

Correspondence between the Bloch's Theorem and the Oyibo Grand Unified Theorem within the Purview of Generic Torus

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Abstract We have shown in [1] that the geometric object for the geometrical and quantization foundation of the Oyibo grand unified theorem (GUT) is the torus. This torus is generic meaning that its nature depends on the periodic boundary conditions of the system and other physical constraint conditions. By observing that the Bloch's theorem also has this torus as its generic geometric object, we show that there is correspondence between the Bloch's theorem and the Oyibo GUT. The implication of this correspondence in using the Oyibo GUT to study strongly correlated systems in condensed matter physics is then discussed.

Keywords Bloch's theorem, Oyibo grand unified theorem, Group transformation, Geometric object, Torus

1. Introduction

The foundation of our highly simplified correlated variation approach (HSCVA) for investigating few body correlated systems has been based on the geometrical nature of the lattice structures of such systems [2-5]. In the pioneering paper, Chen and Mei (1989) developed the correlated variational method to study the standard Hubbard model only in 1D and 2D systems with singlet states [2]. To prevent boundary effects, their CVA was applied to finite 1D lattices sites arranged in 1D ring and 2D planar lattices arranged in a torus in 2D. In advancing the work, we have simplified the approach by first formulating the statistical equivalent of the Hubbard model in all three dimensions [4]. The remarkable correspondence between our statistical formulation and the Hubbard model makes it easy to modify the latter for systems having triplet states so that our HSCVA is applicable to the Hubbard model and its various extended versions in all three dimensions [3]. Like the CVA of Chen and Mei [2], the arrangement of the lattice sites in 1D is in a ring form and those in 2D and 3D systems are toroidal to avoid the edge effects on some sites. These arrangements are consistent with related works of arranging

lattices in tori so as to obtain lattices with appropriate periodic boundaries [6, 7].

It was therefore quite reassuring to have recently observed and shown that the generic geometric object needed for the geometrical and quantization foundation of the Oyibo grand unified theorem (GUT) formulated as the mathematical basis for the unified force field theory as envisaged by Oyibo [8-9] also known as the theory of everything (TOE) is the torus [1]. In that work, one of us has shown that a torus ring can be constructed from the Pythagoras triangle which is an invariant geometric object in nature. Then using projective geometry, he was able to construct from this Pythagoras triangle the torus invariant within the Lorentz transformation. He then observed that this invariance of the toroidal geometry of the Oyibo GUT and the invariance of the Einstein's principle of relativity within the purview of Lorentz transformation, is responsible for the remarkable correspondence of the Oyibo GUT with the Einstein's unified field equation for conformal invariant field theories. Following that line of thinking, it is postulated here that since the Bloch theorem from which the original Hubbard model was formulated [10-13] has been observed to have a toroidal lattice foundation for appropriate periodic boundaries, then it should have a correspondence with the Oyibo GUT. This is the motivation for this paper because the demonstration of such correspondence implies that the Hubbard model and its extended versions can also be formulated from the Oyibo GUT. The following plan will be used for this study. As a first step, it will be necessary to first re-visit the formation

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of the Hubbard model and its extended versions from the Bloch's theorem in Section 2. The demonstration of the correspondence of the Bloch's theorem and the Oyibo GUT will be done in Section 3. Thereafter, we will discuss the implication of this correspondence and conclude.

2. Revisiting the Formation of the Hubbard Model and Its Extended Versions

The Bloch theorem plays a central role in conduction electron dynamics as it specifies the form of the wave functions that characterize electron energy levels in a periodic crystal [13]. It was therefore useful in the formation of the tight-binding (TB) model which is an approach to the calculation of the electronic band structure of a lattice that relies on the fact that electrons are tightly bound (localized) to the atoms. The kinetic energy is included by allowing electrons to hop from one site to another. These features make the model to be lattice dependent. In its simplest form, the Hubbard model is an extension of the tight-binding model, wherein electrons can hop between lattice sites while the Coulomb interaction provides their localization tendency [14]. This formation makes the Hubbard model and all its extended versions to be lattice structure dependent and therefore emanate from the Bloch's theorem. For as it is well known, the general lattice Hamiltonian model within the occupation number formalism for electrons with spin σ interacting via a spin-dependent interaction $V^{ee}(\vec{r} - \vec{r}')$ in the presence of an ionic lattice potential $V^{ion}(\vec{r})$ has the form [10,12]

$$H = \hat{H}_0 + \hat{H}_{int} \quad (2.1)$$

where

$$\hat{H}_0 = \sum_{\sigma} \int d^3\vec{r} \psi_{\sigma}^+(\vec{r}) \left[-\frac{\hbar^2}{2m} \Delta + V^{ion}(\vec{r}) \right] \psi_{\sigma}(\vec{r}) \quad (2.2)$$

$$\hat{H}_{int} = \frac{1}{2} \sum_{\sigma\sigma'} \int d^3\vec{r} \int d^3\vec{r}' V^{ee}(\vec{r} - \vec{r}') \hat{n}_{\sigma}(\vec{r}) \hat{n}_{\sigma'}(\vec{r}') \quad (2.3)$$

In the above equation, the $\psi_{\sigma}(\vec{r}), \psi_{\sigma}^+(\vec{r})$ are the usual field operators and $\hat{n}_{\sigma}(\vec{r}) = \psi_{\sigma}^+(\vec{r}) \psi_{\sigma}(\vec{r})$ is the local density. It has been observed [12] that the interaction term is diagonal in the space variables \vec{r}, \vec{r}' [i.e. it depends only on the (operator-valued) densities of electrons at site \vec{r}, \vec{r}' which interact via $V^{ee}(\vec{r} - \vec{r}')$]. The lattice potential entering the non-interacting part (Eq.2.2) leads to the splitting of the parabolic dispersion into infinitely many bands which are enumerated by the index α . The non-interacting problem is then characterized by the Bloch wave functions $\phi_{\alpha k}(\vec{r})$ and the band energies $\mathcal{E}_{\alpha k}$. The

Wannier functions $\chi(\vec{r})$ localized at site R_i are often introduced:

$$\chi_{\alpha i}(\vec{r}) = \frac{1}{\sqrt{N}} \sum_k e^{-ik \cdot R_i} \phi_{\alpha k}(\vec{r}) \quad (2.4)$$

where N is the number of lattice sites. Thus one can construct creation and annihilation operators by the method of second quantization [13], $c_{\alpha i \sigma}^+$ and $c_{\alpha i \sigma}$ respectively, for electrons with spin σ in the band α at site R_i as

$$c_{\alpha i \sigma}^+ = \int d^3\vec{r} \chi_{\alpha i}(\vec{r}) \psi_{\sigma}^+(\vec{r}) \quad (2.5a)$$

where

$$\psi_{\sigma}^+(\vec{r}) = \sum_{i\alpha} \chi_{\alpha i}^+(\vec{r}) c_{\alpha i \sigma}^+ \quad (2.5b)$$

Consequently, the Hamiltonian may be written in the lattice representation as

$$H = \sum_{ij\sigma} t_{\alpha ij} c_{\alpha i \sigma}^+ c_{\alpha j \sigma} + \frac{1}{2} \sum_{\alpha\beta\gamma\delta} \sum_{ijmn} \sum_{\sigma\sigma'} v_{ijmn}^{\alpha\beta\gamma\delta} c_{\alpha i \sigma}^+ c_{\beta j \sigma'}^+ c_{\delta n \sigma'} c_{\gamma m \sigma} \quad (2.6)$$

where the matrix elements are given by

$$t_{\alpha ij} = \int d^3\vec{r} \chi_{\alpha i}^*(\vec{r}) \left[-\frac{\hbar^2}{2m} \Delta + V^{ion}(\vec{r}) \right] \chi_{\alpha j}(\vec{r}) \quad (2.7)$$

and

$$v_{ijmn}^{\alpha\beta\gamma\delta} = \int d^3\vec{r} \int d^3\vec{r}' V^{ee}(\vec{r} - \vec{r}') \chi_{\alpha i}^*(\vec{r}) \chi_{\beta j}^*(\vec{r}') \times \chi_{\delta n}(\vec{r}') \chi_{\gamma m}(\vec{r}) \quad (2.8)$$

The Hamiltonian given by Eq.(2.6) is too general to be tractable in dimensions $d > 1$. Therefore, it has to be simplified using physically motivated truncations such as:

- That the Fermi surface (FS) lies within a single conduction band that is well separated from other bands so that the inter band interaction is weak hence we will restrict our study to a single band ($\alpha = \beta = \gamma = \delta = 1$) which is the aforementioned TB band. Thus Eq. (2.6) reduces to

$$H_{1-band} = \sum_{ij\sigma} t_{ij} c_{i\sigma}^+ c_{j\sigma} + \frac{1}{2} \sum_{ijmn} \sum_{\sigma\sigma'} v_{ijmn} c_{i\sigma}^+ c_{j\sigma'}^+ c_{n\sigma'} c_{m\sigma} \quad (2.9)$$

- The above single band Hamiltonian is obviously still too complicated for most purposes because of the complex interaction part, and therefore needs to be simplified further. Taking into account the weak overlap between neighbouring orbitals in a TB description, one expects that the overlap between nearest neighbours (NN) is important hence only NN hopping and interactions are allowed. So the site indices are restricted to only NN positions i and j

leaving us with only NN hopping, a purely local contribution and four NN contributions. NN hopping between sites i and j are controlled by the $c_{i\sigma}^+ c_{j\sigma}$ which are the creation (annihilation) operators at site $i(j)$ having spin $\sigma(\sigma)$ while the hopping term is defined as $t_{ij} = t$ for NN hopping $\langle i,j \rangle$ (i.e. $i \neq j$) and $t_{ij} = 0$ otherwise.

The purely local contribution is the on-site Coulombic interaction term defined as

$$v_{iiii} = U \quad (2.10)$$

while the four NN contributions are:

$$v_{ijij} = V \quad (2.11)$$

which is the NN Coulombic interaction term,

$$v_{ijij} = \Delta t \quad (2.12)$$

which gives rise to an occupation dependent hopping rate,

$$v_{ijji} = J \quad (2.13)$$

which is the NN Heisenberg exchange term and

$$v_{ijij} = J' \quad (2.14)$$

which describes the exchange hopping processes.

This truncated form of the single band Hubbard model in Eq. (2.9) with only the Coulombic interaction (Eq. 2.10) is the famous Hubbard model given by [10-14]

$$H_H = -t \left[\sum_{\langle i,j \rangle, \sigma} c_{i\sigma}^+ c_{j\sigma} + H.C. \right] + U \sum_i n_{i\uparrow} n_{i\downarrow} \quad (2.15)$$

It is pertinent to point out that apart from the four NN contributions (Eqs. 2.11 – 2.14) which can be added to Eq.(2.15), there have been modifications of the Hubbard so that today it has numerous extensions and relatives [14-18] which have diverse application in condensed matter physics. Suffice to re-emphasize here, however, all these models depend on the lattice structure. For it is a common knowledge that lattice structure without periodic boundaries often leads to edge effect making some lattice sites not to be on equal footing with the others. As stated above, one way to resolve this limitation is to investigate the Hubbard lattice structure as torus in all three dimensions [2,4]. Here a torus is simply a lattice geometry with periodic boundary conditions [6] and is generic meaning that its nature depends on these boundary conditions of the system and other physical constraint conditions owing to the generic nature of the Oyibo GUT [1]. In this form of torus, the lattices are scale symmetric which means the lattices do not change their shapes when they are expanded into large lattices to study large densities or contracted into small lattices for few particles systems.

Now in geometrical terms, quantization is achieved by applying the so-called Born-Von Karman cyclic boundary conditions [18-19], which is tantamount to turning the cubic crystal into a torus as sketched in Fig. 1. The proposal in [20]

that the process of implementing cyclic boundary conditions would require going from the usual 3D to 5D background because the latter has in addition local degrees of freedom which become vital in introducing quantum effects, was used to show in [1] that the parametric equations of quantization of the torus into 5-dimensional background lattice is:

$$(ct)^2 + (aT)^2 - x^2 - y^2 - z^2 = 2(ct)(aT)\sin\theta \\ \equiv \begin{cases} \pm 2(ct)(aT), & \text{if } \theta = (n + \frac{1}{2})\pi, \\ 0, & \text{if } \theta = n\pi \end{cases} \quad (2.15)$$

where $n = 1, 2, 3, 4, 5$ by limiting $\theta = 2\pi/n$, needed for only rotations compatibles with translational symmetry of a crystal lattice in 3D space such that $2 \cos \theta - 1 = \text{integer}$.

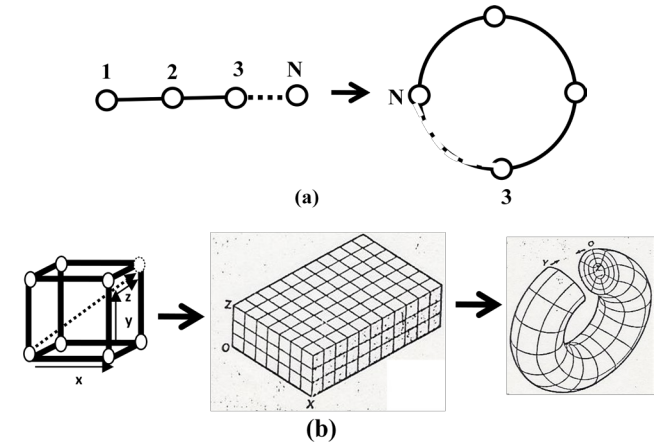


Figure 1. Illustrative sketch of the projection of (a) a linear lattice sites into a 1D ring (b) cubic lattice sites into curve torus

3. Correspondence of Bloch's Theorem and Oyibo GUT

For ease in drawing an analogy between Bloch's theorem and Oyibo GUT, which will enable us to establish their correspondence, a summary is provided in Table 1, adopting Oyibo's notations, as far as necessary, in order to facilitate identification of corresponding transformations and functions. Familiarity with both the Bloch's theorem [19, 21] and the Oyibo GUT (see [22] for pedagogical review) will be assumed.

According to Oyibo [8-9, 22], a function, $G = G(Y_1, Y_2, \dots, Y_p)$ is said to be conformal invariant under a given group transformation

$$T_k : Y_i = f_i(y_1, y_2, \dots, y_p, k) \quad (3.1)$$

if T_k is the group of the transformation and

$$G(Y_1, Y_2, \dots, Y_p) = F_i(y_1, y_2, \dots, y_p, k) \cdot G(y_1, y_2, \dots, y_p) \quad (3.2)$$

where $F_i(y_1, y_2, \dots, y_p, k)$ is a function of y_i and k the single group parameter.

From Table 1, we see that the group of translations in a crystal lattice and the Bloch (electron wave) function, more specifically lattice waves, have analogous properties as Eqs.(3.1) and (3.2). As a result, one needs only to show here that in quantized space-time geometry with given (Planck length) lattice spacing, the definition of conformal invariance in Eqs.(3.1) and (3.2) follows, as in Bloch's theorem, from translational periodicity of any function of position in the lattice.

On the basis of the definition in Eqs.(3.1) and (3.2), Oyibo states that there exists a set of conservation equations

$$(G_{0n})_t + (G_{1n})_x + (G_{2n})_y + (G_{3n})_z = 0, \quad (n = 0, 1, 2, 3, 4) \quad (3.3)$$

which may be rewritten in the Einstein-like form $G_{ij,j} = 0$ where the function defined mathematically by $G_{mn} = G_{mn}(x, y, z, ct, \dot{x}, \dot{y}, \dot{z}, \ddot{x}, \ddot{y}, \ddot{z}, \rho, \mu, T, P, \dots)$ is a set of "generic" quantities which are arbitrary functions of space and time coordinates (x, y, z, ct) , velocities $(\dot{x}, \dot{y}, \dot{z})$, accelerations $(\ddot{x}, \ddot{y}, \ddot{z})$, density (ρ) , fluid or gas viscosity (μ) , temperature (T) , pressure (P) , etc.

Now by applying the conformal transformation defined by Eq.(3.1) to a system of partial differential equations of n^{th} order given by

$$G_j(x^1 \dots x^p, y^1 \dots y^q, \frac{\partial^n y^1}{(\partial x^1)^n}, \dots, \frac{\partial^n y^q}{(\partial x^p)^n}) = 0, \quad (3.4)$$

one derives "solutions" of Eq.(3.2) in terms of the absolute invariants η_n of the subgroup of transformations for the independent coordinate variables in the form:

$$\eta_n = g_{n0}(ct)^{n+1} + g_{n1}x^{n+1} + g_{n2}y^{n+1} + g_{n3}z^{n+1} \quad (3.5)$$

where $n = 0, 1, 2, 3, 4$ are the five degrees of freedom of Oyibo GUT as in the Bloch's theorem shown in Table 1. Observe that we have indicated in Table 1 how the restriction to the *five* values of n in Eq.(3.3) are analogous to the constraints in Bloch's theorem, on translational invariance of the Hamiltonian operator, by rotational symmetry of a crystal lattice in 3-dimensional space represented by angles, $\theta = 2\pi/n$, where $n = 1, 2, 3, 4$ or 6, corresponding to n -fold rotation axis of symmetry [1].

Finally, analogy between Bloch's theorem and Oyibo GUT leads to identification of η_n with the wave number associated with the primitive translation vectors of the reciprocal lattice of a 3D crystal

$$\vec{\eta} = \frac{n_1 \vec{b}_1}{N_1} + \frac{n_2 \vec{b}_2}{N_2} + \frac{n_3 \vec{b}_3}{N_3}, \quad (3.6)$$

(n_1, n_2, n_3) being integers and $(\vec{b}_1, \vec{b}_2, \vec{b}_3)$ the primitive translation vectors of the reciprocal lattice space and $N_i = N_1 N_2 N_3$ the number of primitive cells in the crystal.

For a cubic lattice in \vec{r} -space, $|\vec{b}_i| = 2\pi/a_0$, where a_0 is the lattice constant, which we shall take (in Oyibo GUT representation of general theory of relativity) to be the Planck length.

However, the n^{th} order partial differential equations (3.4) are not quite analogous to the non-relativistic Schrodinger equation of Bloch's theorem in Table 1 except insofar as the system of partial differential in Eqs.(3.4) may be rewritten, in the first order, as follows

$$G_j\left(x^\mu, J^\mu, \frac{\partial J^\mu}{\partial x^\nu}\right) = 0 \quad (3.7)$$

in terms of the variables $(x^\mu, J^\mu) \equiv (x^1 \dots x^p, y^1 \dots y^q)$ where

$$J^\mu \equiv (\rho, \vec{J}) = \left(\psi^* \psi, \psi^* \left(-i \frac{\hbar}{m} \vec{\nabla} \right) \psi \right) \quad (3.8)$$

is the probability current density 4-vector conserved by Schrodinger's equation. It can easily be shown that if we introduce the differential operator $\partial_\mu = \partial/\partial x^\mu = (\partial/\partial t, \nabla)$ into Eq.(3.8), it becomes the current conservation law [15]

$$\partial_\mu J^\mu \equiv 0 \quad (3.9)$$

so that in a very subtle way we have obtained the Noether theorem. This is an important result as it re-enact the expectation of Oyibo when he pointed out in p285 of [8] that, "This new methodology is not totally new in the sense that the conservation equations are field equations... (see Eq.(3.3) here) could be condensed to tensor components similar to the world tensor components in the general theory of relativity of various versions of the many proposed Unified Force Field theories by previous investigators including Einstein himself. Perhaps one of the most significant part of the relationship to the previous formulations and results is the evolution of a characteristic sub-group variable (see Eq.(3.5) here) which seems to be a generalization of a good number of similar variables in the previous formulations."

Table 1. Analogy between Bloch's Theorem and Oyibo GUT

Bloch's Theorem (Lattice translational invariance)	Oyibo GUT (Conformal Transformation)
<p>Define lattice translation group of transformation</p> $T_k : \vec{R} = \vec{r} + \vec{k} \equiv \vec{r} + k_1\vec{a}_1 + k_2\vec{a}_2 + k_3\vec{a}_3 \leftrightarrow \vec{Y},$ <p>(k_1, k_2, k_3) being integers; $(\vec{a}_1, \vec{a}_2, \vec{a}_3)$ primitive lattice translation vectors,</p> <p>& a wave function $\psi \equiv \psi(\vec{R}) \leftrightarrow G(\vec{Y})$;</p> <p>or, for interatomic potential energy function, $\Phi(\vec{R}_1^0 + \vec{u}_1, \vec{R}_2^0 + \vec{u}_2, \dots) \leftrightarrow G(\vec{Y})$ and deformations, $\vec{u}_i (\equiv \delta \vec{x}_i)$ of a lattice from equilibrium positions, $\vec{R}_i^0 = k_i \vec{a}_i$, the force constants $(\partial^2 \Phi / \partial \vec{u}_i \partial \vec{u}_j)_0$ must satisfy</p> $\sum_j (\partial^2 \Phi / \partial \vec{u}_i \partial \vec{u}_j)_0 = 0, \Rightarrow \text{all force constant must depend only on the relative positions } (\vec{R}_i - \vec{R}_j) \text{ of the lattice sites .}$	<p>Define conformal transformation</p> $T_k : Y_i = f_i(y_1, y_2, y_3, k) \leftrightarrow \vec{R}_i \equiv x_i + k_i \vec{a}_i,$ <p>and $G \equiv G(Y_1, Y_2, Y_3) \leftrightarrow \psi(\vec{R})$ (for electron wave); or in terms of the potential energy, $\Phi(\vec{R}_1^0 + \vec{u}_1, \vec{R}_2^0 + \vec{u}_2, \dots)$ for deformation of points in a "Planck" lattice from equilibrium :</p> $G(Y_1, Y_2, Y_3) \leftrightarrow \frac{\partial^2 \Phi}{\partial \vec{u}_i \partial \vec{u}_j}$
<p>Then Bloch's Theorem (for electron wave) states that there exists a wave vector $(\vec{\eta})$ such that $T_k \psi(\vec{r}) \equiv \psi(\vec{R}) = \exp(i\vec{k} \cdot \vec{\eta}) \psi(\vec{r})$ is a solution of Schrodinger's equation for ψ :</p> $H(\vec{r})\psi(\vec{r}) \equiv \left(-\frac{\hbar^2}{2m} \nabla^2 + V(\vec{r}) \right) \psi(\vec{r}) = E \psi(\vec{r}),$ <p>where, $T_k H(\vec{r}) \equiv H(\vec{r} + \vec{k})$ is also invariant under rotation through θ, provided that $(1 - 2 \cos \theta) = \text{integer}$, m say, giving $m = 0, \pm 1, 2, 3$ i.e., $\theta = 2\pi / n$, with $n = 1, 2, 3, 4$ or 6.</p>	<p>Then Oyibo GUT states that there exists a function $F_i \leftrightarrow \exp(i\vec{\eta} \cdot [\vec{R}_i^0 - \vec{R}_j^0])$ such that</p> $T_k G(y_1, y_2, \dots, y_p) \equiv G(Y_1, Y_2, \dots, Y_p) = F_i(y_1, y_2, \dots, y_p, k) \bullet G(y_1, y_2, \dots, y_p),$ <p>\leftrightarrow there exists a dynamical matrix $D_{mn}(\vec{\eta})$</p> $= \exp(i\vec{\eta} \cdot [\vec{R}_i^0 - \vec{R}_j^0]) (\partial^2 \Phi / \partial \vec{u}_i \partial \vec{u}_j)_0$ <p>where D_{nm} is equivalent to the "generic" function $G_{nm} \equiv G_{nm}(\vec{r}, \dot{\vec{r}}, \ddot{\vec{r}}, ct, \rho, T, \dots)$ obeying the Einstein - like conservation equation $G_{ij,j} \equiv$</p> $(G_{0n})_t + (G_{1n})_x + (G_{2n})_y + (G_{3n})_z = 0.$ <p>Note the restriction on n to five values, $n = 0, 1, 2, 3, 4$. as in Bloch's theorem.</p>
<p>Eigenfunctions are $\psi(\vec{r}, \vec{\eta})$ where</p> $\vec{\eta} = \frac{n_1 \vec{b}_1}{N_1} + \frac{n_2 \vec{b}_2}{N_2} + \frac{n_3 \vec{b}_3}{N_3}, \text{ where } (n_1, n_2, n_3)$ <p>are integers, and $(\vec{b}_1, \vec{b}_2, \vec{b}_3)$ primitive translation vectors of a reciprocal lattice containing $N = N_1 N_2 N_3$ primitive cells.</p>	<p>"Solutions" in terms of the absolute invariants η_n of the subgroup of transformations for the independent coordinate variables in the form</p> $\eta_n = g_{n0}(ct)^{n+1} + g_{n1}x^{n+1} + g_{n2}y^{n+1} + g_{n3}z^{n+1}$ <p>$\leftrightarrow \vec{q} ^{n+1}$, where $(n = 0, 1, 2, 3, 4)$.</p>

4. Summary and Conclusions

The Oyibo GUT was formulated from the Navier-Stokes equations which have classical origin and this raised doubt for its applicability to quantum systems. We have shown in a previous study that the Navier-Stokes equation has periodic solutions and then use it to formulate an $O(4,2) \times SU(3) \times U(1)$ gauge theory of quantum gravity [23] and in a more recent study one of us has shown that the geometric object for its geometrical and quantization foundation is the generic torus which has also been used in the literature as the geometric object of the Hubbard model formulated on the basis of the Bloch's theorem. It turns out that there is a remarkable correspondence between Bloch's theorem and the Oyibo GUT. This has open the possibility that the Oyibo GUT can in principle be used to formulate strongly correlated models which emanates from the Bloch's theorem. In particular, we have reviewed the formation of the Hubbard model and its extended versions from the Bloch's theorem. The implication is that the Hubbard model and its extended versions can in principle also be formulated from the Oyibo GUT. Thereafter, since the Hubbard model and its extended versions have very important applications in condensed matter physics, we claim that such applications have been brought under the purview of the Oyibo GUT and therefore constitute this related area in the Oyibo grand unified theory. Thus we have demonstrated an important conjecture of Oyibo in p303 of [8] that, "The generic formulation of the grand unified theorem is in good part based on synthesizing, interpreting and realizing that generic, universal conservation equations should in light of the general and global outlook, be expected to be the source from which the unified force field can be derived [8]."

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