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On the linear representations of the symmetry groups of single-wall carbon nanotubes

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Abstract

The positions of atoms forming a carbon nanotube are usually described by using a system of generators of the symmetry group. Each atomic position corresponds to an element of the set $\mathbb{Z} \times \{0, 1, \ldots, n\} \times \{0, 1\}$, where n depends on the considered nanotube. We obtain an alternative, rather different description by starting from a three-axes description of the honeycomb lattice. In our mathematical model, which is a factor space defined by an equivalence relation in the set $\{(v_0, v_1, v_2) \in \mathbb{Z}^3 \mid v_0+v_1+v_2 \in \{0, 1\}\}$, the neighbours of an atomic position can be described in a simpler way, and the mathematical objects with geometric or physical significance have a simpler and more symmetric form. We present some results concerning the linear representations of the symmetry groups of single-wall carbon nanotubes in order to illustrate the proposed approach.

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1. Introduction

A single-wall carbon nanotube is a cylindrical structure with a diameter of a few nanometres, periodic along its axis, which can be imagined as a rolled up honeycomb lattice. The high symmetry of carbon nanotubes has facilitated the theoretical investigation of the physical phenomena occurring in these materials [1, 8–10, 12–14, 16]. The spatial symmetries (translations, rotations and screw axes, mirror and glide planes, etc) form a line group, which is the maximal subgroup of the Euclidean group that leaves the nanotube invariant. The role of this group is analogous to that of crystallographic space groups in solid-state physics. Some important properties of the band structure (electronic, phonon, etc) can be directly deduced from the symmetry groups.

The symmetry group of a carbon nanotube depends on the diameter of the tubule and on the helical arrangement of the carbon hexagons. The irreducible representations of these

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groups are well known [6, 7, 15], but, generally, equivalent representations offer distinct formal advantages. A calculation very simple in a representation can become much more complicated in an equivalent representation. Therefore, we think it is worth looking for new representations and for new ways to describe the atomic structure of these materials. Our aim is to present an improved version of the mathematical model proposed in [3] and some applications illustrating this new approach.

2. Honeycomb lattice in a three-axes description

The vectors corresponding to the vertices of a regular triangle

$$e_0 = (1, 0), \qquad e_2 = \left(-\tfrac{1}{2}, \tfrac{\sqrt{3}}{2}\right), \qquad e_3 = \left(-\tfrac{1}{2}, -\tfrac{\sqrt{3}}{2}\right) \tag{1}$$

allow us to define the bijecton

$$\mathcal{L} \longrightarrow \mathbb{L} : (v_0, v_1, v_2) \mapsto v_0 e_0 + v_1 e_1 + v_2 e_2 \tag{2}$$

from the set

$$\mathcal{L} = \{v = (v_0, v_1, v_2) \in \mathbb{Z}^3 \mid v_0 + v_1 + v_2 \in \{0; 1\}\} \tag{3}$$

to the set $\mathbb{L}$ of all the vertices of a *honeycomb lattice*. The subset $\mathcal{L}$ of $\mathbb{Z}^3$ becomes in this way a mathematical model for the honeycomb lattice. One can remark that

$$\mathcal{L} = \mathcal{T} \cup (\mathcal{T} + (1, 0, 0)), \tag{4}$$

where

$$\mathcal{T} = \{v = (v_0, v_1, v_2) \in \mathbb{Z}^3 \mid v_0 + v_1 + v_2 = 0\}. \tag{5}$$

The mapping

$$d : \mathcal{L} \times \mathcal{L} \longrightarrow \mathbb{N}, \qquad d(v, u) = |v_0 - u_0| + |v_1 - u_1| + |v_2 - u_2| \tag{6}$$

is a distance on $\mathcal{L}$, and u is a *neighbour of order* l of v if $d(v, u) = l$. The *nearest neighbours* of v are

$$\begin{aligned} v^0 &= (v_0 + \varepsilon(v), v_1, v_2) \\ v^1 &= (v_0, v_1 + \varepsilon(v), v_2) \qquad \text{where} \qquad \varepsilon(v) = (-1)^{v_0+v_1+v_2} \\ v^2 &= (v_0, v_1, v_2 + \varepsilon(v)), \end{aligned} \tag{7}$$

and the six *next-to-nearest neighbours* of v are the points $v^{ij} = (v^i)^j$ corresponding to all the pairs (i, j) with $i \neq j$. The *symmetry group* G of the honeycomb lattice coincides with the group of all the isometries of the metric space $(\mathcal{L}, d)$ and is generated by the transformations

$$\begin{aligned} &\mathcal{L} \longrightarrow \mathcal{L} : \quad (v_0, v_1, v_2) \mapsto (v_1, v_2, v_0) \\ &\mathcal{L} \longrightarrow \mathcal{L} : \quad (v_0, v_1, v_2) \mapsto (v_0, v_2, v_1) \\ &\mathcal{L} \longrightarrow \mathcal{L} : \quad (v_0, v_1, v_2) \mapsto (-v_0 + 1, -v_1, -v_2). \end{aligned} \tag{8}$$

The subgroup of *translations* contained in G corresponds to $\mathcal{T}$:

$$\{u \mid v \in \mathcal{L} \Rightarrow v + u \in \mathcal{L}\} = \{u \in \mathcal{L} \mid \varepsilon(u) = 1\} = \mathcal{T}. \tag{9}$$

We can extend the description based on e_0, e_1, e_2 to the whole plane. Each vector v admits the representation

$$v = a \sum_{i=0}^{2} \langle v, e_i \rangle e_i \qquad \text{with} \quad a = \frac{2}{3} \tag{10}$$

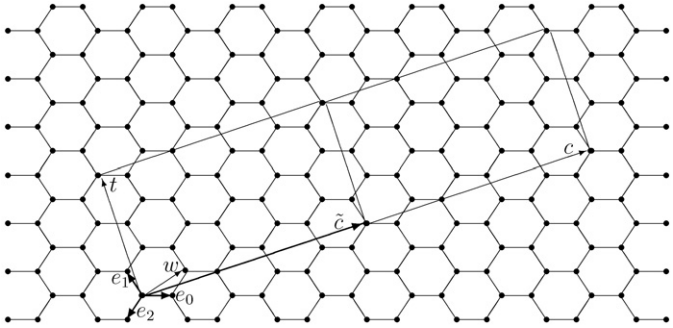

Figure 1. The unit cell of the carbon nanotube with the chiral vector $c = (10, -2, -8)$. In this case $n = 2$, $\tilde{c} = (5, -1, -4)$, $t = (-1, 3, -2)$, $w = (1, 0, -1)$ and $\tilde{q} = 14$.

and the usual scalar product and norm become

$$\langle v, u\rangle = a\sum_{i=0}^{2}\langle v, e_i\rangle\langle u, e_i\rangle, \qquad \|v\| = \sqrt{a\sum_{i=0}^{2}\langle v, e_i\rangle^2}. \tag{11}$$

The vectors e_0, e_1, e_2 form a tight frame [2] and a system of coherent vectors [4] in $\mathbb{R}^2$. For each vector v, the 'canonical coordinates'

$$\hat{v}_0 = \langle v, e_0\rangle, \qquad \hat{v}_1 = \langle v, e_1\rangle, \qquad \hat{v}_2 = \langle v, e_2\rangle \tag{12}$$

satisfy the relation $\hat{v}_0 + \hat{v}_1 + \hat{v}_2 = 0$, and we can identify $\mathbb{R}^2$ with the space

$$\mathcal{K} = \{k = (k_0, k_1, k_2) \mid k_0, k_1, k_2 \in \mathbb{R}, k_0 + k_1 + k_2 = 0\} \tag{13}$$

by using the linear isomorphism

$$\mathcal{K} \longrightarrow \mathbb{R}^2 : \quad (k_0, k_1, k_2) \mapsto a\sum_{i=0}^{2} k_i e_i. \tag{14}$$

The representation of a vector v as a linear combination of e_0, e_1, e_2 is not unique:

$$v = a\sum_{i=0}^{2}\hat{v}_i e_i = a\sum_{i=0}^{2}(\hat{v}_i + \alpha)e_i \tag{15}$$

for any $\alpha \in \mathbb{R}$. All the elements (v_0, v_1, v_2) of the set

$$\{(\hat{v}_0 + \alpha, \hat{v}_1 + \alpha, \hat{v}_2 + \alpha) \mid \alpha \in \mathbb{R}\}$$

correspond to the same point of the plane, and

$$\langle k, v\rangle = a\sum_{i=0}^{2} k_i\hat{v}_i = a\sum_{i=0}^{2} k_i(\hat{v}_i + \alpha) = a\sum_{i=0}^{2} k_i v_i \tag{16}$$

for all $k \in \mathcal{K}$.

3. An alternative mathematical model for carbon nanotubes

A *single-wall carbon nanotube* can be visualized as the structure obtained by rolling a honeycomb lattice (which is a mathematical model for a graphene sheet) such that the endpoints of a translation vector $c \in \mathcal{T}$ are folded one onto the other (see figure 1). The

vector c is called the *chirality* of the tubule. After the graphene sheet rolling, the points $\ldots, v-2c, v-c, v, v+c, v+2c, \ldots$ are folded one onto the other, for any $v \in \mathcal{L}$. Therefore, each element of the set $[v] = v + \mathbb{Z}c$, that is, each element of the set

$$[v_0, v_1, v_2] = \{(v_0, v_1, v_2) + j\,(c_0, c_1, c_2) \mid j \in \mathbb{Z}\} \tag{17}$$

describes the same point of the nanotube. The subset of the factor space $\mathbb{Z}^3/\mathbb{Z}c$,

$$\mathcal{L}_c = \left\{ [v_0, v_1, v_2] \in \frac{\mathbb{Z}^3}{\mathbb{Z}c} \,\middle|\, v_0 + v_1 + v_2 \in \{0; 1\} \right\}, \tag{18}$$

can be regarded as a mathematical model for the nanotube of chirality c. A symmetry transformation $\mathcal{L} \longrightarrow \mathcal{L} : v \mapsto gv$ of the honeycomb lattice satisfying the relation

$$[v] = [u] \quad \Longrightarrow \quad [gv] = [gu] \tag{19}$$

defines a *symmetry transformation* of nanotube $\mathcal{L}_c$, namely

$$g : \mathcal{L}_c \longrightarrow \mathcal{L}_c : [v] \mapsto [gv]. \tag{20}$$

Each nanotube $\mathcal{L}_c$ admits the symmetry transformations:

$$\begin{aligned} &\tau : \mathcal{L}_c \longrightarrow \mathcal{L}_c : [v_0, v_1, v_2] \mapsto [-v_0 + 1, -v_1, -v_2] \\ &g_u : \mathcal{L}_c \longrightarrow \mathcal{L}_c : [v] \mapsto [v + u] \qquad \text{for any} \qquad u \in \mathcal{T}. \end{aligned} \tag{21}$$

Let n be the greatest common divisor of c_0, c_1, c_2, and let $\tilde{c} = (\tilde{c}_0, \tilde{c}_1, \tilde{c}_2)$, where

$$\tilde{c}_0 = \frac{1}{n} c_0, \qquad \tilde{c}_1 = \frac{1}{n} c_1, \qquad \tilde{c}_2 = \frac{1}{n} c_2. \tag{22}$$

The transformation $g_{\tilde{c}}$ represents a *rotation of nanotube* of angle $2\pi/n$ with respect to its axis. Since

$$(c_2 - c_1)c_0 + (c_0 - c_2)c_1 + (c_1 - c_0)c_2 = 0 \tag{23}$$

the vector $(c_2 - c_1, c_0 - c_2, c_1 - c_0)$ is orthogonal to c (it has the direction of nanotube axis). The transformation g_t corresponding to

$$t = \frac{1}{\mathcal{R}}(c_2 - c_1, c_0 - c_2, c_1 - c_0), \tag{24}$$

where

$$\mathcal{R} = \gcd\{c_2 - c_1, c_0 - c_2, c_1 - c_0\} = \begin{cases} 3n & \text{if } \tilde{c}_2 - \tilde{c}_1 \in 3\mathbb{Z} \\ n & \text{if } \tilde{c}_2 - \tilde{c}_1 \notin 3\mathbb{Z} \end{cases} \tag{25}$$

represents the shortest pure translation of nanotube. From $c_0 + c_1 + c_2 = 0$ we get

$$(c_1 - c_2)^2 + (c_2 - c_0)^2 + (c_0 - c_1)^2 = 3\left(c_0^2 + c_1^2 + c_2^2\right) \tag{26}$$

that is, $\mathcal{R}^2 \|t\|^2 = 3\|c\|^2$, whence

$$q = \frac{1}{\mathcal{R}}\left(c_0^2 + c_1^2 + c_2^2\right) \in n\mathbb{Z}. \tag{27}$$

For any $u \in \mathcal{T}$ the projections of u on c and t can be written as

$$\frac{\langle u, c\rangle}{\|c\|^2} c = \left(u_1 \frac{c_1 - c_0}{\mathcal{R}} + u_2 \frac{c_2 - c_0}{\mathcal{R}}\right) \frac{c}{q}, \qquad \frac{\langle u, t\rangle}{\|t\|^2} t = \left(u_1 \frac{c_2}{n} - u_2 \frac{c_1}{n}\right) \frac{t}{\tilde{q}}, \tag{28}$$

where $\tilde{q} = \frac{q}{n}$. Since

$$\gcd\{(c_1 - c_0)/\mathcal{R}, (c_2 - c_0)/\mathcal{R}\} = 1 \qquad \text{and} \qquad \gcd\{c_2/n, c_1/n\} = 1 \tag{29}$$

it follows that the projection of $\mathcal{T}$ on c is $\mathbb{Z}\frac{c}{q}$, and the projection of $\mathcal{T}$ on t is $\mathbb{Z}\frac{t}{\tilde{q}}$. Let $w \in \mathcal{T}$ be the shortest vector with

$$\frac{\langle w, t\rangle}{\|t\|^2} t = \frac{t}{\tilde{q}}. \tag{30}$$

From the relation

$$w = \frac{\langle w, c\rangle}{\|c\|^2} c + \frac{\langle w, t\rangle}{\|t\|^2} t \tag{31}$$

we get $[qw] = [nt]$.

Without loss of generality we can assume $c_0 > c_1 \geqslant c_2$. In the case $c_1 = c_2$ we have an *armchair* nanotube, and in the case $c_1 = 0$ a *zig-zag* nanotube. The nanotubes with $0 \neq c_1 \neq c_2$ are called *chiral nanotubes*. Our approach works for any single-wall carbon nanotube, but in this paper we restrict ourselves to chiral nanotubes. The symmetry group G_c of the nanotube $\mathcal{L}_c$ is generated by the transformations $\varrho = g_{\tilde{c}}$, $\sigma = g_w$ and τ:

$$G_c = \langle \varrho, \sigma, \tau \mid \varrho\sigma = \sigma\varrho, \varrho^n = \tau^2 = (\sigma\tau)^2 = (\varrho\tau)^2 = e\rangle. \tag{32}$$

For each $[v] \in \mathcal{L}_c$, there exist $s \in \mathbb{Z}$, $m \in \{0, 1, \ldots, n-1\}$ and $p \in \{0; 1\}$ uniquely determined such that $[v] = \sigma^s \varrho^m \tau^p [0, 0, 0]$. The usual description [5, 15] of the atomic positions of the atoms forming a carbon nanotube is based on this remark, and the set

$$\{(s, m, p) \mid s \in \mathbb{Z}, m \in \{0, 1, \ldots, n-1\}, p \in \{0, 1\}\} \tag{33}$$

is used as a mathematical model. The alternative mathematical model we present in this paper is $\mathcal{L}_c$.

4. A tight-binding approach to carbon nanotubes

Consider the Hilbert space $(l^2(\mathcal{L}_c), \langle , \rangle)$, where

$$l^2(\mathcal{L}_c) = \left\{ \psi : \mathcal{L}_c \longrightarrow \mathbb{C} \,\middle|\, \sum_{v \in \mathcal{L}_c} |\psi(v)|^2 < \infty \right\} \tag{34}$$

$$\langle \psi_1, \psi_2 \rangle = \sum_{v \in \mathcal{L}_c} \overline{\psi_1(v)} \psi_2(v) \tag{35}$$

and the unitary representation of G_c in $l^2(\mathcal{L}_c)$ is defined by

$$g : l^2(\mathcal{L}_c) \longrightarrow l^2(\mathcal{L}_c) \qquad (g\psi)[v] = \psi(g^{-1}[v]). \tag{36}$$

For each $\kappa \in (0, \infty)$, the linear operator

$$H : l^2(\mathcal{L}_c) \longrightarrow l^2(\mathcal{L}_c) \qquad (H\psi)[v] = \kappa \sum_{j=0}^{2} \psi[v^j] \tag{37}$$

is a self-adjoint G_c-invariant operator.

If $k \in \mathcal{K}$ is such that $\langle k, c\rangle \in 2\pi\mathbb{Z}$ then

$$\mathrm{e}^{\mathrm{i}\langle k, v\rangle} = \mathrm{e}^{\mathrm{i}\langle k, v + jc\rangle} \qquad \text{for any} \quad j \in \mathbb{Z} \tag{38}$$

and hence, the function

$$\mathcal{L}_c \longrightarrow \mathbb{C} : [v] \mapsto \mathrm{e}^{-\mathrm{i}\langle k, v\rangle} \tag{39}$$

is well defined (it does not depend on the representative we choose for $[v]$). The Hamiltonian used in the tight-binding description of π bands has the form (37).

Theorem 1.

(a) For any k such that $\langle k, c\rangle \in 2\pi\mathbb{Z}$ the numbers $\pm E(k)$, where

$$E(k) = \kappa|\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a}| = \kappa\sqrt{3 + 2\cos(k_0 - k_1)a + 2\cos(k_1 - k_2)a + 2\cos(k_2 - k_0)a} \quad (40)$$

belong to the spectrum of H.

(b) The bounded functions belonging to an extension of $l^2(\mathcal{L}_c)$

$$\psi_k^{\pm} : \mathcal{L}_c \longrightarrow \mathbb{C}, \qquad \psi_k^{\pm}[v] = \mathrm{e}^{-\mathrm{i}\langle k, v\rangle}\varphi_k^{\pm}[v], \quad (41)$$

where

$$\varphi_k^{\pm}[v] = \begin{cases} \mathrm{e}^{\mathrm{i}\lambda(k)} & \text{if } \varepsilon(v) = 1 \\ \pm\mathrm{e}^{-\mathrm{i}\lambda(k)} & \text{if } \varepsilon(v) = -1 \end{cases} \quad (42)$$

and

$$\lambda(k) = \begin{cases} -\frac{1}{2}\arg(\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a}) & \text{if } \mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a} \neq 0 \\ 0 & \text{if } \mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a} = 0 \end{cases} \quad (43)$$

are eigenfunctions of H corresponding to the eigenvalues $\pm E(k)$, that is,

$$H\psi_k^{\pm} = \pm E(k)\psi_k^{\pm}. \quad (44)$$

Proof.

(a) The function $\psi : \mathcal{L}_c \longrightarrow \mathbb{C}$, $\psi[v] = \mathrm{e}^{-\mathrm{i}\langle k, v\rangle}\varphi[v]$, where

$$\varphi[v] = \begin{cases} \alpha & \text{if } \varepsilon(v) = 1 \\ \beta & \text{if } \varepsilon(v) = -1 \end{cases} \quad (45)$$

and α, β are two constants, satisfies the relation $H\psi = E\psi$ if and only if (α, β) is a solution of the system of equations

$$\begin{cases} \kappa(\mathrm{e}^{-\mathrm{i}k_0a} + \mathrm{e}^{-\mathrm{i}k_1a} + \mathrm{e}^{-\mathrm{i}k_2a})\beta = E\alpha \\ \kappa(\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a})\alpha = E\beta. \end{cases} \quad (46)$$

This system has non-trivial solutions if and only if

$$\begin{vmatrix} -E & \kappa(\mathrm{e}^{-\mathrm{i}k_0a} + \mathrm{e}^{-\mathrm{i}k_1a} + \mathrm{e}^{-\mathrm{i}k_2a}) \\ \kappa(\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a}) & -E \end{vmatrix} = 0, \quad (47)$$

that is, if and only if E is one of the numbers $\pm E(k)$.

(b) If $\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_0a} \neq 0$ then the equation $\kappa(\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a})\alpha = \pm E(k)\beta$ leads to

$$\beta = \pm\alpha\frac{\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a}}{|\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_2a}|} = \pm\alpha\,\mathrm{e}^{-2\mathrm{i}\lambda(k)}. \quad (48)$$

Choosing $\alpha = \mathrm{e}^{\mathrm{i}\lambda(k)}$ we get $\beta = \pm\mathrm{e}^{-\mathrm{i}\lambda(k)}$, and hence, up to a multiplicative constant, the solution of $H\psi = \pm E(k)\psi$ is $\psi = \psi_k^{\pm}$. If $\mathrm{e}^{\mathrm{i}k_0a} + \mathrm{e}^{\mathrm{i}k_1a} + \mathrm{e}^{\mathrm{i}k_0a} = 0$ then $E(k) = 0$, and the functions

$$\psi_k^{\pm}[v] = \mathrm{e}^{-\mathrm{i}\langle k, v\rangle}\begin{cases} 1 & \text{if } \varepsilon(v) = 1 \\ \pm 1 & \text{if } \varepsilon(v) = -1 \end{cases} \quad (49)$$

are eigenfunctions. □

The relation $\langle k, c\rangle \in 2\pi\mathbb{Z}$ defines a family of equidistant straight lines orthogonal to c with the distance between neighbouring lines equal to $2\pi/\|c\|$.

Theorem 2.

(a) The function

$$E : \mathcal{K} \longrightarrow [0, 3\kappa] \qquad E(k) = \kappa |e^{ik_0 a} + e^{ik_1 a} + e^{ik_2 a}| \tag{50}$$

is even and periodic

$$E(k) = E(-k) \qquad E(k) = E(k + b_0) = E(k + b_1) = E(k + b_2) \tag{51}$$

where

$$\begin{aligned} b_0 &= \left(\frac{4\pi}{3a}, -\frac{2\pi}{3a}, -\frac{2\pi}{3a}\right), \\ b_1 &= \left(-\frac{2\pi}{3a}, \frac{4\pi}{3a}, -\frac{2\pi}{3a}\right), \\ b_2 &= \left(-\frac{2\pi}{3a}, -\frac{2\pi}{3a}, \frac{4\pi}{3a}\right). \end{aligned} \tag{52}$$

(b) The functions $\psi_k = \psi_k^+$ and ψ_k^- are eigenfunctions of ϱ and σ

$$\varrho \psi_k^\pm = e^{i\langle k, \tilde{c}\rangle} \psi_k^\pm \qquad \sigma \psi_k^\pm = e^{i\langle k, w\rangle} \psi_k^\pm \tag{53}$$

and satisfy the relation

$$\tau \psi_k^\pm = \pm e^{-ik_0 a} \psi_{-k}^\pm. \tag{54}$$

(c) The eigenspaces

$$\begin{aligned} \mathcal{E}_k &= \{\alpha\psi_k + \beta\psi_{-k} \mid \alpha, \beta \in \mathbb{C}\} \\ \mathcal{E}_k^- &= \{\alpha\psi_k^- + \beta\psi_{-k}^- \mid \alpha, \beta \in \mathbb{C}\} \end{aligned} \tag{55}$$

corresponding to $E(k)$ and $-E(k)$ are G_c-invariant, $\mathcal{E}_k = \mathcal{E}_{-k}$ and $\mathcal{E}_k^- = \mathcal{E}_{-k}^-$.

(d) If k is such that $E(k) \neq 0$ then

$$\psi_{k+b_i}^\pm = e^{i\pi/3} \psi_k^\pm \qquad \mathcal{E}_{k+b_i} = \mathcal{E}_k \tag{56}$$

for any $i \in \{0, 1, 2\}$.

Proof. Since $e^{i4\pi/3} = e^{-i2\pi/3}$ we have

$$E(k + b_0) = \kappa |e^{ik_0 a}\, e^{i4\pi/3} + e^{ik_1 a}\, e^{-i2\pi/3} + e^{ik_2 a}\, e^{-i2\pi/3}| = E(k).$$

$$\lambda(k + b_0) = -\frac{1}{2} \arg[(e^{ik_0 a} + e^{ik_1 a} + e^{ik_2 a})\, e^{-i2\pi/3}] = \lambda(k) + \frac{\pi}{3}$$

for any k with $E(k) \neq 0$. If v is such that $\varepsilon(v) = 1$ then $\langle b_0, v\rangle = 2\pi v_0$ and

$$\psi_{k+b_0}[v] = e^{-i\langle k+b_0, v\rangle}\, e^{i\lambda(k+b_0)} = e^{-i\langle k, v\rangle}\, e^{-i2\pi v_0}\, e^{i\lambda(k)}\, e^{i\pi/3} = e^{i\pi/3} \psi_k[v].$$

If v is such that $\varepsilon(v) = -1$ then $\langle b_0, v\rangle = 2\pi v_0 - 2\pi/3$ and

$$\psi_{k+b_0}[v] = e^{-i\langle k+b_0, v\rangle}\, e^{-i\lambda(k+b_0)} = e^{-i\langle k, v\rangle}\, e^{-i2\pi v_0}\, e^{i2\pi/3}\, e^{-i\lambda(k)}\, e^{-i\pi/3} = e^{i\pi/3} \psi_k[v]. \qquad \square$$

From the periodicity of $E(k)$ and $\mathcal{E}_k$ it follows that we can restrict our analysis to the case $k \in \mathcal{B}_c$, where

$$\mathcal{B}_c = \{k \in \mathcal{B} \mid \langle k, c\rangle \in 2\pi\mathbb{Z}\} \tag{57}$$

and $\mathcal{B}$ is the hexagonal domain (see figure 2)

$$\mathcal{B} = \left\{ k \in \mathcal{K} \,\middle|\, -\frac{2\pi}{3a} \leqslant k_i \leqslant \frac{2\pi}{3a} \text{ for any } i \in \{0, 1, 2\} \right\}. \tag{58}$$

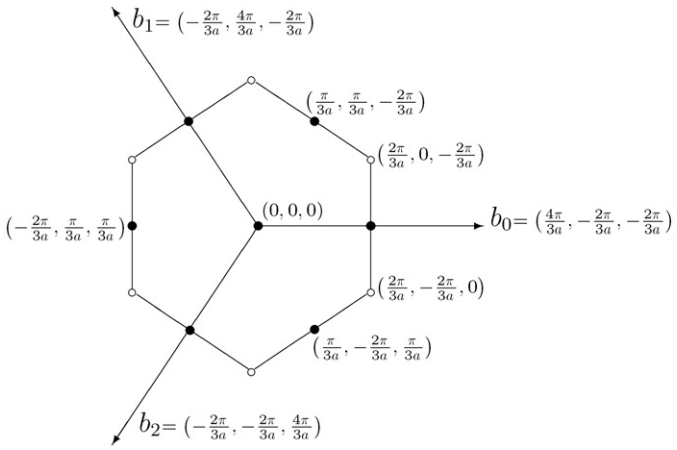

Figure 2. The hexagonal domain $\mathcal{B}$, the points of Λ (indicated by $\bullet$) and the points k with $E(k) = 0$, called K points (indicated by $\circ$).

Theorem 3. *For even n the space $\mathcal{E}_k$ is one dimensional if and only if $k \in \Lambda$, where*

$$\Lambda = \left\{(0,0,0), \pm\left(-\frac{2\pi}{3a}, \frac{\pi}{3a}, \frac{\pi}{3a}\right), \pm\left(\frac{\pi}{3a}, -\frac{2\pi}{3a}, \frac{\pi}{3a}\right), \pm\left(\frac{\pi}{3a}, \frac{\pi}{3a}, -\frac{2\pi}{3a}\right)\right\}.$$

Proof. If $\mathcal{E}_k$ is one dimensional then there is a constant C such that

$$\psi_{-k}[v] = C\psi_k[v] \qquad \text{for any} \quad v \in \mathcal{L}_c.$$

For $v = (0,0,0)$ we get $C = \mathrm{e}^{-2\mathrm{i}\lambda(k)}$. Therefore,

$$\psi_{-k} = C\psi_k \quad \Longleftrightarrow \quad \begin{cases} \mathrm{e}^{2\mathrm{i}\langle k,v\rangle} = 1 & \text{if } \varepsilon(v) = 1 \\ \mathrm{e}^{2\mathrm{i}\langle k,v\rangle} = C^2 & \text{if } \varepsilon(v) = -1. \end{cases} \tag{59}$$

Particularly, we must have $\mathrm{e}^{2\mathrm{i}k_0 a} = \mathrm{e}^{2\mathrm{i}k_1 a} = \mathrm{e}^{2\mathrm{i}k_2 a}$. For $k \in \mathcal{B}$ this relation is possible only if there are $\alpha, \beta \in \{0, \pm 1\}$ such that $k_1 = k_0 + \alpha\pi/a$ and $k_2 = k_0 + \beta\pi/a$. From $k_0 + k_1 + k_2 = 0$ we get

$$k = \left(-\frac{\alpha+\beta}{3}\frac{\pi}{a}, \frac{2\alpha-\beta}{3}\frac{\pi}{a}, \frac{-\alpha+2\beta}{3}\frac{\pi}{a}\right). \tag{60}$$

All these points lie on the family of straight lines defined by $\langle k, c\rangle \in 2\pi\mathbb{Z}$ since

$$\langle k, c\rangle = k_0 c_0 a + \left(k_0 + \alpha\frac{\pi}{a}\right)c_1 a + \left(k_0 + \beta\frac{\pi}{a}\right)c_2 a = (\alpha c_1 + \beta c_2)\pi \in 2\pi\mathbb{Z}. \tag{61}$$

In the case $\alpha = \beta = 0$ we get $k = (0,0,0)$, $\psi_k[v] = 1$ and

$$\varrho\psi_k = \psi_k \qquad \sigma\psi_k = \psi_k \qquad \tau\psi_k = \psi_k. \tag{62}$$

In the case $\alpha = \beta = 1$ we get $k = (-2\pi/3a, \pi/3a, \pi/3a)$, $\psi_k[v] = (-1)^{v_0}\,\mathrm{e}^{-\mathrm{i}\pi/6}$ and

$$\varrho\psi_k = (-1)^{\tilde{c}_0}\psi_k \qquad \sigma\psi_k = (-1)^{w_0}\psi_k \qquad \tau\psi_k = -\psi_k. \tag{63}$$

All the other points $k \in \mathcal{B}_c$ with $\dim \mathcal{E}_k = 1$ can be obtained from the analysed points by using permutations of coordinates and/or multiplication by (-1). □

In the case when n is odd only a part of the points of Λ lies on the family of straight lines defined by $\langle k, c\rangle \in 2\pi\mathbb{Z}$.

5. Two-dimensional representations of G_c

If $k \in \mathcal{B}_c\backslash\Lambda$ then $\dim \mathcal{E}_k = 2$ and the matrices of ϱ, σ and τ in the basis $\{\psi_k, \psi_{-k}\}$ are

$$\varrho = \begin{pmatrix} \mathrm{e}^{\mathrm{i}\langle k,\tilde{c}\rangle} & 0 \\ 0 & \mathrm{e}^{-\mathrm{i}\langle k,\tilde{c}\rangle} \end{pmatrix}, \qquad \sigma = \begin{pmatrix} \mathrm{e}^{\mathrm{i}\langle k,w\rangle} & 0 \\ 0 & \mathrm{e}^{-\mathrm{i}\langle k,w\rangle} \end{pmatrix}, \qquad \tau = \begin{pmatrix} 0 & \mathrm{e}^{\mathrm{i}k_0a} \\ \mathrm{e}^{-\mathrm{i}k_0a} & 0 \end{pmatrix}. \tag{64}$$

Theorem 4. *For each $k \in \mathcal{B}_c\backslash\Lambda$ the representations of G_c in $\mathcal{E}_k$ and $\mathcal{E}_k^-$ are equivalent. They are reducible if and only if the numbers $\langle k, \tilde{c}\rangle$ and $\langle k, w\rangle$ belong to $\mathbb{Z}\pi$.*

Proof. The linear transformation $\mathcal{E}_k \longrightarrow \mathcal{E}_k^- : \alpha\psi_k + \beta\psi_{-k} \mapsto \mathrm{i}\alpha\psi_k^- - \mathrm{i}\beta\psi_{-k}^-$ is an isomorphism of representations. If

$$P = \begin{pmatrix} \alpha & \beta \\ \gamma & \delta \end{pmatrix}$$

is the matrix in the basis $\{\psi_k, \psi_{-k}\}$ of the projector corresponding to a G_c-invariant subspace of $\mathcal{E}_k$ then $P^2 = P$ and

$$\begin{pmatrix} \alpha & \beta \\ \gamma & \delta \end{pmatrix}\begin{pmatrix} \mathrm{e}^{\mathrm{i}\langle k,\tilde{c}\rangle} & 0 \\ 0 & \mathrm{e}^{-\mathrm{i}\langle k,\tilde{c}\rangle} \end{pmatrix} = \begin{pmatrix} \mathrm{e}^{\mathrm{i}\langle k,\tilde{c}\rangle} & 0 \\ 0 & \mathrm{e}^{-\mathrm{i}\langle k,\tilde{c}\rangle} \end{pmatrix}\begin{pmatrix} \alpha & \beta \\ \gamma & \delta \end{pmatrix} \tag{65}$$

$$\begin{pmatrix} \alpha & \beta \\ \gamma & \delta \end{pmatrix}\begin{pmatrix} \mathrm{e}^{\mathrm{i}\langle k,w\rangle} & 0 \\ 0 & \mathrm{e}^{-\mathrm{i}\langle k,w\rangle} \end{pmatrix} = \begin{pmatrix} \mathrm{e}^{\mathrm{i}\langle k,w\rangle} & 0 \\ 0 & \mathrm{e}^{-\mathrm{i}\langle k,w\rangle} \end{pmatrix}\begin{pmatrix} \alpha & \beta \\ \gamma & \delta \end{pmatrix} \tag{66}$$

$$\begin{pmatrix} \alpha & \beta \\ \gamma & \delta \end{pmatrix}\begin{pmatrix} 0 & \mathrm{e}^{\mathrm{i}k_0a} \\ \mathrm{e}^{-\mathrm{i}k_0a} & 0 \end{pmatrix} = \begin{pmatrix} 0 & \mathrm{e}^{\mathrm{i}k_0a} \\ \mathrm{e}^{-\mathrm{i}k_0a} & 0 \end{pmatrix}\begin{pmatrix} \alpha & \beta \\ \gamma & \delta \end{pmatrix}. \tag{67}$$

From the last three relations it follows that

$$\beta\,\mathrm{e}^{\mathrm{i}\langle k,\tilde{c}\rangle} = \beta\,\mathrm{e}^{-\mathrm{i}\langle k,\tilde{c}\rangle} \qquad \beta\,\mathrm{e}^{\mathrm{i}\langle k,w\rangle} = \beta\,\mathrm{e}^{-\mathrm{i}\langle k,w\rangle} \qquad \alpha\,\mathrm{e}^{\mathrm{i}k_0a} = \delta\,\mathrm{e}^{\mathrm{i}k_0a} \tag{68}$$

$$\gamma\,\mathrm{e}^{\mathrm{i}\langle k,\tilde{c}\rangle} = \gamma\,\mathrm{e}^{-\mathrm{i}\langle k,\tilde{c}\rangle} \qquad \gamma\,\mathrm{e}^{\mathrm{i}\langle k,w\rangle} = \gamma\,\mathrm{e}^{-\mathrm{i}\langle k,w\rangle} \qquad \beta\,\mathrm{e}^{-\mathrm{i}k_0a} = \gamma\,\mathrm{e}^{\mathrm{i}k_0a}. \tag{69}$$

Since $\mathrm{e}^{\mathrm{i}k_0a} \neq 0$ we obtain $\alpha = \delta$. If either $\langle k, \tilde{c}\rangle \notin \mathbb{Z}\pi$ or $\langle k, w\rangle \notin \mathbb{Z}\pi$ then $\beta = \gamma = 0$ and the representation (64) is irreducible. If the numbers $\langle k, \tilde{c}\rangle$ and $\langle k, w\rangle$ belong to $\mathbb{Z}\pi$ then imposing the condition $P^2 = P$ we obtain

$$P = \frac{1}{2}\begin{pmatrix} 1 & \mathrm{e}^{\mathrm{i}k_0a} \\ \mathrm{e}^{-\mathrm{i}k_0a} & 1 \end{pmatrix} \qquad \text{or} \qquad P = \frac{1}{2}\begin{pmatrix} 1 & -\mathrm{e}^{\mathrm{i}k_0a} \\ -\mathrm{e}^{-\mathrm{i}k_0a} & 1 \end{pmatrix}. \tag{70}$$

These complementary projectors correspond to the decomposition

$$\mathcal{E}_k = \{\alpha(\psi_k + \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) \mid \alpha \in \mathbb{C}\} \oplus \{\alpha(\psi_k - \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) \mid \alpha \in \mathbb{C}\} \tag{71}$$

of $\mathcal{E}_k$ into direct sum of one-dimensional G_c-invariant subspaces, and

$$\begin{aligned} \varrho(\psi_k \pm \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) &= (-1)^m(\psi_k \pm \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) \\ \sigma(\psi_k \pm \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) &= (-1)^p(\psi_k \pm \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) \\ \tau(\psi_k \pm \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) &= \pm(\psi_k \pm \mathrm{e}^{-\mathrm{i}k_0a}\psi_{-k}) \end{aligned} \tag{72}$$

where $m, p \in \mathbb{Z}$ are such that $\langle k, \tilde{c}\rangle = m\pi$ and $\langle k, w\rangle = p\pi$. □

Relations (64) define for each k belonging to the set

$$\mathcal{B}_c^{\mathrm{irred}} = \{k \in \mathcal{B}_c\backslash\Lambda \mid \langle k, \tilde{c}\rangle \notin \mathbb{Z}\pi \text{ or } \langle k, w\rangle \notin \mathbb{Z}\pi\} \tag{73}$$

a two-dimensional irreducible representation $\mathcal{D}_c(k)$. Some of these representations are equivalent. Particularly, $\mathcal{D}_c(k) = \mathcal{D}_c(-k)$.

6. Clebsch–Gordan coefficients

Let $k, k' \in \mathcal{B}_c^{\text{irred}}$ be such that $k^+ = k + k'$ and $k^- = k - k'$ belong to $\mathcal{B}_c^{\text{irred}}$. The direct product of the representations $\mathcal{D}_c(k)$ and $\mathcal{D}_c(k')$ admits the decomposition

$$\mathcal{D}_c(k) \otimes \mathcal{D}_c(k') = \mathcal{D}_c(k^+) \oplus \mathcal{D}_c(k^-)$$

and the matrices corresponding to ϱ, σ and τ are

$$\begin{pmatrix} e^{i\langle k,\tilde{c}\rangle} & 0 \\ 0 & e^{-i\langle k,\tilde{c}\rangle} \end{pmatrix} \otimes \begin{pmatrix} e^{i\langle k',\tilde{c}\rangle} & 0 \\ 0 & e^{-i\langle k',\tilde{c}\rangle} \end{pmatrix} = \begin{pmatrix} e^{i\langle k^+,\tilde{c}\rangle} & 0 & 0 & 0 \\ 0 & e^{i\langle k^-,\tilde{c}\rangle} & 0 & 0 \\ 0 & 0 & e^{-i\langle k^-,\tilde{c}\rangle} & 0 \\ 0 & 0 & 0 & e^{-i\langle k^+,\tilde{c}\rangle} \end{pmatrix}$$

$$\begin{pmatrix} e^{i\langle k,w\rangle} & 0 \\ 0 & e^{-i\langle k,w\rangle} \end{pmatrix} \otimes \begin{pmatrix} e^{i\langle k',w\rangle} & 0 \\ 0 & e^{-i\langle k',w\rangle} \end{pmatrix} = \begin{pmatrix} e^{i\langle k^+,w\rangle} & 0 & 0 & 0 \\ 0 & e^{i\langle k^-,w\rangle} & 0 & 0 \\ 0 & 0 & e^{-i\langle k^-,w\rangle} & 0 \\ 0 & 0 & 0 & e^{-i\langle k^+,w\rangle} \end{pmatrix} \tag{74}$$

$$\begin{pmatrix} 0 & e^{ik_0a} \\ e^{-ik_0a} & 0 \end{pmatrix} \otimes \begin{pmatrix} 0 & e^{ik'_0a} \\ e^{-ik'_0a} & 0 \end{pmatrix} = \begin{pmatrix} 0 & 0 & 0 & e^{ik_0^+a} \\ 0 & 0 & e^{ik_0^-a} & 0 \\ 0 & e^{-ik_0^-a} & 0 & 0 \\ e^{-ik_0^+a} & 0 & 0 & 0 \end{pmatrix},$$

respectively. The unitary matrix

$$M = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 1 & 0 & 0 \end{pmatrix} \tag{75}$$

satisfies the relations

$$M^{-1} \begin{pmatrix} e^{i\langle k^+,\tilde{c}\rangle} & 0 & 0 & 0 \\ 0 & e^{i\langle k^-,\tilde{c}\rangle} & 0 & 0 \\ 0 & 0 & e^{-i\langle k^-,\tilde{c}\rangle} & 0 \\ 0 & 0 & 0 & e^{-i\langle k^+,\tilde{c}\rangle} \end{pmatrix} M = \begin{pmatrix} e^{i\langle k^+,\tilde{c}\rangle} & 0 & 0 & 0 \\ 0 & e^{-i\langle k^+,\tilde{c}\rangle} & 0 & 0 \\ 0 & 0 & e^{i\langle k^-,\tilde{c}\rangle} & 0 \\ 0 & 0 & 0 & e^{-i\langle k^-,\tilde{c}\rangle} \end{pmatrix}$$

$$M^{-1} \begin{pmatrix} e^{i\langle k^+,w\rangle} & 0 & 0 & 0 \\ 0 & e^{i\langle k^-,w\rangle} & 0 & 0 \\ 0 & 0 & e^{-i\langle k^-,w\rangle} & 0 \\ 0 & 0 & 0 & e^{-i\langle k^+,w\rangle} \end{pmatrix} M = \begin{pmatrix} e^{i\langle k^+,w\rangle} & 0 & 0 & 0 \\ 0 & e^{-i\langle k^+,w\rangle} & 0 & 0 \\ 0 & 0 & e^{i\langle k^-,w\rangle} & 0 \\ 0 & 0 & 0 & e^{-i\langle k^-,w\rangle} \end{pmatrix}$$

$$M^{-1} \begin{pmatrix} 0 & 0 & 0 & e^{ik_0^+a} \\ 0 & 0 & e^{ik_0^-a} & 0 \\ 0 & e^{-ik_0^-a} & 0 & 0 \\ e^{-ik_0^+a} & 0 & 0 & 0 \end{pmatrix} M = \begin{pmatrix} 0 & e^{ik_0^+a} & 0 & 0 \\ e^{-ik_0^+a} & 0 & 0 & 0 \\ 0 & 0 & 0 & e^{ik_0^-a} \\ 0 & 0 & e^{-ik_0^-a} & 0 \end{pmatrix}. \tag{76}$$

Therefore, the entries of M are Clebsch–Gordon coefficients [11] corresponding to the considered direct product. In this case, the only non-null coefficients are

$$(kk'11|k^+1) = (kk'12|k^-1) = (kk'21|k^-2) = (kk'22|k^+2) = 1. \tag{77}$$

More details concerning the Clebsch–Gordon coefficients and their applications in carbon nanotube physics can be found in the articles of Damnjanović *et al* [5, 7, 15].

7. Concluding remarks

The present paper can be regarded as a pure mathematical exercise. We have defined the factor sets $\mathcal{L}_c$, the groups G_c acting on $\mathcal{L}_c$ as groups of permutations, and we have studied certain representations of these groups defined on some spaces of functions $\psi : \mathcal{L}_c \longrightarrow \mathbb{C}$. We have proved that the groups G_c are isomorphic to the symmetry groups of single-wall carbon nanotubes, and the considered representations are directly related to some representations used in carbon nanotube physics.

The present paper can also be regarded as presenting an alternative mathematical model for carbon nanotubes. We think that this alternative approach offers some formal advantages, namely, certain calculations may be simpler in this approach than in the usual one.

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