(x)^f] \quad (1.4)$$

where $(\partial_x = \frac{\partial}{\partial x})$ and x is the spatial coordinate. Position-dependent mass, in quantum mechanics, has significant consequences across a broad spectrum of physical systems. For instance, in semiconductor physics, variations in the effective mass of charge carriers can occur spatially, often due to impurities or heterostructures, resulting in intricate quantum transport phenomena. Similarly, in the realm of atomic and molecular physics, fluctuations in mass within distinct regions of molecules or nanoscale structures can markedly influence their electronic and vibrational characteristics. Moreover, position-dependent mass is essential for investigation unusual phenomena such as localization effects in disordered media, quantum tunneling in spatially variable potentials, and the formation of topological phases in manufactured materials [4–7]. Physicists can learn more about the behavior of quantum systems and open up new possibilities for scientific and technological advancement by including position-dependent mass in theoretical models and numerical simulations.

The Schrödinger equation with position-dependent mass, introduced by von Roos [5],

offers a theoretical basis for characterizing quantum systems where the particle's mass changes depending on its position.

The PDM von-Roos kinetic energy operator [5]; the PDM can rewrite $m(x) = m_0 f(x)$, where m_0 is the usual constant mass of the Schrödinger particle. The ordering ambiguity parameters f, g , and h align with the von-Roos restriction that $f + g + h = -1$ [5], yet the continuity conditions at heterojunctions imply that $f = h$. In this regard, Mustafa-Mazharimousavi's ordering [7–9] (i.e., $f = -1/4 = h$ and $g = -1/2$) satisfies such conditions.

In the current thesis proposal, we shall present a transformation in the radial coordinate of the global monopole spacetime and prove that the associated Schrödinger equation becomes a one-dimensional PDM von Roos Schrödinger equation. In so doing, we are motivated by Tymchyshyn's and Khlevniuk [10] observation that a point mass in motion within the curved space transforms to a PDM in Euclidean Space. By employing Mustafa Mazharimousavi's ordering, the PDM kinetic energy operator collapses into [7]:

$$\hat{T}_{MM} = -\frac{\hbar^2}{2} [m(x)^{-\frac{1}{4}} \partial_x m(x)^{-\frac{1}{2}} \partial_x m(x)^{-\frac{1}{4}}] \quad (1.5)$$

The organization of this thesis as follows: In Chapter 2, we introduce global monopole spacetime and demonstrate that a deformation or transformation in the radial coordinate of the global monopole spacetime would produce an efficacious PDM Schrödinger equation. Specifically, we show that a point canonical transformation of the radial coordinate of the PGM yields what is called in the literature Mustafa-Mazharimousavi's kinetic energy operator \hat{T}_{MM} (1.5). We begin with the Schrödinger equation in the background of transformed global monopole

spacetime. We then combine our results with the PDM von Roos Schrödinger equation to obtain Mustafa-Mazharimousavi's kinetic energy operator (1.5). Next, in Chapter 3, we consider the PDM Schrödinger equation with a Kratzer potential in global monopole spacetime. in Chapter 4, We discuss the consequences of such a PDM Schrödinger equation (with \hat{T}_{MM}) using the Kratzer potential in global monopole spacetime and a Wu-Yang magnetic monopole [12]. Our conclusion is given in Chapter 5.

Chapter 2

SCHRÖDINGER PARTICLES IN A PGM-SPACETIME

BACKGROUND

2.1 Global Monopole Spacetime

One of the intriguing phenomena that plays a significant role in current physics is the global monopole, which arises from the breaking of global $O(3)$ symmetry [20]. Other examples of such phenomena include cosmic strings and domain walls [4, 13–15]. Generally, a global monopole is a topological defect that can form due to the phase changes in the early universe [20]. This topic has been the subject of several notable investigations. Harari and Loust'o discovered that the gravitational potential of topological defects is negative [21].

Einstein's equations can be exactly solved to obtain the PGM spacetime. The metric for PGM spacetime emerges as a fascinating synthesis of gravitational curvature and topological properties [22], capturing the subtle interplay between geometry and topology on cosmic scales. The metric for global monopole spacetime is characterized by features that reflect the gravitational and topological effects caused by these cosmic relics. The presence of a global monopole, in particular, induces anisotropic curved and non-trivial topological structures, which manifest as deviations from the typical flat Minkowski spacetime metric [23].

The metric characterizing global monopole spacetime can be represented as:

$$ds^2 = g_{ab}dx^a dx^b = -dt^2 + B(r)dr^2 + r^2 d\Omega^2 \quad (2.1)$$

where g_{ab} is the metric tensor, r is the radial separation from the PGM's center, and $d\Omega^2$ represents the solid angle. The function $B(r)$ represents the gravitational and topological effects caused by the global monopole, showing the divergence from spherical symmetry and the presence of a deficit angle, which is a characteristic of topological defects. Moreover, the function $B(r)$ exhibits different behaviors in the vicinity of and away from the global monopole center, representing the transition from asymptotically flat spacetime to localized gravitational effects. The metric shows a notable deviation from the Minkowski metric [20, 24] in the central region, indicating robust topological and gravitational effects due to the global monopole. As one moves away from the center, the metric approximates the asymptotically flat form typical of locations far from large objects [25, 26].

Vilenkin and Barriola [20] have found that:

$$B(r) = \frac{1}{1 - 8\pi G\mu^2} \quad , \quad (2.2)$$

where G is the Newton's gravitational constant and μ is a constant represents the energy scale [24, 25]. We can rewrite the metric as:

$$ds^2 = -dt^2 + \frac{1}{\alpha^2} dr^2 + r^2 (d\theta^2 + \sin^2\theta d\phi^2) \quad (2.3)$$

Here $1 \geq \alpha^2 = 1 - 8\pi G\mu^2 > 0$, α is called the global monopole parameter [24].

Then, the metric tensor g_{ab} can be written as

$$g_{ab} = \begin{pmatrix} \frac{1}{\alpha^2} & 0 & 0 \\ 0 & r^2 & 0 \\ 0 & 0 & r^2 \sin^2\theta \end{pmatrix} ; \quad a, b = r, \theta, \phi,$$

and it's inverse is

$$g^{ab} = \begin{pmatrix} \alpha^2 & 0 & 0 \\ 0 & \frac{1}{r^2} & 0 \\ 0 & 0 & \frac{1}{r^2 \sin^2 \theta} \end{pmatrix}$$

Essentially, the metric characterizing global monopole spacetime captures the complex interaction between topological defects and gravitational curvature, providing insights into the significant impacts of global monopoles on the evolution of the universe [21, 23]. We can better comprehend the gravitational and topological features that constitute the cosmic environment by clarifying the mathematical formalism and physical consequences associated with this metric [27].

2.2 Schrödinger Particle in a Deformed or Transformed PGM Spacetime

Let's consider Schrödinger particles interact with a PGM spacetime metric as in (2.3) and subjected to a point canonical transformation (PCT) in the form of:

$$r = \int \sqrt{f(\rho)} d\rho = \sqrt{q(\rho)} \rho \quad (2.4)$$

where $f(\rho)$ and $q(\rho)$ are two non-zero positive valued functions that are correlated through

$$\sqrt{f(\rho)} = \sqrt{q(\rho)} \left[1 + \frac{q'(\rho)}{2q(\rho)} \rho \right].$$

To show this correlation, we just take the derivative for both sides of the integral that is given in (2.4). According to this transformation, the PGM metric becomes:

$$ds^2 = -dt^2 + \frac{f(\rho)}{\alpha^2} d\rho^2 + q(\rho) \rho^2 (d\theta^2 + \sin^2 \theta d\phi^2) \quad (2.5)$$

Then, the transformed metric tensor of the global monopole spacetime takes the form:

$$g_{ab} = \begin{pmatrix} \frac{f(\rho)}{\alpha^2} & 0 & 0 \\ 0 & \rho^2 q(\rho) & 0 \\ 0 & 0 & q(\rho) \rho^2 \sin^2 \theta \end{pmatrix}; \quad a, b = \rho, \theta, \phi \quad (2.6)$$

The determinant of which is:

$$g = |g_{ab}| = \frac{f(\rho)}{\alpha^2} q(\rho)^2 \rho^4 \sin^2 \theta$$

The inverse metric tensor is :

$$g^{ab} = \begin{pmatrix} \frac{\alpha^2}{f(\rho)} & 0 & 0 \\ 0 & \frac{1}{\rho^2 q(\rho)} & 0 \\ 0 & 0 & \frac{1}{q(\rho) \rho^2 \sin^2 \theta} \end{pmatrix} \quad (2.7)$$

The general form of the Schrödinger equation in any spacetime can be written as:

$$\left[\left(-\frac{\hbar^2}{2m_0} \frac{1}{\sqrt{g}} \partial_a \sqrt{g} g^{ab} \partial_b \right) + V(\rho) \right] \Psi(\rho, \theta, \phi, t) = i \hbar \partial_t \Psi(\rho, \theta, \phi, t) \quad (2.8)$$

Here, $V(\rho)$ is the potential energy, and $\Psi(\rho, \theta, \phi, t)$ is the wave function of the particle, which depends on the spacetime coordinates. Equation (2.8) can be extended to give:

$$\left[-\frac{\hbar^2}{2m_0} \frac{1}{\sqrt{g}} \left(\partial_\rho \sqrt{g} g^{\rho\rho} \partial_\rho + \partial_\theta \sqrt{g} g^{\theta\theta} \partial_\theta + \partial_\phi \sqrt{g} g^{\phi\phi} \partial_\phi \right) + V(\rho) \right] \Psi = i \hbar \partial_t \Psi.$$

By substituting the elements of the metric tensor of the transformed PGM spacetime in the Schrödinger equation above we get

$$\left[-\frac{\hbar^2}{2m_0} \left(\frac{\alpha^2}{q(\rho) \rho^2 \sqrt{f(\rho)}} \partial_\rho \frac{\rho^2 q(\rho)}{\sqrt{f(\rho)}} \partial_\rho + \frac{1}{\sin \theta \rho^2 q(\rho)} \partial_\theta \sin \theta \partial_\theta \right. \right. \\ \left. \left. + \frac{1}{q(\rho) \rho^2 \sin^2 \theta} \partial_\phi^2 \right) + V(\rho) \right] \Psi = i \hbar \partial_t \Psi$$

Using separation of variable, we use $\Psi(\rho, \theta, \phi, t) = \psi(\rho, \theta, \phi)f(t)$ to get:

$$\left[-\frac{\hbar^2}{2m_0} \left(\frac{\alpha^2}{q(\rho)\rho^2\sqrt{f(\rho)}} \partial_\rho \frac{\rho^2 q(\rho)}{\sqrt{f(\rho)}} \partial_\rho \psi + \frac{1}{\sin\theta \rho^2 q(\rho)} \partial_\theta \sin\theta \partial_\theta \psi + \frac{1}{q(\rho)\rho^2 \sin^2\theta} \partial_\phi^2 \psi \right) + V(\rho) \psi \right] \frac{1}{\psi} = i\hbar \frac{1}{f(t)} f'(t)$$

Note that the R.H.S. of this equation is a function of t only, and the L.H.S. is a function of ρ , θ and ϕ only. Thus, the two sides must be equal to the same constant, which we will call E . Then we have:

$$f'(t) = \frac{-iE}{\hbar} f(t) \Rightarrow f(t) = e^{-\frac{iEt}{\hbar}} \quad (2.9)$$

and

$$\left[-\frac{\hbar^2}{2m_0} \left(\frac{\alpha^2}{q(\rho)\rho^2\sqrt{f(\rho)}} \partial_\rho \frac{\rho^2 q(\rho)}{\sqrt{f(\rho)}} \partial_\rho \psi + \frac{1}{\sin\theta \rho^2 q(\rho)} \partial_\theta \sin\theta \partial_\theta \psi + \frac{1}{q(\rho)\rho^2 \sin^2\theta} \partial_\phi^2 \psi \right) + V(\rho) \psi \right] = E \psi. \quad (2.10)$$

Next, for equation (2.10) we assume that $\psi(\rho, \theta, \phi) = \psi(\rho)Y(\theta, \phi)$, to obtain:

$$\left[-\frac{\hbar^2}{2m_0} \left(\frac{\alpha^2}{\rho^2 q(\rho)\sqrt{f(\rho)}\psi(\rho)} \partial_\rho \frac{\rho^2 q(\rho)}{\sqrt{f(\rho)}} \partial_\rho \psi(\rho) + \frac{1}{\sin\theta \rho^2 q(\rho)Y} \partial_\theta \sin\theta \partial_\theta Y + \frac{1}{q(\rho)\rho^2 \sin^2\theta Y} \partial_\phi^2 Y \right) + V(\rho) \right] = E \quad (2.11)$$

Multiplying by $\frac{-2m_0 q(\rho)\rho^2}{\hbar^2}$, from the left, we get:

$$\frac{1}{\psi(\rho)} \left[\frac{\alpha^2}{\sqrt{f(\rho)}} \partial_\rho \frac{\rho^2 q(\rho)}{\sqrt{f(\rho)}} \partial_\rho \psi(\rho) - \frac{2m_0 q(\rho)\rho^2}{\hbar^2} (V(\rho) - E) \right] + \frac{1}{Y} \left[\frac{1}{\sin\theta} \partial_\theta \sin\theta \partial_\theta Y + \frac{1}{\sin^2\theta} \partial_\phi^2 Y \right] = 0$$

The term in the first square brackets depends on ρ only. but the remainder depends on θ and ϕ only. Thus, each must be constant. we will write the separation constant in the form $l(l+1)$, so we have:

$$\frac{1}{Y} \left[\frac{1}{\sin\theta} \partial_\theta \sin\theta \partial_\theta Y + \frac{1}{\sin^2\theta} \partial_\phi^2 Y \right] = -l(l+1) \quad (2.12)$$

and

$$\frac{1}{\psi(\rho)} \left[\frac{\alpha^2}{\sqrt{f(\rho)}} \partial_\rho \frac{\rho^2 q(\rho)}{\sqrt{f(\rho)}} \partial_\rho \psi(\rho) - \frac{2m_0 q(\rho) \rho^2}{\hbar^2} (V(\rho) - E) \right] = l(l+1) \quad (2.13)$$

Equation (2.12) represents the dependence of $\Psi(r, \theta, \phi)$ on θ and ϕ , and multiplying it from the left by $-\hbar^2 Y$, it becomes:

$$-\hbar^2 \left[\frac{1}{\sin\theta} \partial_\theta \sin\theta \partial_\theta Y + \frac{1}{\sin^2\theta} \partial_\phi^2 Y \right] = \hbar^2 l(l+1) Y \quad (2.14)$$

The given equation is the angular momentum equation, which may be expressed as $L^2 Y_{lm}(\theta, \phi) = \hbar^2 l(l+1) Y_{lm}(\theta, \phi)$. Where \hat{L} is the angular momentum operator in spherical coordinates. The spherical harmonics, denoted as $Y_{lm}(\theta, \phi)$, represent a set of functions used in the study of spherical symmetry. The angular momentum quantum number is represented by l .

According to our PCT (2.4) we have $\partial_\rho = \left(\frac{dr}{d\rho} \right) \partial_r = \left(\sqrt{f(\rho)} \right) \partial_r$, and $r^2 = q(\rho) \rho^2$, so equation (2.13) in terms of r is now given by:

$$\left[\alpha^2 \partial_r r^2 \partial_r \psi(r(\rho)) - \frac{2m_0 r^2}{\hbar^2} (V(r(\rho)) - E) \psi(r(\rho)) \right] = l(l+1) \psi(r(\rho))$$

If we multiply by $\frac{-\hbar^2}{2m_0 r^2 \alpha^2}$, we obtain

$$\left[-\frac{\hbar^2}{2m_0} \frac{1}{r^2} \partial_r r^2 \partial_r \psi(r) + \frac{1}{\alpha^2} (V(r) - E) \psi(r) \right] = -\frac{\hbar^2}{2m_0} \frac{l(l+1)}{\alpha^2 r^2} \psi(r)$$

This can be written in the form:

$$\frac{\hbar^2}{2m_0} \left(-\frac{1}{r^2} \partial_r r^2 \partial_r \psi + \frac{l(l+1)}{\alpha^2 r^2} \psi \right) + \frac{1}{\alpha^2} V(r) \psi = \frac{1}{\alpha^2} E \psi$$

Now, by substituting $\frac{E}{\alpha^2} = \mathcal{E}$ and $\frac{l(l+1)}{\alpha^2} = \beta(\beta+1)$, last equation becomes:

$$\frac{\hbar^2}{2m_0} \left(-\frac{1}{r^2} \partial_r r^2 \partial_r \psi + \frac{\beta(\beta+1)}{r^2} \psi \right) + \frac{1}{\alpha^2} V(r) \psi = \mathcal{E} \psi \quad (2.15)$$

This can be simplified using $\psi(r) = \frac{U(r)}{r}$ to obtain :

$$\frac{\hbar^2}{2m_0} \left(-\partial_r^2 U(r) + \frac{\beta(\beta+1)}{r^2} U(r) \right) + \frac{1}{\alpha^2} V(r) U(r) = \mathcal{E} U(r) \quad (2.16)$$

To express β in terms of l and α , we use the definition $\frac{l(l+1)}{\alpha^2} = \beta(\beta+1)$

$$\beta = -\frac{1}{2} + \frac{\sqrt{\alpha^2 + 4l(l+1)}}{2\alpha}$$

This equation calculates the value of the regular angular momentum quantum number l for flat Minkowski spacetime when $\alpha = 1$. Furthermore, the two quantum mechanical systems described in equations (2.13) and (2.15) have the same spectrum and remain unchanged. In other words, if you know the solution to one of them, you can quickly determine the solution to the other. However, they both possess identical energy.

It should be noted that the angular component of the wave function $Y(\theta, \phi)$, remains constant for all spherically symmetric potentials. The specific form of the potential, $V(r)$, only influences the radial component of the wave function, $\psi(\rho)$, which is determined by Equation 2.15, Thus $\psi(r, \theta, \phi, t) = \psi(r) Y_{lm}(\theta, \phi) e^{-i Et/\hbar}$.

2.3 Position-Dependent Mass in the Radially Transformed PGM Spacetime Metric

When dealing with situations in which the particle's mass changes depending on its position, it is necessary to adjust the kinetic energy operator in (1.4) to accommodate this variability. In Chapter 1, it has been shown that by choosing MM-ordering, $f = h = -\frac{1}{4}$ and $g = -\frac{1}{2}$, the PDM von-Roos kinetic energy operator takes the form of Mustafa-Mazharimousavi's PDM one, \hat{T}_{MM} given in (1.5) [8, 9].

We begin with recalling the radial equation in (2.16):

$$\frac{\hbar^2}{2m_0} \left(-\partial_r^2 U(r) + \frac{\beta(\beta+1)}{r^2} U(r) \right) + \frac{1}{\alpha^2} V(r) U(r) = \mathcal{E} U(r)$$

and substituting:

$$U(r) = U(r(\rho)) = f(\rho)^{-\frac{1}{4}} \phi(\rho) \quad (2.17)$$

Moreover, our PCT $r = \int \sqrt{f(\rho)} d\rho$ allows us to write $\partial_r U(r)$ as

$$\partial_r U(r) = \left(\frac{d\rho}{dr} \right) \partial_\rho U = f(\rho)^{-\frac{1}{2}} \partial_\rho \left(f(\rho)^{-\frac{1}{4}} \phi(\rho) \right)$$

and

$$\begin{aligned} \partial_r^2 U(r) &= \partial_r (\partial_r U) \\ &= \partial_r \left[f(\rho)^{-\frac{1}{2}} \partial_\rho \left(f(\rho)^{-\frac{1}{4}} \phi(\rho) \right) \right] \\ &= \frac{d\rho}{dr} \partial_\rho \left[f(\rho)^{-\frac{1}{2}} \partial_\rho \left(f(\rho)^{-\frac{1}{4}} \phi(\rho) \right) \right] \\ &= f(\rho)^{-\frac{1}{2}} \left(\partial_\rho \left[f(\rho)^{-\frac{1}{2}} \partial_\rho \left(f(\rho)^{-\frac{1}{4}} \phi(\rho) \right) \right] \right), \end{aligned}$$

to obtain

$$\begin{aligned} \left(-\frac{\hbar^2}{2m_0} f(\rho)^{-\frac{1}{2}} \partial_\rho f(\rho)^{-\frac{1}{2}} \partial_\rho + \frac{\hbar^2}{2m_0} \frac{\beta(\beta+1)}{q(\rho)\rho^2} + \frac{1}{\alpha^2} V(r) \right) f(\rho)^{-\frac{1}{4}} \phi(\rho) \\ = \mathcal{E} f(\rho)^{-\frac{1}{4}} \phi(\rho) \end{aligned} \quad (2.18)$$

Next, we multiply this equation, from the left, by $f(\rho)^{\frac{1}{4}}$ and get:

$$\left(-\frac{\hbar^2}{2m_0} f(\rho)^{-\frac{1}{4}} \partial_\rho f(\rho)^{-\frac{1}{2}} \partial_\rho f(\rho)^{-\frac{1}{4}} + V_{eff}(\rho) \right) \phi(\rho) = \mathcal{E} \phi(\rho) \quad (2.19)$$

Where the effective potential V_{eff} is now given by:

$$V_{eff} = \frac{\hbar^2}{2m_0} \frac{\beta(\beta+1)}{q(\rho)\rho^2} + \frac{1}{\alpha^2} V(\rho)$$

that includes the central repulsive term. The particle is often pushed far from the origin, similar to the repulsive force seen in classical mechanics. Thus, the resulting operator for the effective PDM kinetic energy is:

$$\hat{T}_{MM} = -\frac{\hbar^2}{2m_0} f(\rho)^{-\frac{1}{4}} \partial_\rho f(\rho)^{-\frac{1}{2}} \partial_\rho f(\rho)^{-\frac{1}{4}} \quad (2.20)$$

Which is known as Mustafa-Mazharimousavi's (MM) PDM-kinetic-energy operator [8] Then, the two systems, identified in (2.16) and (2.19), exhibit

isospectrality, meaning they possess identical energy levels and remain unchanged under certain transformations.

Chapter 3

PDM SCHRÖDINGER PARTICLE IN A KRATZER POTENTIAL IN A PGM SPACETIME

3.1 Kratzer Potential

Molecular potentials are used to describe the interactions between atoms or molecules. They provide a foundation for studying the behaviour and properties of matter on a molecular scale. They are essential in the field of physics, allowing us to explore the intricacies of molecular behaviour and acquire crucial knowledge about the fundamental forces that control matter. Some examples of various molecular potentials and their usage also include Yukawa potential [35–38, 40], Morse potential [44–49], Hulthén potential [43, 50–54], and Pöschl-Teller potential [55–60]. The Kratzer potential [41], which we are going to study, demonstrates a distinctive property of tending towards infinity as the distance between nuclei collapses to zero [61, 62]. This behaviour is a result of the repulsive interactions that exist between the individual molecules within the potential.

The Kratzer potential is defined as [41, 63]:

$$V_K(r) = -2D \left(\frac{A}{r} - \frac{A^2}{2r^2} \right) \quad (3.1)$$

where D is the energy required to separate two atoms in a diatomic molecule, when they are at an equilibrium inter-nuclear distance A . Both A and D are positive values. The potential under consideration is a combination of the attractive Coulomb-like term potential at long distances and a repulsive centrifugal like barrier (the second term)

at small distances. Their combination of which introduces an effective potential well, which functions as a model for studying the molecular structure, molecular energy spectra, chemical interactions, and inter-nuclear vibration of diatomic molecules.

3.2 PDM Schrödinger Particle with Kratzer Potential and in a PGM Spacetime Background

We have shown, in the previous chapter, that the radial the Schrödinger equation takes the form of (2.16). If we substitute the Kratzer potential in (2.16) we obtain:

$$\frac{\hbar^2}{2m_0} \left(-\partial_r^2 U(r) + \frac{\beta(\beta+1)}{r^2} U(r) \right) - \frac{2D}{\alpha^2} \left(\frac{A}{r} - \frac{A^2}{2r^2} \right) U(r) = \mathcal{E} U(r) \quad (3.2)$$

We aim to find the solution for $U(r)$, and determine the allowed energies. To simplify this equation we will use $\hbar = 2m_0 = 1$ it becomes:

$$\left[-\partial_r^2 + \frac{\beta(\beta+1)}{r^2} - \frac{2D}{\alpha^2} \left(\frac{A}{r} - \frac{A^2}{2r^2} \right) \right] U(r) = \mathcal{E} U(r) \quad (3.3)$$

We can rewrite it as:

$$\left(-\partial_r^2 + \frac{L(L+1)}{r^2} - \frac{\omega}{r} \right) U(r) = \mathcal{E} U(r) \quad (3.4)$$

where $\omega = 2DA/\alpha^2$ and

$$L(L+1) = \beta(\beta+1) + \frac{DA^2}{\alpha^2} = \frac{l(l+1) + DA^2}{\alpha^2}.$$

Thus:

$$L = -\frac{1}{2} + \sqrt{\frac{1}{4} + \frac{l(l+1) + DA^2}{\alpha^2}}$$

Next, we define a new variable $z = \sqrt{-\mathcal{E}}r$ so that:

$$\left(\partial_z^2 - \frac{L(L+1)}{z^2} - \frac{\tilde{\omega}}{z} - 1 \right) U = 0 \quad (3.5)$$

where $\tilde{\omega} = \frac{\omega}{\sqrt{-\mathcal{E}}}$.

Next, we will study the asymptotic behavior. As z goes to infinity the constant term dominates, so:

$$\partial_z^2 U = U \quad (3.6)$$

The general solution is: $U = A_1 e^{-z} + B_1 e^z$ but e^z is unbounded as z goes to infinity, so B must be zero. Thus, for large z we have:

$$U = A_1 e^{-z}$$

On the other hand, as z goes to zero the second term in equation 3.5 dominates, then:

$$\partial_z^2 U - \frac{L(L+1)}{z^2} U = 0 \quad (3.7)$$

The general solution is

$$U = C z^{L+1} + D z^{-L}$$

But U blows up as z goes to zero, so $D = 0$, Thus, for small z :

$$U = C z^{L+1}$$

Finally, we can write $U(z)$ as:

$$U(z) = z^{L+1} e^{-z} F(z)$$

In the hope that the new function $F(z)$ will simplify the equation.

$$\partial_z U = z^L e^{-z} [(L+1-z) F + z \partial_z F]$$

and

$$\partial_z^2 U = z^L e^{-z} \left[\left(-2L - 2 + z + \frac{L(L+1)}{z} \right) F + 2(L+1-z) \partial_z F + z \partial_z^2 F \right]$$

Then the radial equation in terms of $F(z)$ becomes:

$$z \partial_z^2 F + 2(L+1-z) \partial_z F + [\tilde{\omega} - 2(L+1)] F = 0 \quad (3.8)$$

Equation (3.8) is a confluent hypergeometric differential equation with the confluent hypergeometric function/series solution ${}_1F_1(a, b, x)$:

$$F(z) = C_{nl} {}_1F_1 \left(L - \frac{\tilde{\omega}}{2} + 1, 2 + 2L, 2z \right) \quad (3.9)$$

If this solution is to be a physically acceptable one, it should be finite and square integrable. This would once again necessitate that the confluent hypergeometric series be terminated to a polynomial of degree $n = 0, 1, 2, 3, \dots$ so that:

$$L - \frac{\tilde{\omega}}{2} + 1 = -n \quad (3.10)$$

We can easily solve it to get the value of \mathcal{E}

$$\mathcal{E} = -\left(\frac{\omega}{\tilde{\omega}}\right)^2 = -\frac{\omega^2}{4(1+L+n)^2} \quad (3.11)$$

which implies the energy levels:

$$E_{n,l} = -\frac{\omega^2 \alpha^2}{4(1+L+n)^2} \quad (3.12)$$

Then the solution is written as:

$$\Psi_{nl}(z) = C_{nl} z^{L+1} e^{-z} {}_1F_1\left(L - \frac{\tilde{\omega}}{2} + 1, 2 + 2L, 2z\right) \quad (3.13)$$

Obviously, equation (3.12) suggests that the energies tend to converge to $E_{n,l} \sim -D\alpha^2$ as the coupling constant $A \rightarrow \infty$. This would indicate that all energy levels cluster at the value $E_{n,l} \sim -D\alpha^2$ as $A \rightarrow \infty$. Such a tendency is observed in Figures 3.1(a) and 3.1(b), where the energies $E_{n,l} \sim -1/4$ for the parametric values used therein.

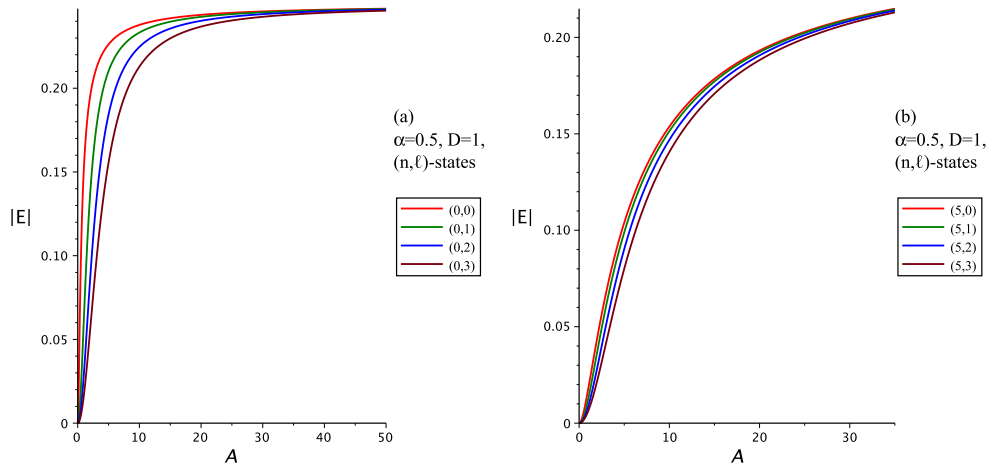


Figure 3.1: The energy levels for (n, l) -states against the coupling constant A for $l = 0, 1, 2, 3$ so that Fig. 3.1(a) for $n = 0$, and Fig. 3.1(b) for $n=5$.

In Figure 3.2, we show the energies $|E_{n,l}|$ against different values of the monopole parameter and within the range $0 < \alpha \leq 1$. The comparison between Figures 3.2(a), (b), and (c) suggests that the energies $|E_{n,l}|$ decrease as the radial quantum number increases. Indicating a general tendency of the energy levels towards the value $E_{n,l} \rightarrow -1/n^2$ for $n \gg 1$. This trend can be clearly observed in Fig. 3.2 (d). Moreover, the spacing between the energy levels narrows as the radial quantum number increases.

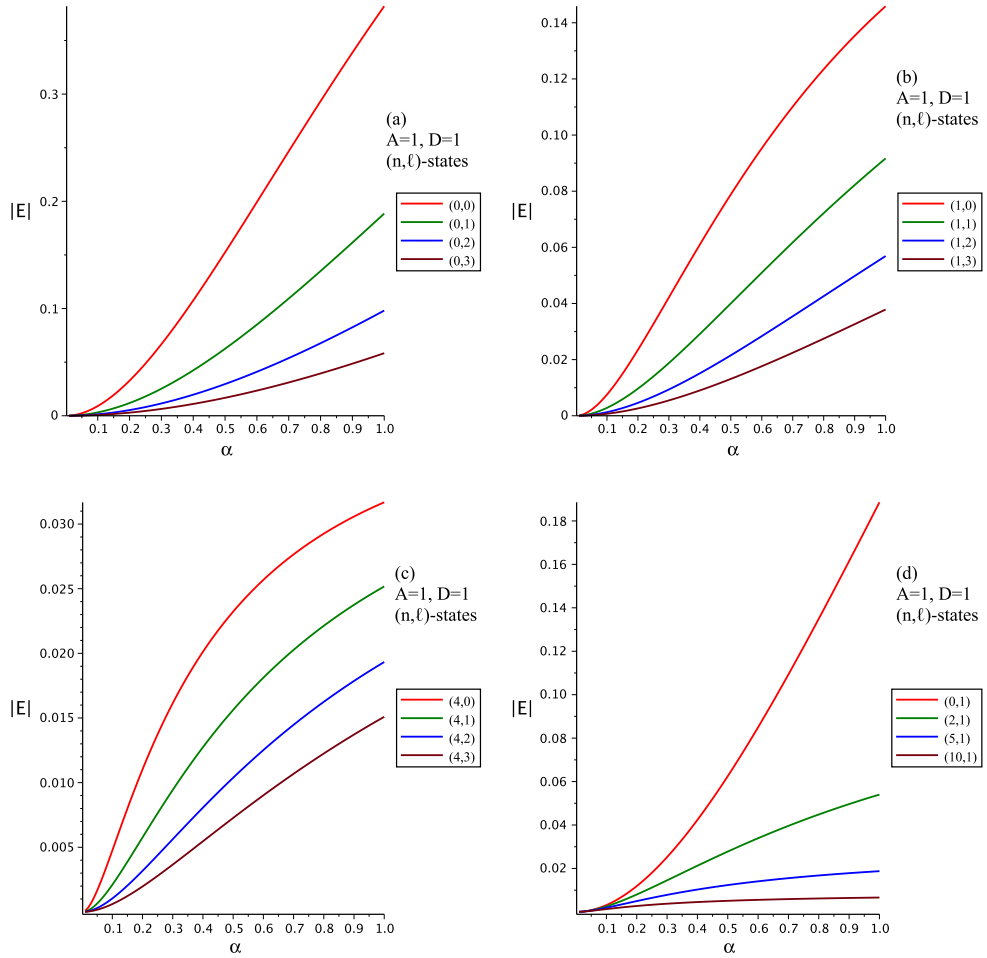


Figure 3.2: The energy levels for (n, l) -states against the monopole parameter α for $l = 0, 1, 2, 3$ so that Fig. 3.2(a) for $n = 0$, Fig. 3.2(b) for $n=1$, Fig. 3.2(c) for $n=4$, and Fig. 3.2(d) for $l = 1$ and $n = 0, 2, 5, 10$.

Chapter 4

PDM SCHRÖDINGER PARTICLES WITH KRATZER POTENTIAL IN A PGM SPACETIME AND A WU-YANG MAGNETIC MONOPOLE

4.1 Wu- Yang Magnetic Monopole

The choice of the vector potential $\mathbf{A} = (0, 0, A_\phi)$ around a global monopole necessarily involves singularities. This situation is analogous to selecting a coordinate system on the surface of a sphere, such as longitude and latitude. Even though the sphere itself is free of intrinsic singularities, any coordinate system will have some singularities. To avoid introducing singularities, one can divide the sphere into overlapping regions and define a coordinate system without singularities in each region [64–66].

Following this strategy, Wu and Yang divided the space outside a magnetic monopole into two regions, R_A and R_B . They defined a vector potential A_A in R_A and a vector potential A_B in R_B , covering the entire space exterior to the magnetic monopole. These regions overlap in R_{AB} with the monopole at the origin as they have chosen [64]

$$\begin{aligned} R_A : 0 \leq \theta < \frac{\pi + 2\eta}{2} \quad , r > 0, \quad 0 \leq \phi < 2\pi \\ R_B : \frac{\pi - 2\eta}{2} < \theta \leq \pi \quad , r > 0, \quad 0 \leq \phi < 2\pi \\ R_{AB} : \frac{\pi - 2\eta}{2} < \theta \leq \frac{\pi + 2\eta}{2} \quad , r > 0, \quad 0 \leq \phi < 2\pi \end{aligned}$$

Where $0 < \eta \leq \frac{\pi}{2}$. The vector potential components are chosen to be:

$$A_{r,A} = A_{\theta,A} = 0, \quad A_{\phi,A} = g(1 - \cos\theta) \quad (4.1)$$

$$A_{r,B} = A_{\theta,B} = 0, \quad A_{\phi,B} = -g(1 + \cos\theta) \quad (4.2)$$

where g is the strength of the Wu-Yang magnetic monopole. To simplify the notation and make it quite handy, we use the form $A_\phi = sg - g \cos\theta$, such that $s = 1$ for $A_{\phi,A}$ and $s = -1$ for $A_{\phi,B}$.

4.2 PDM Schrödinger Equation with Kratzer Potential and Wu-Yang Magnetic Monopole

The Schrödinger equation with potential $V(r)$, in a vector potential $\mathbf{A} = (0, 0, A_\phi)$ is given by:

$$\left[\left(-\frac{1}{\sqrt{g}} (\partial_a - ieA_a) \sqrt{g} g^{ab} (\partial_b - ieA_b) \right) + V(r) \right] \Psi(r, \theta, \phi, t) = i \partial_t \Psi(r, \theta, \phi, t) \quad (4.3)$$

Equation (4.3) can be simplified :

$$\left[-\frac{1}{\sqrt{g}} \left((\partial_r - ieA_r) \sqrt{g} g^{rr} (\partial_r - ieA_r) + (\partial_\theta - ieA_\theta) \sqrt{g} g^{\theta\theta} (\partial_\theta - ieA_\theta) \right. \right. \\ \left. \left. + (\partial_\phi - ieA_\phi) \sqrt{g} g^{\phi\phi} (\partial_\phi - ieA_\phi) \right) + V(r) \right] \Psi(r, \theta, \phi, t) = i \partial_t \Psi(r, \theta, \phi, t)$$

By substituting the elements of the metric tensor of global monopole spacetime and the Wu-Yang vector potential the Schrödinger equation becomes:

$$\left[-\frac{\alpha^2}{r^2} \partial_r (r^2 \partial_r) - \frac{1}{r^2} \left(\frac{1}{\sin\theta} \partial_\theta \sin\theta \partial_\theta + \frac{1}{\sin^2\theta} [\partial_\phi - ieA_\phi]^2 \right) \right. \\ \left. + V(r) \right] \Psi(r, \theta, \phi, t) = i \partial_t \Psi(r, \theta, \phi, t) \quad (4.4)$$

Where r is given by (2.4). We will use the separation of variables and assume that $\Psi(r, \theta, \phi, t) = \psi(r) Y_{\tilde{q}lm}(\theta, \phi) e^{-iEt}$, where $q = eg \geq 0$ and $\tilde{q} = sq$. Then, the time-independent Schrödinger equation becomes:

$$\frac{1}{\psi(r)} \left[-\alpha^2 \partial_r (r^2 \partial_r) + r^2 (V(r) - E) \right] \psi(r) \\ - \frac{1}{Y_{\tilde{q}lm}(\theta, \phi)} \left(\frac{1}{\sin\theta} \partial_\theta \sin\theta \partial_\theta + \frac{1}{\sin^2\theta} [\partial_\phi - ieA_\phi]^2 \right) Y_{\tilde{q}lm}(\theta, \phi) = 0$$

The terms in the first square brackets depends only on r , whereas the remainder depends on θ and ϕ . Hence, we may take:

$$\left(\frac{1}{\sin\theta} \partial_\theta \sin\theta \partial_\theta + \frac{1}{\sin^2\theta} [\partial_\phi - i e A_\phi]^2 \right) Y_{\bar{q}lm}(\theta, \phi) = -\lambda Y_{\bar{q}lm}(\theta, \phi) \quad (4.5)$$

and

$$\left[-\frac{\alpha^2}{r^2} \partial_r (r^2 \partial_r) + \frac{\lambda}{r^2} + V(r) \right] \psi(r) = E \psi(r) \quad (4.6)$$

Now, we have to find the values of λ by using the substitution $Y_{\bar{q}lm}(\theta, \phi) = e^{i(m+\bar{q})\phi} \mathcal{G}_{qlm}(\theta)$ into Equation (4.6) to obtain:

$$\left(\frac{1}{\sin\theta} \partial_\theta \sin\theta \partial_\theta - \frac{1}{\sin^2\theta} [m + q \cos\theta]^2 \right) \mathcal{G}_{qlm}(\theta) = -\lambda \mathcal{G}_{qlm}(\theta) \quad (4.7)$$

The above equation is independent of the values of s , thus the solution for $\mathcal{G}_{qlm}(\theta)$ are the same in the two regions A and B. This means

$$[\mathcal{G}_{qlm}(\theta)]_A = \mathcal{G}_{qlm}(\theta) = [\mathcal{G}_{qlm}(\theta)]_B.$$

To solve Equation 4.7 we are going to use the substitution $x = \cos\theta$, it becomes:

$$\left([1-x^2] \partial_x^2 - 2x \partial_x - \frac{[m+qx]^2}{1-x^2} \right) \mathcal{G}_{qlm}(x) = -\lambda \mathcal{G}_{qlm}(x) \quad (4.8)$$

This differential equation has singularities at $x = 1$ and $x = -1$. We define now

$$\mathcal{G}_{qlm}(x) = (1+x)^{a/2} (1-x)^{b/2} P_{qlm}(x) \quad (4.9)$$

where $a = (|m|-q)$ and $b = (|m|+q)$.

By using the substitution (4.9) Equation (4.8) becomes:

$$(x^2 - 1) P_{qlm}''(x) + 2[q + (m+1)x] P_{qlm}'(x) + (m+m^2 - \lambda - q^2) P_{qlm} = 0 \quad (4.10)$$

The solution of the above equation is the hypergeometric functions/series:

$$P_{qlm} = C_1 F_1 \left(|m| + \frac{1}{2} \pm \frac{1}{2} \sqrt{4\lambda + 4q^2 + 1}, |m| - q + 1, \frac{1}{2}(1+x) \right) \quad (4.11)$$

For the sake of having a finite and square integrable wave functions, the power series of the hypergeometric function must be terminated, thus:

$$-\tilde{n} = |m| + \frac{1}{2} \pm \frac{1}{2} \sqrt{4q^2 + 4\lambda + 1}; \quad \tilde{n} = 0, 1, 2, 3, \dots$$

We solve this equation to get λ :

$$\lambda = (|m| + \tilde{n} + 1)(|m| + \tilde{n}) - q^2 = l(l+1) - q^2 \quad (4.12)$$

Where $l = |m| + \tilde{n} = 0, 1, 2, 3, \dots$ is the angular momentum quantum number. That is, when $q = 0$ (this means no Wu-Yang magnetic monopole) then $\lambda = l(l+1)$ represents the regular eigenvalues for the textbook spherical harmonics. Now, we can write the solution Y_{qlm} as:

$$Y_{qlm}(\theta, \phi) = \begin{cases} (1 + \cos \theta)^{\frac{a}{2}} (1 - \cos \theta)^{\frac{b}{2}} P_{qlm}(\cos \theta) e^{i(m+q)\phi} & , \text{in region A} \\ (1 + \cos \theta)^{\frac{a}{2}} (1 - \cos \theta)^{\frac{b}{2}} P_{qlm}(\cos \theta) e^{i(m-q)\phi} & , \text{in region B} \end{cases} \quad (4.13)$$

Wu and Yang have called Y_{qlm} as the monopole harmonics.

Now, we are going to write the radial equation 4.6, with the Kratzer potential $V_K(r) = -2D \left(\frac{A}{r} - \frac{A^2}{2r^2} \right)$ and $\lambda = l(l+1) - q^2$. Then the radial equation becomes:

$$\left[-\frac{\alpha^2}{r^2} \partial_r (r^2 \partial_r) + \frac{l(l+1) - q^2}{r^2} - 2D \left(\frac{A}{r} - \frac{A^2}{2r^2} \right) \right] \psi(r) = E \psi(r) \quad (4.14)$$

We may rewrite it as:

$$\left[-\frac{1}{r^2} \partial_r (r^2 \partial_r) + \frac{\zeta(\zeta+1)}{r^2} - \frac{\omega}{r} \right] \psi(r) = \mathcal{E} \psi(r) \quad (4.15)$$

where $\mathcal{E} = \frac{E}{\alpha^2}$, $\omega = \frac{2DA}{\alpha^2}$ and $\zeta(\zeta+1) = \frac{l(l+1) - q^2 + DA^2}{\alpha^2}$. If we solve it for ζ we can easily get

$$\zeta = -\frac{1}{2} + \sqrt{\frac{1}{4} + \frac{A^2 D + l(l+1) - q^2}{\alpha^2}} \quad (4.16)$$

Furthermore, the solution to Equation (4.15) would read:

$$\psi(\rho(r)) = C_{nl} r^\zeta e^{-\sqrt{-\mathcal{E}}r} {}_1F_1\left(1 + \zeta - \frac{\omega}{2\sqrt{-\mathcal{E}}}, 2 + 2\zeta, 2\sqrt{-\mathcal{E}}r\right) \quad (4.17)$$

Hereby, the condition that the wavefunction be finite and square integrable would necessitate that the confluent hypergeometric series be terminated to a polynomial of

degree $n \geq 0$ to obtain:

$$1 + \zeta - \frac{\omega}{2\sqrt{-\mathcal{E}}} = -n \quad (4.18)$$

We can easily solve it to get the value of \mathcal{E}

$$\mathcal{E} = -\frac{\omega^2}{4(1 + \zeta + n)^2} \quad (4.19)$$

Which implies the energy levels:

$$E_{n,l,q} = -\frac{\omega^2 \alpha^2}{4(1 + \zeta + n)^2} \quad (4.20)$$

At this point, one should notice that such energy levels collapse into those without the Wu-Yang magnetic monopole (i.e., $q = 0$) given in Eq.(3.12). In Figures 4.1(a) and (b), we show the energies $|E_{n,l,q}|$ at $\alpha = 0.5$ for different values of the strengths q of the Wu-Yang magnetic monopole. The values of q are chosen so that the square root in Eq. (4.16) is a real value. We observe that the Wu-Yang magnetic monopole yields non-equally spaced energy levels. Additionally, the spacing between the energy levels increases as the strength of the Wu-Yang magnetic monopole increases. Moreover, the spacing between the energy levels decreases as the radial quantum number increases. Furthermore, in Figures 4.1 (c) and (d) we plot the energy levels against the coupling parameter A for $\alpha = 0.5$ and $\alpha = 0.1$, respectively at $q = 1$. It is clear that as the value of the coupling parameter A increases, the energy values increase until reaching their maximum. This maximum energy value depends on the value of α . This tendency is evident from equation (4.20), which suggests that the energies tend to converge to $E_{n,l} \rightarrow -D\alpha^2$ as the coupling constant $A \rightarrow \infty$.

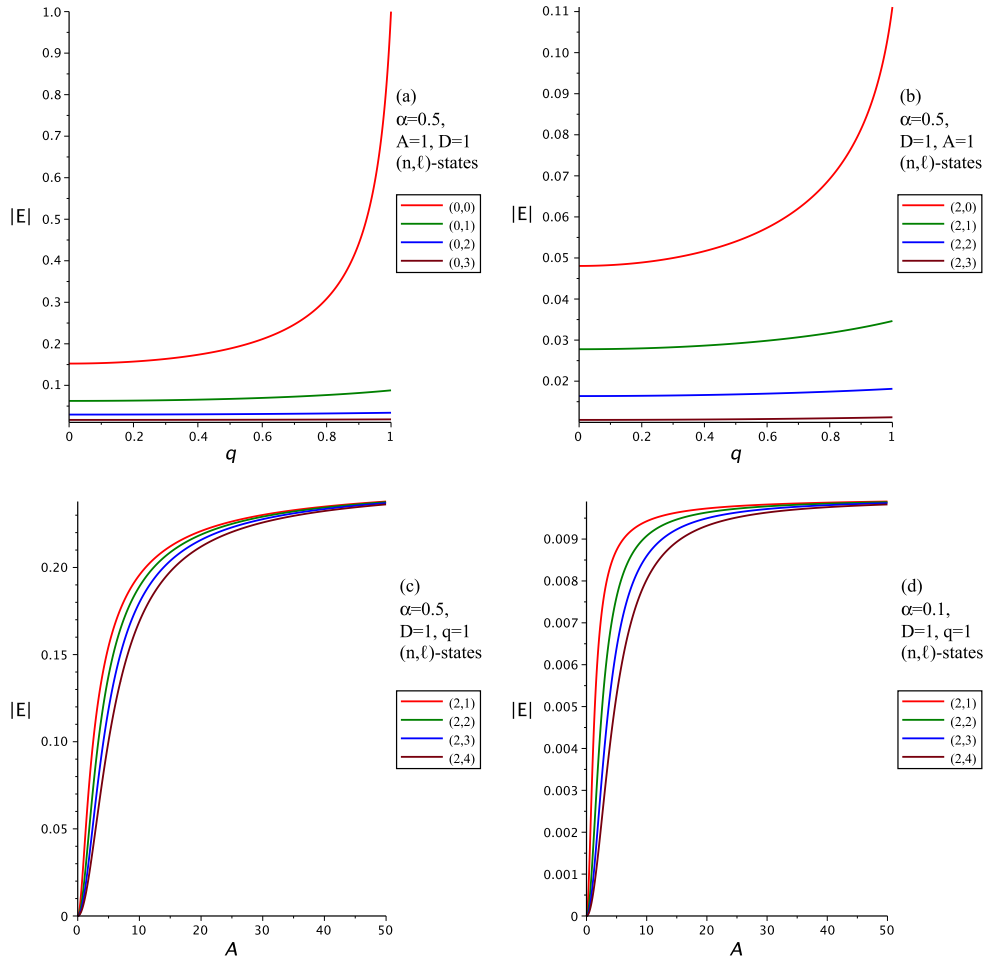


Figure 4.1: The energy levels for (n, l) -states against the WY-magnetic monopole parameter q for $D = 1 = A$ so that Fig. 4.1(a) for $n = 0$ and $\alpha = 0.5$, Fig. 4.1(b) for $n = 2$ and $\alpha = 0.5$, and the energy levels against the coupling parameter A for $q = 1 = D$ so that Fig. 4.1(c) for $n = 2$, $\alpha = 0.5$, and Fig. 4.1(d) for $n = 2$, $\alpha = 0.1$

Chapter 5

CONCLUSION

This work demonstrates that a particular transformation (2.4) of a global monopole spacetime yields a von Roos position-dependent mass Schrödinger equation (2.19). Additionally, we have solved the Schrödinger equation with Kratzer potential, a significant molecular potential describing the vibration-rotation spectra of diatomic molecules in global monopole spacetime, to obtain the energy levels spectrum. Here, the radial wave equation is derived in terms of the confluent hypergeometric function ${}_1F_1(a, b, x)$, and the angular part is expressed in terms of spherical harmonics.

Subsequently, we examined the Schrödinger equation with the Kratzer potential and Wu-Yang magnetic monopole in global monopole spacetime. Here, the radial wave function is still given in terms of the confluent hypergeometric function, but the angular part is expressed in terms of the Wu-Yang monopole harmonics.

Regarding the energy levels, when we plot the energy levels resulting from solving the radial part of the Schrödinger equation with the Kratzer potential equation (3.3) against the global monopole parameter α for all allowed values of α (i.e. $1 \geq \alpha > 0$), we observe that the energies $|E_{n,l}|$ decrease as the radial quantum number increases (documented in Figures 3.2(a), and (b)). Moreover, in Figure 3.2(d), we notice that the spacing between the energy levels narrows as the radial quantum number n increases.

Next, we plot the energy levels resulting by solving the radial part of the Schrödinger

equation with Kratzer potential and Wu-Yang magnetic (equation 4.15) against the strength q of the Wu-Yang magnetic at $\alpha = 0.5$. We observe that the energy levels are not equally spaced. Additionally, the gap between the energy levels widens as the strength of the Wu-Yang magnetic monopole increases. Furthermore, the gap between the energy levels narrows as the radial quantum number grows as shown in Figure 4.1(a) and (b).

When we plot the energies against the parameter A as shown in Figure 4.1 (c) and (d) for $\alpha = 0.5$ and $\alpha = 0.1$, respectively, at a constant q (we set $q=1$). We observe that as the coupling parameter A increases, the energies also increase until they reach their maximum value. Additionally, the highest possible value of the energy depends on the value of α . This trend is evident from equation (4.20), indicating that the energies approach convergence towards $E_{n,l} \rightarrow -D\alpha^2$ as the coupling constant $A \rightarrow \infty$.

Overall, this study improves our comprehension of the substantial impacts on the movement of charged particles subjected to topological defects and interacting via the Kratzer potential, with or without Wu-Yang magnetic monopole, specifically within the framework of a point-like global monopole.

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