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Graphene and non-Abelian quantization

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Abstract

In this paper, we employ a simple nonrelativistic model to describe the low energy excitation of graphene. The model is based on a deformation of the Heisenberg algebra which makes the commutator of momenta proportional to the pseudo-spin. We solve the Landau problem for the resulting Hamiltonian, which reduces in the large mass limit while keeping the Fermi velocity fixed, to the usual linear one employed to describe these excitations as massless Dirac fermions. This model, extended to negative mass, allows us to reproduce the leading terms in the low energy expansion of the dispersion relation for both nearest and next-to-nearest-neighbor interactions. Taking into account the contributions of both Dirac points, the resulting Hall conductivity, evaluated with a ζ -function approach, is consistent with the anomalous integer quantum Hall effect found in graphene. Moreover, when considered in first order perturbation theory, it is shown that the next-to-leading term in the interaction between nearest neighbor produces no modifications in the spectrum of the model while an electric field perpendicular to the magnetic field produces just a rigid shift of this spectrum.

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(Some figures may appear in colour only in the online journal)

1. Introduction

Several theoretical ideas have produced a fruitful exchange of interpretations and methods between fundamental physics and condensed matter theory, such as spontaneously broken symmetry [1–4] or renormalization group methods [5–9].

The recent experimental construction of graphene [10] opens a new connection between condensed matter and quantum field theory, since its low energy excitations can be represented as massless planar fermions and described by means of a (pseudo) relativistic theory.

Gaphene (see [11, 12] and references therein) is a two-dimensional, one-atom-thick, allotrope of carbon which has attracted great attention in the past few years. The carbon atoms are arranged on a honeycomb structure made out of hexagons, a structure with a great versatility. Carbon nanotubes, for example, can be obtained by rolling the graphene plane along a given direction and reconnecting the carbon bonds at the boundaries, giving rise to an essentially one-dimensional object. Also, the replacement of a hexagon by a pentagon in this lattice introduces a positive curvature defect; this allows us to wrap-up graphene to give fullerenes, molecules where carbon atoms are arranged on spherical structures. Note that the accumulation of graphene layers, weakly coupled by van der Waals forces, constitute the well-known graphite.

This peculiar material was theoretically predicted by Semenoff in 1984 [13] (see also [14]) and experimentally produced in the lab in 2004 [10].

The electronic properties of graphene are the result of the sp^2 hybridization between one s and two p orbitals, which leads to a trigonal planar lattice with a σ bond between carbon atoms separated by 1.42 Angstrom. This filled band gives the lattice its robustness. The third p orbital of the carbon atom, oriented perpendicularly to the plane of the graphene, gives rise to the π band through covalent bonds with neighboring atoms. Since this p orbital contributes with only one electron, the π band is half-filled in neutral graphene.

The low energy excitations of graphene accept a description as states of chiral massless Dirac fermions with a pseudo-relativistic linear dispersion relation, in which the speed of light *c* is replaced by the *Fermi velocity*, $v_F \approx 10^{-3}c$. Then, the Lagrangian describing these low energy states in the presence of an electromagnetic field is similar to that of QED for massless fermions, shearing therefore some of its peculiarities. In particular, when a magnetic field is applied perpendicularly to the plane of graphene, an anomalous integer quantum Hall effect [15, 16] takes place, which has been experimentally measured [17, 18].

It is not the goal of this paper to discuss the physics of graphene from first principles. Rather, we will consider a simple effective non-relativistic Hamiltonian, suggested by a particular deformation of the Heisenberg algebra (non-commutativity of momenta, consistent with the introduction of an external constant non-Abelian magnetic field), which could be useful to describe the low energy excitations produced by the dominant nearest and next-tonearest-neighbor interactions in graphene.

The non-commutativity of spacetime is an old idea [19], the first example of which was probably discussed by Landau in 1930 [20]. It has been revived in recent years within the context of string theory⁶ and since then, non-commutative field theories have attracted much attention in various fields such as mathematics, theoretical physics [22–25] and phenomenology [26].

The breaking of commutativity of the position operators and the representations of the algebra of the non-commutative spacetime coordinates has been studied in [27]. The non-commutativity in the momenta algebra we are interested in can be related to the deformation

⁶ The literature is vast in this field; for a general review, see [21].

quantization of Poissonian structures developed in [28] and considered as a kind of magnetic quantization [29, 30].

All this research has stimulated the construction of new models in quantum mechanics⁷, which has opened new routes to explore, for example in superconductivity [6]. Also a massless Dirac-like Hamiltonian in a generalized noncommutative space and its relation with graphene has been considered in [31].

Recently, some models based on a kind of nonstandard deformation of the Heisenberg algebra, which can be realized by shifting the dynamical variables with the spin, have been studied in [32, 33]. In the following, we will consider a similar deformation, but concentrated in the commutators among momenta, which can be interpreted as the introduction of a constant non-Abelian magnetic field [34].

In the following section, we present the model and derive the Hamiltonian. In section 3, we study the free case and in section 4, we introduce a constant magnetic field and solve the corresponding Landau problem. In section 5, we apply our results to describe the low energy states associated with the leading nearest and next-to-nearest interactions in graphene.

In section 6, we evaluate the associated Hall conductivity employing a ζ -function approach, finding that the result is consistent with the anomalous integer quantum Hall effect present in this material. In section 7, we are also able to show that the next-to-leading (quadratic) term in the nearest-neighbor interaction does not change the spectrum at first order in perturbation theory. In section 8, we comment on the case where crossed electric and magnetic fields are present and, finally, in section 9 we establish our conclusions.

At the end of the paper, an appendix is dedicated to the Lagrangian of the model and studies its symmetry, conserved current and the generating functional of Green's functions. It also comments on the weak field and gradient expansion of the generating functional, its relation with the Hall conductivity and the topological considerations involved.

2. Deformation of the Heisenberg algebra

We consider particles leaving on a plane whose dynamical variables satisfy the deformed Heisenberg algebra given by

$$[X_i, X_j] = 0, \quad [X_i, P_j] = \iota \delta_{ij}, \quad [P_i, P_j] = 2\iota \theta^2 \epsilon_{ij3} \sigma_3, \quad i, j = 1, 2,$$
(2.1)

where the momenta commutator is proportional to the *pseudospin* σ_3 and θ is a parameter with dimensions of momentum. (For convenience, we take $\hbar = 1$ and return to full units when necessary.)

Note that these particles are described by wavefunctions with two components, $\psi = {\varphi \choose \chi} \in \mathbf{L}_2(\mathbb{R}^2) \otimes \mathbb{C}^2$, and these operators have the structure of 2×2 matrices on \mathbb{C}^2 .

The deformed algebra in equation (2.1) can be realized by defining

$$X_i := x_i \otimes \mathbf{1}_2 \qquad P_i := p_i \otimes \mathbf{1}_2 + \theta \, \mathbf{1}_{\mathbf{L}_2} \otimes \sigma_i, \tag{2.2}$$

with x_i , i = 1, 2, the usual (commutative) coordinates on the plane and $p_i = -\iota \partial_i$, operators on $\mathbf{L}_2(\mathbb{R}^2)$, and σ_i , i = 1, 2, the first two Pauli matrices. We will also write $\boldsymbol{\sigma} := (\mathbf{1}_{\mathbf{L}_2} \otimes \sigma_1, \mathbf{1}_{\mathbf{L}_2} \otimes \sigma_2)$. For notational convenience, from now on we will avoid the explicit indication of the symbol \otimes , which can lead to no confusion.

Our aim is to consider the direct generalization of the Hamiltonian of a (nonrelativistic) particle of charge *e* and mass *m*, minimally coupled to an external ((2+1)-dimensional) electromagnetic field, $\{A_0, \mathbf{A} := (A_1, A_2)\}$, which is constructed by means of the replacements $x_i \rightarrow X_i$, $p_i \rightarrow P_i$.

 $^{^{7}}$ The literature is very extensive; a partial list can be found in the references of [33].

For the time being, we make $A_0 = 0$. Then,

$$H_0 = \frac{(\mathbf{p} - e\mathbf{A})^2}{2m} \quad \rightarrow \quad H_\theta = \frac{(\mathbf{p} - e\mathbf{A} + \theta\sigma)^2}{2m}, \tag{2.3}$$

where the electromagnetic field is taken in the Coulomb gauge, $\nabla \cdot \mathbf{A} = 0$.

The Hamiltonian can also be written as

$$H = \frac{(\mathbf{p} - e\mathbf{A})^2}{2m} + v_F \,\boldsymbol{\sigma} \cdot (\mathbf{p} - e\mathbf{A}), \qquad (2.4)$$

where we have defined the Fermi velocity

$$v_F := \frac{\theta}{m} \tag{2.5}$$

and subtracted the constant θ^2/m .

In the $m \to \infty$ limit, with fixed v_F , the resulting linear Hamiltonian is appropriate to describe the conducting effective particles in graphene around the *Fermi points* [11–13, 15], which justify our proposal. Note that this limit does not correspond to a small but rather to a large deformation of the commutator $[P_1, P_2]$.

Note also that the modification of the Hamiltonian in equation (2.4) can also be interpreted as the introduction of an SU(2) non-Abelian constant and uniform 'magnetic' field $\sim \theta^2$. Indeed, the SU(2) transformations relate both components of the wavefunctions, and the commutator of *covariant* derivatives gives

$$[p_1 - eA_1 + \theta\sigma_1, p_2 - eA_2 + \theta\sigma_2] = \iota e (\partial_1 A_2 - \partial_2 A_1) + 2\iota \theta^2 \sigma_3.$$
(2.6)

In this sense, if we take a constant magnetic field $B = (\partial_1 A_2 - \partial_2 A_1)$, the system we are considering is a kind of non-Abelian version of the Landau problem [34].

In the appendix, we describe the Lagrangian of this model, study its symmetry and conserved current, and discuss the relation of the asymptotic expansion of its generating functional with the Hall conductivity.

3. The free case

In this section, we consider the *free* case, with A = 0. The Hamiltonian reduces to

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$$H = \frac{\mathbf{p}^2}{2m} + v_F \,\boldsymbol{\sigma} \cdot \mathbf{p}. \tag{3.1}$$

We propose solutions of the form

$$\psi_{\mathbf{k}}(\mathbf{x}) = e^{\imath \mathbf{k} \cdot \mathbf{x}} \chi(\mathbf{k}), \qquad (3.2)$$

with $\chi(\mathbf{k}) \in \mathbb{C}^2$, which replaced in

$$[H - \mathcal{E}(\mathbf{k})] \psi_{\mathbf{k}}(\mathbf{x}) = 0 \tag{3.3}$$

leads to

$$\left\{\frac{\mathbf{k}^2}{2m} + v_F \,\boldsymbol{\sigma} \cdot \mathbf{k} - \mathcal{E}(\mathbf{k})\right\} \chi(\mathbf{k}) = 0.$$
(3.4)

Nontrivial solutions require

$$\det\left\{\frac{\mathbf{k}^2}{2m} + v_F \,\boldsymbol{\sigma} \cdot \mathbf{k} - \mathcal{E}(\mathbf{k})\right\} = 0,\tag{3.5}$$

which gives

$$v_F^2 \mathbf{k}^2 = \left[\mathcal{E}(\mathbf{k}) - \frac{\mathbf{k}^2}{2m} \right]^2.$$
(3.6)



Figure 1. The dispersion relations for the two branches of solutions for the free case.

Then, we get the following dispersion relation (approximately linear for small $|\mathbf{k}|$, see figure 1):

$$\mathcal{E}(\mathbf{k}) = \frac{\mathbf{k}^2}{2m} \pm v_F |\mathbf{k}| \tag{3.7}$$

which, replaced in equation (3.4) for $\mathbf{k} \neq 0$, shows that the pseudo-spinor $\chi(\mathbf{k})$ has definite pseudo-helicity

$$\left(\boldsymbol{\sigma}\cdot\hat{\mathbf{k}}\right)\chi_{\pm}(\mathbf{k}) = \pm\chi_{\pm}(\mathbf{k}). \tag{3.8}$$

On the other hand, for $\mathbf{k} = \mathbf{0}$, the two linearly independent vectors $\begin{pmatrix} 1 \\ 0 \end{pmatrix}$ and $\begin{pmatrix} 0 \\ 1 \end{pmatrix}$ are just constants solutions with vanishing eigenvalue.

The Hamiltonian in equation (3.1) commutes with the *effective angular momentum*, the generator of a U(1) symmetry,

$$J := -\iota \partial_{\varphi} + \frac{1}{2} \sigma_3, \quad [H, J] = 0.$$
(3.9)

Under a rotation on the plane, the wavefunction in equation (3.2) changes into

$$\mathcal{U}(\vartheta)\psi(\mathbf{x}) := e^{i\frac{\vartheta}{2}\sigma_3}\psi(R(\vartheta)^{-1}\mathbf{x}) = e^{i\mathbf{k}\cdot\left(R(\vartheta)^{-1}\mathbf{x}\right)}e^{i\frac{\vartheta}{2}\sigma_3}\chi(\mathbf{k}).$$
(3.10)

For $\vartheta = 2\pi$, we have $R(2\pi) = \mathbf{1}_3$ and $e^{i\pi\sigma_3} = -\mathbf{1}_2$. So we get

$$\mathcal{U}(2\pi)\psi(\mathbf{x}) = -\psi(\mathbf{x}). \tag{3.11}$$

Therefore, these particles can be considered as fermions [11].

4. Constant magnetic field perpendicular to the plane

In this section, we consider the electromagnetic vector potential of a constant magnetic field orthogonal to the plane

$$\mathbf{A} = Bx_1 \hat{\mathbf{e}}_2 \quad \Rightarrow \quad \partial_1 A_2 - \partial_2 A_1 = B \quad \text{and} \quad \nabla \cdot \mathbf{A} = 0. \tag{4.1}$$

In this case, the Hamiltonian can be written as

$$2mH = p_1^2 + (p_2 - eBx_1)^2 + 2mv_F\sigma_1p_1 + 2mv_F\sigma_2(p_2 - eBx_1), \qquad (4.2)$$

which clearly commutes with p_2 . This allows us to look for generalized eigenfunctions of the form

$$\psi(\mathbf{x}) = \frac{e^{ikx_2}}{\sqrt{2\pi}} \Phi(x_1), \tag{4.3}$$

where $\Phi = \begin{pmatrix} \varphi \\ \chi \end{pmatrix}$.

The eigenvalue equation for the Hamiltonian, $(H - \mathcal{E})\psi = 0$, reduces to the pair of coupled differential equations

$$\begin{cases} \left[p_1^2 + (eB)^2 \left(x_1 - \frac{k}{eB} \right)^2 \right] - \lambda \end{cases} \varphi = -2mv_F \left\{ p_1 + \iota eB \left(x_1 - \frac{k}{eB} \right) \right\} \chi,$$

$$\begin{cases} \left[p_1^2 + (eB)^2 \left(x_1 - \frac{k}{eB} \right)^2 \right] - \lambda \end{cases} \chi = -2mv_F \left\{ p_1 - \iota eB \left(x_1 - \frac{k}{eB} \right) \right\} \varphi,$$
(4.4)

with $\lambda = 2m\mathcal{E}$.

At this point, it is convenient to change the remaining variable x_1 in favor of $q := \sqrt{|eB|} \left(x_1 - \frac{k}{eB}\right)$. So $p_1 = \sqrt{|eB|}p$, with $p := -i\frac{\partial}{\partial q}$, and the system in equation (4.4) is written as

$$\{|eB|[p^{2} + q^{2}] - \lambda\}\varphi(q) = -2mv_{F}\sqrt{|eB|} \{p + \iota \operatorname{sgn}(eB) q\} \chi(q), \{|eB|[p^{2} + q^{2}] - \lambda\}\chi(q) = -2mv_{F}\sqrt{|eB|} \{p - \iota \operatorname{sgn}(eB) q\}\varphi(q).$$
(4.5)

For simplicity, in the following we will take eB > 0. The case eB < 0 can be obtained from the previous one by simply interchanging the components of the solutions, $\varphi(q) \leftrightarrow \chi(q)$, as can be easily seen from equation (4.5). So we consider

$$\{eB[p^2+q^2]-\lambda\}\begin{pmatrix}\varphi(q)\\\chi(q)\end{pmatrix} = -2mv_F\sqrt{eB}\{p\sigma_1-q\sigma_2\}\begin{pmatrix}\varphi(q)\\\chi(q)\end{pmatrix}.$$
(4.6)

The presence of twice the Hamiltonian of a harmonic oscillator of frequency 1 in the brackets on the left-hand side of equations (4.6), operator with eigenfunctions $\phi_n(q) = e^{-q^2/2}H_n(q)$ (H_n the Hermite polynomials) and eigenvalues 2n + 1, for n = 0, 1, ..., suggests that $\varphi(q) \sim \phi_n(q)$ and $\chi(q) \sim \phi_{n'}(q)$, for some n, n'. Moreover, since the Hermite functions satisfy

$$\begin{cases} \phi_n'(q) + q\phi_n(q) = 2n\phi_{n-1}(q) \\ \phi_n'(q) - q\phi_n(q) = -\phi_{n+1}(q), \end{cases}$$
(4.7)

one concludes that these solutions are of the form

$$\varphi(q) = c_1 \phi_{n+1}(y), \qquad \chi(q) = c_2 \phi_n(q)$$
(4.8)

for n = 0, 1, 2, ..., with $c_1, c_2 \in \mathbb{C}$. Replaced in equation (4.5), we get the homogeneous system of algebraic equations

$$M\begin{pmatrix}c_1\\c_2\end{pmatrix} = \begin{pmatrix}0\\0\end{pmatrix},\tag{4.9}$$

with

$$M = \begin{pmatrix} eB(2n+3) - \lambda & i2mv_F\sqrt{eB} \\ -i2mv_F\sqrt{eB}(2n+2) & eB(2n+1) - \lambda \end{pmatrix}.$$
(4.10)

The eigenvalues are determined by the condition

$$\det M = [2eB(n+1) - \lambda]^2 - (eB)^2 - 8m^2 v_F^2 eB(n+1) = 0,$$
(4.11)

which implies that

$$\mathcal{E}_{n,s} = \frac{\lambda_{n,s}}{2m} = (v_F \sqrt{eB}) \frac{1}{z} \left[n + 1 + \frac{s}{2} \sqrt{1 + 8z^2(n+1)} \right], \tag{4.12}$$

with $s = \pm 1$ and

$$z := \frac{mv_F}{\sqrt{eB}}.\tag{4.13}$$

Appropriately normalized, the generalized eigenfunctions are written as

$$\psi_{k,n,s}(\mathbf{x}) = \frac{e^{\imath k x_2}}{\sqrt{2\pi}} K_{n,s} e^{-\frac{q^2}{2}} \begin{pmatrix} H_{n+1}(q) \\ \frac{\imath}{2z} [1 - s\sqrt{1 + 8z^2(n+1)}] H_n(q) \end{pmatrix},$$
(4.14)

where

$$K_{n,s} = \sqrt[4]{eB} \frac{2^{-\frac{n}{2}-1} \sqrt{\left(\frac{s}{2} + \sqrt{1+8z^2(n+1)}\right)^2 - \frac{1}{4}}}{\sqrt[4]{\pi} \sqrt{(n+1)!} \sqrt{1+8z^2(n+1)}}$$
(4.15)

and

$$q = \sqrt{eB} \left(x_1 - \frac{k}{eB} \right). \tag{4.16}$$

Indeed, taking into account the orthogonality relations for the Hermite functions, it can be easily verified that

$$(\psi_{k,n,s},\psi_{k',n',s'}) = \delta_{n,n'}\delta_{s,s'}\delta(k-k').$$
(4.17)

Finally, note that there is another solution Φ_0 of equation (4.6) whose components are given by

$$\varphi(q) = \left(\frac{eB}{\pi}\right)^{1/4} \phi_0(q), \qquad \chi(q) = 0.$$
 (4.18)

Indeed, since $H_0(q) = 1$, we have

$$(p - \iota q)\phi_0(q) = -\iota(\partial_q + q)e^{-q^2/2} = 0$$
(4.19)

and equation (4.5) reduces to

$$\left\{\frac{eB}{m}\frac{1}{2}[p^2+q^2] - \mathcal{E}_0\right\}\phi_0(q) = \left\{\frac{eB}{m}\frac{1}{2} - \mathcal{E}_0\right\}\phi_0(q) = 0.$$
(4.20)

This implies that $\mathcal{E}_0 = eB/2m$, which is independent of v_F .

One can also verify that

$$\psi_{k,0}(\mathbf{x}) = \frac{\mathrm{e}^{\imath k x_2}}{\sqrt{2\pi}} \left(\frac{eB}{\pi}\right)^{1/4} \begin{pmatrix} \phi_0(q)\\ 0 \end{pmatrix} \tag{4.21}$$

satisfy

$$(\psi_{k,n,s},\psi_{k',0})=0, \quad (\psi_{k,0},\psi_{k',0})=\delta(k-k').$$
(4.22)

As mentioned in section 2, these results can be interpreted as the solution of a non-Abelian version of the Landau problem, in which we have also introduced a constant and uniform non-Abelian magnetic field. Moreover, in the $m \to \infty$ limit, these eigenfunctions are similar to the solutions found [35] for the Dirac equation in a constant magnetic background.

Note that one can take appropriate linear combinations of the generalized eigenfunctions in equations (4.14) and (4.21) so as to construct a manifestly complete set of generalized vectors in our space,

$$\left\{\frac{\mathrm{e}^{\mathrm{i}kx_2}}{\sqrt{2\pi}}\phi_n(q)\begin{pmatrix}1\\0\end{pmatrix}, \frac{\mathrm{e}^{\mathrm{i}kx_2}}{\sqrt{2\pi}}\phi_n(q)\begin{pmatrix}0\\1\end{pmatrix}; k \in \mathbb{R}, n = 0, 1, 2, \ldots\right\},\tag{4.23}$$



Figure 2. The energy levels $\mathcal{E}_{n,s}$ for negative mass, in units of $v_F \sqrt{eB}$ and for w = 100.

where $\phi_n(q)$ are the Hermite functions. Then, the set of generalized eigenvectors of *H* that we found is also complete, which ensures that in our analysis we got the whole spectrum of the Hamiltonian.

In the following, we consider the large-z (large-m) limit we are interested in⁸.

For $z \gg 1$, the energies are given by

$$\mathcal{E}_{n,s} = v_F \sqrt{eB} \left\{ s \sqrt{2(n+1)} + \frac{(n+1)}{z} + O(z^{-2}) \right\}$$
(4.27)

and the eigenfunctions reduce to

$$\psi_{k,n,s}(\mathbf{x}) = \frac{e^{ikx_2}}{\sqrt{2\pi}} \frac{2^{-\frac{n}{2}-1} e^{-\frac{q^2}{2}}}{\sqrt[4]{\pi}\sqrt{(n+1)!}} \left(\frac{H_{n+1}(q) + O(z^{-1})}{-\frac{is}{2}\sqrt{8(n+1)}} H_n(q) + O(z^{-1}) \right).$$
(4.28)

Negative mass. Since we are interested in the description of low energy states and the construction of the solutions of equation (4.4) is independent of the sign of *m*, we can explore the behavior of the system for m < 0. Indeed, it is safe to change the sign of *m* in the solutions, which corresponds to the replacement $z \rightarrow -w$, with $w = |m|v_F/\sqrt{eB} > 0$, in the last line in equation (4.12) maintaining $v_F > 0$. We get (see figure 2)

⁸ For $z \ll 1$, we have equally spaced levels in each branch, with slightly different slopes,

$$\mathcal{E}_{n,s} = \frac{eB}{m} \left\{ \left(n + 1 + \frac{s}{2} \right) + 2sz^2(n+1) + O(z^4) \right\}.$$
(4.24)

For the eigenfunctions, we get

$$\psi_{k,n,+1}(\mathbf{x}) = \frac{e^{ikx_2}}{\sqrt{2\pi}} \frac{2^{-\frac{n}{2} - \frac{1}{2}}}{\sqrt[4]{\pi}\sqrt{(n+1)!}} e^{-\frac{q^2}{2}} \begin{pmatrix} [1 - z^2(n+1) + O(z^4)]H_{n+1}(q) \\ [-2iz(n+1) + 6iz^3(n+1)^2 + O(z^4)]H_n(q) \end{pmatrix}$$
(4.25)

and

$$\psi_{k,n,-1}(\mathbf{x}) = \frac{e^{ikx_2}}{\sqrt{2\pi}} \frac{2^{-n/2}}{\sqrt[4]{\pi}\sqrt{n!}} e^{-\frac{q^2}{2}} \begin{pmatrix} [z - 3z^3(n+1) + O(z^5)]H_{n+1}(q) \\ \iota \begin{bmatrix} 1 - z^2(n+1) + \frac{11}{2}z^4(n+1)^2 + O(z^5) \end{bmatrix} H_n(q) \end{pmatrix}.$$
(4.26)

The $z \rightarrow 0$ -limit is consistent with the usual Landau level problem.

$$\mathcal{E}_{n,s} = -(v_F \sqrt{eB}) \frac{1}{w} \left[n + 1 + \frac{s}{2} \sqrt{1 + 8w^2(n+1)} \right]$$

= $(v_F \sqrt{eB}) \left\{ -s \sqrt{2(n+1)} - \frac{n+1}{w} + O(w^{-2}) \right\}.$ (4.29)

On the other hand⁹, $\mathcal{E}_0 = -eB/2|m| < 0$.

5. The relation with graphene

The structure of graphene [11] can be seen as a triangular lattice with a basis of two atoms per unit cell. The lattice vectors are

$$\mathbf{a}_1 = \frac{3}{2}a\begin{pmatrix}1\\\frac{1}{\sqrt{3}}\end{pmatrix}, \quad \mathbf{a}_2 = \frac{3}{2}a\begin{pmatrix}1\\-\frac{1}{\sqrt{3}}\end{pmatrix}, \tag{5.1}$$

where $a \approx 1.42$ Å is the lattice constant (distance between nearest-neighbor carbon atoms). The vectors characterizing the reciprocal lattice (whose elementary cell is the Brillouin zone) are given by

$$\mathbf{b}_1 = \frac{2\pi}{3a} \begin{pmatrix} 1\\\sqrt{3} \end{pmatrix}, \qquad \mathbf{b}_2 = \frac{2\pi}{3a} \begin{pmatrix} 1\\-\sqrt{3} \end{pmatrix}. \tag{5.2}$$

The superposed triangular lattices form a hexagonal honeycomb array of carbon atoms where the nearest-neighbor vectors are

$$\boldsymbol{\delta}_1 = \frac{a}{2} \begin{pmatrix} 1\\\sqrt{3} \end{pmatrix}, \qquad \boldsymbol{\delta}_2 = \frac{a}{2} \begin{pmatrix} 1\\-\sqrt{3} \end{pmatrix}, \qquad \boldsymbol{\delta}_3 = a \begin{pmatrix} -1\\0 \end{pmatrix}. \tag{5.3}$$

The tight-binding model for graphene, where it is assumed that electrons can only hop to both nearest $(\langle i, j \rangle)$ and next-to-nearest $(\langle \langle i, j \rangle)$ neighbor atoms, is described by the Hamiltonian [11]

$$H = -t \sum_{\langle i,j \rangle} \sum_{s=\pm} \left(a_{i,s}^{\dagger} b_{j,s} + \text{h.c.} \right) - t' \sum_{\langle \langle i,j \rangle \rangle} \sum_{s=\pm} \left(a_{i,s}^{\dagger} a_{j,s} + b_{i,s}^{\dagger} b_{j,s} + \text{h.c.} \right),$$
(5.4)

where *t* and *t'* are the hopping energies, $a_{i,s}^{\dagger}$ and $a_{i,s}$ are the creation and annihilation operators of electrons in the site *i* with spin *s* belonging to the triangular sublattice *A*, and similarly for the sublattice *B*. By Fourier transforming these operators and diagonalizing the resulting expression, one gets the band structure (dispersion relations) [11]

$$E_{\pm}(\mathbf{k}) = \pm t \sqrt{f(\mathbf{k})} - t' \left[f(\mathbf{k}) - 3 \right]$$
(5.5)

with

$$f(\mathbf{k}) = 3 + 4\cos\left(\frac{3k_1a}{2}\right)\cos\left(\frac{\sqrt{3}}{2}k_2a\right) + 2\cos(\sqrt{3}k_2a) \ge 0.$$
(5.6)

The minima of $f(\mathbf{k})$ in the Brillouin zone correspond to the *Dirac points*

$$\mathbf{K} = \frac{2\pi}{3a} \begin{pmatrix} 1\\ \frac{1}{\sqrt{3}} \end{pmatrix}, \quad \mathbf{K}' = \frac{2\pi}{3a} \begin{pmatrix} 1\\ \frac{-1}{\sqrt{3}} \end{pmatrix}, \tag{5.7}$$

where $f(\mathbf{K}) = 0 = f(\mathbf{K}')$. Then, the bands corresponding to each sign in the dispersion relations, equation (5.5), touch each other at the Dirac points. Note that these two independent Dirac points have been chosen so that they are related by the reflection of \mathbf{k} about the k_1 -axis $(k_2 \rightarrow -k_2)$ [11].

⁹ Note that $\mathcal{E}_0 \to 0$ for $m \to \infty$. This is consistent with the well-known existence of zero modes in the spectrum of massless Dirac fermions coupled to gauge fields, discussed in different contexts in [36], [37] and [38], for example.

A series expansion of $E_{\pm}(\mathbf{k})$ around $\mathbf{k} = \mathbf{K}$ leads to

$$E_{s}(\mathbf{K} + \mathbf{k}) = st \left[\frac{3}{2}a|\mathbf{k}| - \frac{3}{8}a^{2}\mathbf{k}^{2}\sin(3\theta)\right] + t' \left[-\frac{9}{4}a^{2}\mathbf{k}^{2} + 3\right] + O(|\mathbf{k}|^{3}),$$
(5.8)

where $tan(\theta) = k_2/k_1$ and $s = \pm 1$.

The expansion around the second Fermi point, **K**', leads to the same expression with **k** reflected about the k_1 -axis. This corresponds to the change $\theta \rightarrow -\theta$, which changes the sign of the second term in the first bracket on the right-hand side of equation (5.8).

We now turn to the comparison with the model developed in the previous sections. From equations (5.8) and (3.7), it is seen that we can identify $v_F \equiv \frac{3}{2} at$. Moreover, the next-to-nearest-neighbor contribution can be represented in our (free) model by the *mass* term through the identification $-\frac{9}{4}t'a^2 \equiv \frac{1}{2m}$, up to a rigid displacement of the spectrum in an energy 3t' (which can be subtracted to restore the particle–hole symmetry at linear order in $|\mathbf{k}|$). This means that we must consider a *negative mass* in our model, $m = -2/(9t'a^2) < 0.^{10}$

On the other hand, the second (quadratic) term in the nearest-neighbor contribution depends on the direction of \mathbf{k} . We cannot reproduce this behavior within the framework of our model since it is rotationally invariant. Rather, we must treat it as a perturbation on the solutions we found for our model. This will be considered later.

Then, we will assume that the effective Hamiltonian for the low energy states of this system around **K** in the presence of an electromagnetic field, $H_{\mathbf{K}}$, is obtained though minimal coupling and coincides with the one given in equation (2.4).

Since the application of our model to graphene requires $w \approx 10^3$ (see footnote 10), the large w expansion in equation (4.29) gives a good description of the self-energies. So we get for quasi-particle states

$$\mathcal{E}_{n,-} = (v_F \sqrt{eB}) \left\{ \sqrt{2(n+1)} - \frac{n+1}{w} + O(w^{-2}) \right\},$$
(5.12)

and for hole states

$$\mathcal{E}_{n,+} = (v_F \sqrt{eB}) \left\{ -\sqrt{2(n+1)} - \frac{n+1}{w} + O(w^{-2}) \right\}.$$
(5.13)

Moreover, there is an additional (hole) state at energy $\mathcal{E}_0 = -\frac{v_F \sqrt{eB}}{2w} < 0$, slightly below the zero energy level (see equation (4.20)). Indeed, the difference in energy for the two first hole states is $v_F \sqrt{eB}(\sqrt{2} + O(w^{-1}))$.¹¹

¹⁰ According to the values of these parameters reported in [11], we have

$$a = 1.42 \text{ Å}, \quad t = 2.8 \text{ eV}, \quad t' = 0.1 \text{ eV}.$$
 (5.9)

Then,

$$m_F = 600 \text{ km s}^{-1}, \quad mc^2 = -4.3 \times 10^6 \text{ eV},$$
 (5.10)

or, in natural units $(v_F/c \rightarrow v_F, mc/\hbar \rightarrow m)$,

$$v_F = 2 \times 10^{-3}, \quad m = -8.53 \times 10^{13} \text{ m}^{-1}.$$
 (5.11)

This is a rather large value for the (negative) mass of the quasiparticles, since $|m| \simeq 32.9 m_e$. For example, for an external magnetic field B = 10 Tesla, we get $w = \frac{|m|_c}{c} \frac{v_F}{c} \sqrt{\frac{\hbar}{eB}} = 1385$.

¹¹ For $w = 10^3$ and B = 10 Tesla, we have in full units

$$v_F \sqrt{eB} \to \hbar c \, \frac{v_F}{c} \sqrt{\frac{eB}{\hbar}} = 4.87 \times 10^{-2} \, \mathrm{eV}.$$
 (5.14)

This leads to $\mathcal{E}_0 = -2.43 \times 10^{-5} \text{ eV}$ and $\mathcal{E}_{0,+} = -6.89 \times 10^{-2} \text{ eV}$.

5.1. The second Fermi point contribution

As previously mentioned, the dispersion relations near the Fermi points are related by the reflection of momenta about the k_1 -axis. Then, for the linear terms in the effective free Hamiltonian describing the states around the second Fermi point, we can write [11]

$$v_F (\sigma_1 p_1 - \sigma_2 p_2) = v_F \boldsymbol{\sigma}^* \cdot \mathbf{p} = -\sigma_2 (v_F \boldsymbol{\sigma} \cdot \mathbf{p}) \sigma_2.$$
(5.15)

Accordingly, we will assume that the effective Hamiltonian describing the quasi-particles near the second Fermi point in our model, $H_{\mathbf{K}'}$, can be obtained from $H_{\mathbf{K}}$ through the transformation

$$H_{\mathbf{K}'} = -\sigma_2 H_{\mathbf{K}} \sigma_2 = -\frac{(\mathbf{p} - e\mathbf{A})^2}{2m} + v_F \,\boldsymbol{\sigma}^* \cdot (\mathbf{p} - e\mathbf{A}). \tag{5.16}$$

Note that both $H_{\mathbf{K}}$ and $H_{\mathbf{K}'}$ are left invariants under time-reversal times parity transformations, \mathcal{TP} (see the discussion in [12] about the symmetries of this system). Indeed [12],

$$(\mathcal{TP})H_{\mathbf{K}}(\mathcal{TP})^{\dagger} = \sigma_{1}H_{\mathbf{K}}^{*}\sigma_{1} = \sigma_{1}\left\{\frac{(\mathbf{p}-e\mathbf{A})^{2}}{2m} + v_{F}\boldsymbol{\sigma}^{*}\cdot(\mathbf{p}-e\mathbf{A})\right\}\sigma_{1} = H_{\mathbf{K}}$$
(5.17)

and similarly

$$(\mathcal{TP}) H_{\mathbf{K}'} (\mathcal{TP})^{\dagger} = -\sigma_2^* (\mathcal{TP}) H_{\mathbf{K}} (\mathcal{TP})^{\dagger} \sigma_2^*$$

= $-\sigma_2 H_{\mathbf{K}} \sigma_2 = H_{\mathbf{K}'}.$ (5.18)

Evidently, the spectrum of $H_{\mathbf{K}'}$ is obtained from that of $H_{\mathbf{K}}$ by a reflection about the origin and its (generalized) eigenfunctions are just σ_2 times those of $H_{\mathbf{K}}$. This interchanges quasi-particles with holes. Indeed,

$$H_{\mathbf{K}'}(\sigma_2\psi_{k,n,s}) = -\sigma_2(H_{\mathbf{K}}\psi_{k,n,s}) = -\mathcal{E}_{n,s}(\sigma_2\psi_{k,n,s}).$$
(5.19)

Moreover,

1

$$H_{\mathbf{K}'}(\sigma_2\psi_0) = -\sigma_2 (H_{\mathbf{K}}\psi_0) = -\mathcal{E}_0(\sigma_2\psi_0), \qquad (5.20)$$

with an energy $\mathcal{E}'_0 = -\mathcal{E}_0 = \frac{v_F \sqrt{eB}}{2w} > 0$. This state corresponds to a quasi-particle of energy slightly above zero.

Therefore, taking into account the eigenstates of both $H_{\mathbf{K}}$ and $H_{\mathbf{K}'}$, we get, for quasiparticles and for holes, an almost doubly degenerate spectrum, except for one state of quasiparticle and one state of hole with energies near zero. For positive energy states, we have

$$\mathcal{E}_{n,-} = (v_F \sqrt{eB}) \left\{ \sqrt{2(n+1)} - \frac{n+1}{w} + O(w^{-2}) \right\}, -\mathcal{E}_{n,+} = (v_F \sqrt{eB}) \left\{ \sqrt{2(n+1)} + \frac{n+1}{w} + O(w^{-2}) \right\},$$
(5.21)

where the gap between contiguous states is

$$\Delta \mathcal{E}_n = 2(v_F \sqrt{eB}) \left\{ \frac{n+1}{w} + O(w^{-2}) \right\}.$$
(5.22)

As we discuss in the next section, this spectrum reproduces qualitatively the anomalous integer quantum Hall effect found in graphene [15–18], which shows a nonvanishing Hall conductivity for a small (positive or negative) Fermi level.

6. The Hall conductivity

As discussed in the appendix, the Hall conductivity has a topological character and can be calculated from the weak field and gradient expansion of the effective action of the system. However, since we know the exact eigenvalues of energy of our model, we will employ a more direct evaluation method, based on the relation established between the conserved current and the external electromagnetic field, equations (A.13) and (A.14).

For our purposes, it will be sufficient to consider the mean value of the density $j_0(x)$. Then, as in the previous sections, we will take $\mathbf{E} = 0$ and $B \neq 0$ perpendicular to the plane of the system.

Let us start by writing down the partition function of the particles around one Fermi point at inverse temperature β and chemical potential μ . It can be obtained through a Wick rotation of the generating functional Z[A] (see equation (A.8)) to the 2+1 Euclidean space by means of the replacements $x^0 = t \rightarrow -\iota\tau$, $A_0 \rightarrow \iota A_3$, maintaining the other coordinates and components of the gauge field unaltered.

Since, in our case, $A_0 = 0 = A_1$, $A_2 = Bx_1$, from equations (A.1) and (A.8), this Wick rotation leads to the partition function

$$\mathcal{Z}(\beta,\mu,B) := \int \mathcal{D}\psi^{\dagger} \mathcal{D}\psi \,\mathrm{e}^{\int_{0}^{\beta} \mathrm{d}t \int \mathrm{d}^{2}x \,\psi^{\dagger} \{-\frac{\partial}{\partial\tau} + \mu - H\}\psi},\tag{6.1}$$

where *H* is $H_{\mathbf{K}}$ (or $H_{\mathbf{K}'}$) and the functional integral is performed on the set of configurations of the fermionic field which satisfy anti-periodic boundary conditions on $[0, \beta]$. Then,

$$\frac{\partial \log \mathcal{Z}}{\partial \mu}(\beta, \mu, B) = \int_0^\beta dt \int d^2 x \langle \psi^{\dagger} \psi \rangle$$
$$= \beta \int d^2 x \, \frac{1}{e} J_0(\beta, \mu, B) = \beta \int d^2 x \, \frac{\sigma_{xy}}{e} B, \tag{6.2}$$

where we used equation (A.14).

Our goal is now to evaluate the partition function as the functional determinant

$$\mathcal{Z}(\beta,\mu,B) = \operatorname{Det}(D), \tag{6.3}$$

where $D = -\frac{\partial}{\partial \tau} + \mu - H$ is a differential operator defined on a domain of anti-periodic functions of $\tau \in [0, \beta]$. Even though *D* is not symmetric, since *H* does not depend on τ this operator has a complete set of orthogonal generalized eigenfunctions constructed as

$$\Psi_{l,k,n,s} = \frac{1}{\sqrt{\beta}} e^{-\iota\omega_l \tau} \psi_{k,n,s}, \qquad \Psi_{l,k,0} = \frac{1}{\sqrt{\beta}} e^{-\iota\omega_l \tau} \psi_{k,0}, \tag{6.4}$$

with $l \in \mathbb{Z}, k \in \mathbb{R}, n = 0, 1, 2, ..., s = \pm 1$, and where the Matsubara frequencies

$$\omega_l = \frac{2\pi}{\beta} \left(l + \frac{1}{2} \right), \quad \text{with} \quad \omega_l = -\omega_{-l-1}.$$
 (6.5)

Then, the eigenvalues of D are given by

$$D\Psi_{l,k,n,s} = \lambda_{l,n,s}\Psi_{l,k,n,s}, \qquad \lambda_{l,n,s} = \iota\omega_l + \mu - \mathcal{E}_{n,s}, D\Psi_{l,k,0} = \lambda_{l,0}\Psi_{l,k,0}, \qquad \lambda_{l,0} = \iota\omega_l + \mu - \mathcal{E}_0,$$
(6.6)

where $\mathcal{E}_{n,s}$ and \mathcal{E}_0 are the eigenvalues of *H*, studied in the previous sections.

Note that $\lambda_{l,n,s}$ and $\lambda_{l,0}$ are independent of k. Therefore, in evaluating Det(D) we will forget about the index k and, at the end, take into account the degeneracy it introduces *per unit area*, given by the number of flux quanta per unit area [39, 40], $\Delta = eB/2\pi$ (or $\Delta = eB/h$ in full units).

We define Det(D) as the (gauge invariant) ζ -function determinant [41, 42],

$$\log \operatorname{Det}(D) := -\frac{\mathrm{d}}{\mathrm{d}u} \operatorname{Tr}\left\{ \left(\frac{D}{\Lambda} \right)^{-u} \right\} \bigg|_{u \to 0}, \tag{6.7}$$

where Λ is an arbitrary mass scale, the trace is evaluated for $\Re(u)$ sufficiently large and $u \to 0$ stands for the analytic continuation of its derivative to a neighborhood of u = 0. In our case,

$$\log \mathcal{Z}(\beta, \mu, B) = -\frac{\mathrm{d}}{\mathrm{d}u} \left\{ \sum_{l,n,s} \left(\frac{\lambda_{l,n,s}}{\Lambda} \right)^{-u} + \sum_{l} \left(\frac{\lambda_{l,0}}{\Lambda} \right)^{-u} \right\} \bigg|_{u \to 0}.$$
 (6.8)

Let us first consider the contribution of the second term on the right-hand side and evaluate

$$\sum_{l=-\infty}^{\infty} \left(\frac{\lambda_{l,0}}{\Lambda}\right)^{-u} = \sum_{l=0}^{\infty} \left(\frac{\iota\omega_l + \mu - \mathcal{E}_0}{\Lambda}\right)^{-u} + \sum_{l=0}^{\infty} \left(\frac{-\iota\omega_l + \mu - \mathcal{E}_0}{\Lambda}\right)^{-u},\tag{6.9}$$

where we used the properties of the Matsubara frequencies stated in equation (6.5). We can also write

$$\sum_{l=-\infty}^{\infty} \left(\frac{\lambda_{l,0}}{\Lambda}\right)^{-u} = \left(\frac{2\pi}{\beta\Lambda}\right)^{-u} \left\{ \sum_{l=0}^{\infty} \left[\iota \left(l + \frac{1}{2}\right) + \frac{\beta}{2\pi} \left(\mu - \mathcal{E}_{0}\right) \right]^{-u} + \sum_{l=0}^{\infty} \left[-\iota \left(l + \frac{1}{2}\right) + \frac{\beta}{2\pi} \left(\mu - \mathcal{E}_{0}\right) \right]^{-u} \right\}.$$
(6.10)

These series can be expressed in terms of the Hurwitz ζ -function [43, 44]. Taking into account that this function has branch cut discontinuities in the complex plane of its second argument running from 0 to $-\infty$, we get

$$\sum_{l=-\infty}^{\infty} \left(\frac{\lambda_{l,0}}{\Lambda}\right)^{-u} = \left(\frac{2\pi}{\beta\Lambda}\right)^{-u} \left\{ e^{-i\frac{\pi}{2}u} \zeta \left(u, \frac{1}{2} + e^{-i\frac{\pi}{2}\operatorname{sign}(\mu - \mathcal{E}_0)} \frac{\beta}{2\pi} |\mu - \mathcal{E}_0| \right) + e^{i\frac{\pi}{2}u} \zeta \left(u, \frac{1}{2} + e^{i\frac{\pi}{2}\operatorname{sign}(\mu - \mathcal{E}_0)} \frac{\beta}{2\pi} |\mu - \mathcal{E}_0| \right) \right\}.$$
(6.11)

The expression in braces has a vanishing analytic continuation at u = 0. As a consequence, the derivative of the right-hand side at u = 0 does not depend on the arbitrary scale A. Then, the contribution of this Landau level to log DetD per unit area and at low temperature is given by [43, 44]

$$\Delta \left\{ -\frac{\mathrm{d}}{\mathrm{d}u} \sum_{l=-\infty}^{\infty} \left(\frac{\lambda_{l,0}}{\Lambda} \right)^{-u} \Big|_{u \to 0} \right\} = \frac{eB}{2\pi} \left\{ \beta \left(\mathcal{E}_0 - \mu \right) \Theta \left(\mathcal{E}_0 - \mu \right) + o\left(\frac{1}{\beta \left| \mu - \mathcal{E}_0 \right|} \right) \right\}.$$
(6.12)

`

Taking into account that we are interested in the mean number of particles with respect to the neutral material (i.e. with all the Landau levels with negative energy filled, which corresponds to $\mu = 0$), from equation (6.2), we finally get for the contribution of the Landau level with energy \mathcal{E}_0 to the Hall conductivity at zero temperature

$$\frac{B}{e} \sigma_{xy}\Big|_{\mathcal{E}_{0}} = \frac{Be}{2\pi} \left\{ \frac{\partial}{\partial \mu} \left[(\mathcal{E}_{0} - \mu) \Theta (\mathcal{E}_{0} - \mu) \right] - \frac{\partial}{\partial \mu} \left[(\mathcal{E}_{0} - \mu) \Theta (\mathcal{E}_{0} - \mu) \right] \Big|_{\mu=0} \right\}$$

$$= \frac{Be}{2\pi} \left\{ -\Theta (\mathcal{E}_{0} - \mu) + \Theta (\mathcal{E}_{0}) \right\}$$

$$= \frac{Be}{2\pi} \left\{ \Theta (\mathcal{E}_{0}) \Theta (\mu - \mathcal{E}_{0}) - \Theta (-\mathcal{E}_{0}) \Theta (\mathcal{E}_{0} - \mu) \right\}.$$
(6.13)



Figure 3. The Hall conductivity for a realistic value of $w = 10^3$.

Therefore, for positive \mathcal{E}_0 , we get a contribution $(+\frac{e^2}{2\pi})$ to σ_{xy} if $\mu > \mathcal{E}_0$ and zero otherwise. On the other hand, for negative \mathcal{E}_0 , we get a contribution $(-\frac{e^2}{2\pi})$ if $\mu < \mathcal{E}_0$ and zero otherwise.

At this point, it is clear that a similar result is obtained for any other Landau level. Indeed, the same considerations can be made for any other energy eigenvalue \mathcal{E} . Then, for a given chemical potential μ , there is only a finite number of nonvanishing contributions and the Hall conductivity reduces to

$$\sigma_{xy}\Big|_{H_{\mathbf{K}}} = \frac{e^2}{2\pi} \left\{ \Theta(\mu) \left(\sum_{0 < \mathcal{E} < \mu} 1 \right) - \Theta(-\mu) \left(\sum_{\mu < \mathcal{E} < 0} 1 \right) \right\}, \tag{6.14}$$

where the sums run over the energy levels. Indeed, this is a rather general result for the Landau problem of a fermionic system.

For the model of interest in this paper, the spectrum to be considered is the union of those of $H_{\mathbf{K}}$ and $H_{\mathbf{K}'}$, $\mathcal{E} \in \text{Spec}H_{\mathbf{K}} \bigcup \text{Spec}H_{\mathbf{K}'}$, which takes into account the states around both Fermi points. But they must be taken with an additional degeneracy corresponding to the two polarizations of the electron spin, which has played no role up to now in this analysis.

In full units, we get

$$\sigma_{xy} = \frac{2e^2}{h} \left\{ \Theta(\mu) \left(\sum_{0 < \mathcal{E} < \mu} 1 \right) - \Theta(-\mu) (\sum_{\mu < \mathcal{E} < 0} 1) \right\}.$$
(6.15)

Note that

$$\left|\frac{h}{4e^2}\,\sigma_{xy}\right| = \frac{1}{2}\tag{6.16}$$

for small $|\mu|$ (but $|\mu| > |\mathcal{E}_0| = \frac{v_F \sqrt{eB}}{2w}$). This is the characteristic of the *anomalous integer quantum Hall effect* found in graphene [15–18], which shows a nonvanishing Hall conductivity for small (positive or negative) Fermi level.

In figure 3, the Hall conductivity of this model as a function of μ (equal to the Fermi level ϵ_F at zero temperature) has been plotted for $w = 10^3$. Note that it is not possible to appreciate the structure of each step as a double step (due to the small degeneracy breaking remarked in equations (5.21) and (5.22)) with such a realistic value of w (see figure 4). Also note that the



Figure 4. The structure of the step in the Hall conductivity from $\frac{1}{2}$ to $\frac{3}{2}$, with $w = 10^3$.

Hall conductivity of our model vanishes in a small neighborhood of the origin for $|\mu| < \frac{v_F \sqrt{eB}}{2w}$.

7. Perturbation with the quadratic nearest-neighbor interaction term

Note that the quadratic term in the nearest-neighbor dispersion relation in equation (5.8) can be obtained from the solutions of the free problem, equations (3.2) and (3.8), through the (rotational symmetry breaking) perturbation

$$\Delta H = -\frac{a}{4} v_F p_2 [4p_1^2 - \mathbf{p}^2] (\mathbf{p}^2)^{-1} (\boldsymbol{\sigma} \cdot \mathbf{p}).$$
(7.1)

Indeed,

$$\Delta H \,\psi_{\mathbf{k}}(\mathbf{x}) = -\frac{a}{4} \,v_F \mathrm{e}^{i\mathbf{k}\cdot\mathbf{x}} \,k_2 \left[4k_1^2 - \mathbf{k}^2\right] (\mathbf{k}^2)^{-1/2} \left(\boldsymbol{\sigma}\cdot\hat{\mathbf{k}}\right) \chi_s(\mathbf{k}) \\ = -s\frac{a}{4} \,v_F \,k_2 \left[4k_1^2 - \mathbf{k}^2\right] (\mathbf{k}^2)^{-1/2} \,\psi_s(\mathbf{k}) = -s\frac{3}{8} \,t \,a^2 \mathbf{k}^2 \sin\left(3\theta\right) \,\psi_s(\mathbf{k}). \tag{7.2}$$

We will assume that in the presence of an electromagnetic field **A**, derivatives change into covariant derivatives and the perturbation just turns into the Hermitian expression

$$\Delta H = -\frac{a}{8} v_F \{ (p_2 - eA_2) [4(p_1 - eA_1)^2 - (\mathbf{p} - e\mathbf{A})^2] [(\mathbf{p} - e\mathbf{A})^2]^{-1} \boldsymbol{\sigma} \cdot (\mathbf{p} - e\mathbf{A}) + \boldsymbol{\sigma} \cdot (\mathbf{p} - e\mathbf{A}) [(\mathbf{p} - e\mathbf{A})^2]^{-1} [4(p_1 - eA_1)^2 - (\mathbf{p} - e\mathbf{A})^2] (p_2 - eA_2) \}.$$
(7.3)

Note that there is an order indeterminacy in this definition. We will come back later to this point.

We now consider a constant magnetic field as in equation (4.1). Then,

$$p_1 - eA_1 = p_1 = \sqrt{eB} p, \quad p_2 - eA_2 = p_2 - eBx_1 = -\sqrt{eB} q,$$
 (7.4)

and $\triangle H$ reduces to

$$\Delta H = \mathfrak{H} + \mathfrak{H}^{\dagger}, \tag{7.5}$$

where

$$\mathfrak{H} = -\frac{a}{8} v_F \sqrt{eB} q [4p^2 - (p^2 + q^2)] [p^2 + q^2]^{-1} [p\sigma_1 - q\sigma_2].$$
(7.6)

When \mathfrak{H} is applied to the functions in equation (4.14), taking into account that (see equation (4.6))

$$[p\sigma_1 - q\sigma_2] \psi_{k,n,s}(\mathbf{x}) = -\frac{1}{w} \left\{ \frac{m}{eB} \mathcal{E}_{n,s} - \frac{1}{2} \left(p^2 + q^2 \right) \right\} \psi_{k,n,s}(\mathbf{x}), \tag{7.7}$$

where $\mathcal{E}_{n,s}$ is given in equation (4.29), we get

$$\mathfrak{H}\psi_{k,n,s}(\mathbf{x}) = \frac{a}{8} v_F \sqrt{eB} \frac{K_{n,s}}{w} \frac{e^{ikx_2}}{\sqrt{2\pi}} q e^{-\frac{q^2}{2}} \\ \times \left(\frac{\left\{ \frac{m}{eB} \mathcal{E}_{n,s} - \frac{1}{2} \left(2n + 3 \right) \right\}}{\left[2n + 3 \right]} \left[4q^2 - 3(2n + 3) \right] H_{n+1}(q) \\ - \iota \left[\frac{1 - s\sqrt{1 + 8w^2(n+1)}}{2w} \right] \frac{\left\{ \frac{m}{eB} \mathcal{E}_{n,s} - \frac{1}{2} \left(2n + 1 \right) \right\}}{\left[2n + 1 \right]} \left[4q^2 - 3(2n + 1) \right] H_n(q) \right).$$

$$(7.8)$$

The first order perturbations on the eigenvalues of the Hamiltonian are given by the matrix elements

$$\begin{aligned} (\psi_{k',n,s}, \Delta H\psi_{k,n,s}) &= \left(\psi_{k',n,s}, \mathfrak{H}\psi_{k,n,s}\right) + \left(\mathfrak{H}\psi_{k',n,s}, \psi_{k,n,s}\right) \\ &= \delta(k-k') \, a \, v_F \sqrt{eB} \, \frac{K_{n,s}^2}{w} \\ &\times \int_{-\infty}^{\infty} \mathrm{d}q \, \mathrm{e}^{q^2} q \left\{ \frac{\left\{\frac{m}{eB} \, \mathcal{E}_{n,s} - \frac{1}{2} \, (2n+3)\right\}}{[2n+3]} [4q^2 - 3(2n+3)] H_{n+1}^2(q) \\ &+ \left[\frac{1 - s\sqrt{1+8w^2(n+1)}}{2w}\right]^2 \frac{\left\{\frac{m}{eB} \, \mathcal{E}_{n,s} - \frac{1}{2} \, (2n+1)\right\}}{[2n+1]} [4q^2 - 3(2n+1)] H_n^2(q) \right\} = 0, \end{aligned}$$

$$(7.9)$$

since it is the integral on the whole line of an odd function of q.

As previously noted, there is an order indeterminacy in the expression of \mathfrak{H} in equation (7.6), since the operators on the right-hand side of equation (7.1) commute. But any order one chooses will lead to an odd in *q* integrand in the expression of $(\psi_{k',n,s}, \Delta H \psi_{k,n,s})$, giving then a vanishing result.

It is also immediate to show that $\mathfrak{H}\psi_0 = 0$, with ψ_0 given in equation (4.21).

Therefore, we get no modification of the energy eigenvalues at first order in perturbation theory with ΔH in equation (7.3) (or with any permutation of the factors in \mathfrak{H} in equation (7.6)). This means that, in the presence of a constant magnetic field orthogonal to the plane, the spectrum of the Hamiltonian of our model seems to be rather stable against the rotational symmetry breaking perturbation in equation (7.5).

In particular, the Hall conductivity of the model described in section 6 and the explanation of the anomalous integer quantum Hall effect based on it is not modified by the first order perturbation theory in ΔH , operator which incorporates the (next to leading) quadratic term in the nearest-neighbor interactions.

8. Crossed magnetic and electric fields

Let us also consider an electric field in the direction of x_1 . So we can take the electromagnetic field A_{μ} as

$$\mathbf{A} := Bx_1 \,\hat{\mathbf{e}}_2, \quad A_0 := -Ex_1, \tag{8.1}$$

also in the Coulomb gauge.

In this case, the Hamiltonian is

$$H = \frac{(\mathbf{p} - e\mathbf{A})^2}{2m} + v_F \,\boldsymbol{\sigma} \cdot (\mathbf{p} - e\mathbf{A}) - eA_0. \tag{8.2}$$

In the $m \to \infty$ limit, a pseudo-Lorentz transformation of the linear Hamiltonian allows us to solve the eigenvalue equation for crossed electric and magnetic fields in terms of the solutions of the case in which there is only a magnetic field applied perpendicularly to the plane of graphene [45–47]. But, in this nonrelativistic model, the quadratic mass term prevents us to get the solution in this way.

Translation invariance in the x_2 direction suggests that the solutions are again of the form $\Psi(\mathbf{x}) = \frac{e^{i k x_2}}{\sqrt{2\pi}} {\varphi(x_1) \choose \chi(x_1)}$, and the eigenvalue equation reduces to

$$\begin{cases} \left[p_1^2 + (eB)^2 \left(x_1 - \frac{k}{eB} \right)^2 \right] - \lambda + 2meEx_1 \end{cases} \varphi = -2mv_F \left\{ p_1 + \iota eB \left(x_1 - \frac{k}{eB} \right) \right\} \chi, \\ \left\{ \left[p_1^2 + (eB)^2 \left(x_1 - \frac{k}{eB} \right)^2 \right] - \lambda + 2meEx_1 \right\} \chi = -2mv_F \left\{ p_1 - \iota eB \left(x_1 - \frac{k}{eB} \right) \right\} \varphi, \end{cases}$$

$$(8.3)$$

with $\lambda = 2m\mathcal{E}$.

It is convenient to define a new variable as $q := \sqrt{eB} \left[x_1 - \frac{k}{eB} + \frac{mE}{eB^2} \right]$ and, correspondingly, $p := -\iota \frac{\partial}{\partial q} = \left(\frac{-\iota}{\sqrt{eB}} \right) \frac{\partial}{\partial x_1}$, in terms of which the eigenvalue equation is written as

$$\left\{p^{2}+q^{2}-\frac{\Lambda}{eB}\right\}\varphi(q) = 2w\left\{[p+\iota q]+\iota w\frac{E}{v_{F}B}\right\}\chi(q),$$

$$\left\{p^{2}+q^{2}-\frac{\Lambda}{eB}\right\}\chi(q) = 2w\left\{[p-\iota q]-\iota w\frac{E}{v_{F}B}\right\}\varphi(q),$$
(8.4)

where

$$\Lambda := \lambda + \frac{mE}{B} \left[\frac{mE}{B} - 2k \right]$$
(8.5)

with $\lambda = 2m\mathcal{E}$.

For small electric fields, the last terms on the right-hand sides of equation (8.4) can be treated as the perturbation $\left(\frac{w^2 E}{mv_F B}\right)\sigma_2$ on a zero order Hamiltonian whose solutions are the same functions of q as in section 4, with the equilibrium position of the oscillator in the x_1 -direction displaced by an amount $\left(-\frac{mE}{B}\right)$. The zero-order energy eigenvalues are then those of section 4 but shifted by an amount $\frac{E}{B}\left[k - \frac{mE}{2B}\right]$.

On the other hand, since the perturbation is proportional to σ_2 , it has a vanishing mean value due to the orthogonality of Hermite functions (see equation (4.14)). Then, the spectrum is just rigidly shifted by the presence of an electric field perpendicular to the magnetic field, up to first order in perturbation theory in the small parameter $(E/v_F B)$.

9. Conclusions and outlook

In this paper, we have studied a simple nonrelativistic model, suggested by a non-Abelian magnetic quantization approach (based on a Heisenberg algebra deformed by the introduction of an external constant non-Abelian magnetic field), to describe the low energy excitations of graphene. As previously mentioned, these states admit an effective description as excitations of a (2+1)-dimensional pseudo-relativistic field theory of massless Dirac fermions.

Our effective model is based on a deformation of the Heisenberg algebra which makes the commutator of momenta proportional to the pseudo-spin. This can be interpreted as the commutator of two U(2) covariant derivatives in the case of a constant background non-Abelian magnetic field. The modification introduced in the Heisenberg algebra leads to wavefunctions of two components and a Hamiltoninan operator in the form of a 2 × 2 matrix, which reduces to the usually considered linear one in the large mass limit.

We have solved the Landau problem for this planar system, with a constant and uniform magnetic field applied perpendicularly to the plane, explicitly obtaining the Hamiltonian (generalized) eigenfunctions and eigenvalues. We have pointed out that this construction can be directly extended to the case of negative values of the mass parameter. This is important since a negative mass allows us to reproduce with our Hamiltonian the low energy expansion of the dispersion relation around the first Dirac point of graphene, \mathbf{K} , when the leading order terms from both nearest and next-to-nearest-neighbor interactions (linear and quadratic terms in the quasi-momentum expansion, respectively) are retained.

The effective Hamiltonian for the excitations around the second Fermi point, $H_{\mathbf{K}'}$, is proposed to be obtained from $H_{\mathbf{K}}$ through a transformation which maintains the Hamiltonian time reversal times parity invariant.

Through this transformation, the spectrum of the low energy states for the second Dirac point **K**' is simply obtained from that of **K** by a reflection around the origin. This leads to an almost doubly degenerate spectrum for quasi-particles and holes, where the degeneracy is broken by $O(w^{-1})$ terms, with $w = \frac{|m|v_F}{\sqrt{eB}} \simeq 10^3$ for realistic values of the parameters. Moreover, there is one quasi-particle state of energy $\mathcal{E}'_0 = eB/2|m|$ and one hole state of energy $\mathcal{E}_0 = -eB/2|m|$.

The $|m| \rightarrow \infty$ limit reproduces the well-known doubly degenerate spectrum of massless Dirac fermions, with one quasi-particle zero mode and one hole zero mode.

The Hall conductivity of the model has been evaluated from the partition function employing the ζ -function approach to the associated functional determinant. This led us to a rather general expression valid for the Landau problem of a fermionic theory. In the case under study, the result is consistent with the anomalous integer quantum Hall effect found in graphene, in which the (reduced) Hall conductivity as a function of the Fermi energy takes half-integer values (see figure 3). In particular, $\frac{h}{4e^2}\sigma_{xy} = \pm \frac{1}{2}$ for small (but $|\epsilon_F| > \frac{v_F \sqrt{eB}}{2w}$) positive or negative Fermi level, respectively.

Moreover, we have proposed an additional rotational symmetry breaking term to be treated as a perturbation on the Hamiltonian, which reproduces the next-to-leading (quadratic) term in the nearest-neighbor interaction contribution to the dispersion relation with no external field. With the perpendicular magnetic field present, we have shown that this term does not modify the spectrum at first order in perturbation theory. Neither does it change the behavior of the Hall conductivity described above.

Finally, we have shown that an electric field perpendicular to the magnetic field produces just a rigid shift of the spectrum at first order in perturbation theory.

An interesting problem under study refers to the case in which an Aharonov–Bohm singular flux is also present, which can simulate a defect in the lattice. This will be the subject of a forthcoming paper.

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Appendix. Lagrangian, conserved current and generating functional

In this appendix, we construct the Lagrangian for our model and deduce the conserved current associated with its U(1) gauge symmetry. We also discuss the behavior of the generating functional in relation with the Hall conductivity.

The system under consideration can be described by the Lagrangian

$$\mathcal{L} := \frac{l}{2} [\psi^{\dagger} \partial_t \psi - \partial_t \psi^{\dagger} \psi] + \psi^{\dagger} e A_0 \psi - \frac{1}{2m} \{ [(\mathbf{p} - e\mathbf{A} + \theta \sigma) \psi]^{\dagger} \cdot [(\mathbf{p} - e\mathbf{A} + \theta \sigma) \psi] - 2\theta^2 \psi^{\dagger} \psi \}.$$
(A.1)

The variation of the independent dynamical variables ψ^{\dagger} and ψ leads to the Euler–Lagrange equations in the usual way:

$$\frac{\partial \mathcal{L}}{\partial \psi^{\dagger}} - \partial_t \left(\frac{\partial \mathcal{L}}{\partial \left(\partial_t \psi^{\dagger} \right)} \right) - \nabla \cdot \left(\frac{\partial \mathcal{L}}{\partial \left(\nabla \psi^{\dagger} \right)} \right) = \iota \partial_t \psi - \frac{1}{2m} [(\mathbf{p} - e\mathbf{A} + \theta \sigma)^2 - 2\theta^2] \psi = 0,$$
(A.2)

and similarly for ψ .

The Lagrangian in equation (A.1) has a U(1) gauge symmetry. Indeed, it remains invariant against the following local change in the dynamical variables and the electromagnetic field

$$\psi(x) \to e^{ie\alpha(x)}\psi(x) \implies \delta\psi(x) = ie\alpha(x)\psi(x),$$

$$\psi^{\dagger}(x) \to \psi^{\dagger}(x)e^{-ie\alpha(x)} \implies \delta\psi^{\dagger}(x) = -ie\alpha(x)\psi^{\dagger}(x),$$

$$A_{\mu}(x) \to A_{\mu}(x) + \partial_{\mu}\alpha(x).$$
(A.3)

Then, Nother's theorem implies the existence of a locally conserved current given by

$$\alpha j^{\mu} := -\delta \psi^{\dagger} \left(\frac{\partial \mathcal{L}}{\partial \left(\partial_{\mu} \psi^{\dagger} \right)} \right) - \left(\frac{\partial \mathcal{L}}{\partial \left(\partial_{\mu} \psi \right)} \right) \delta \psi.$$
(A.4)

Therefore, the charge density is

$$j^0 = e \,\psi^{\dagger} \psi \tag{A.5}$$

and the current density

$$\mathbf{j} = \frac{e}{2m} \{ \iota (\nabla \psi^{\dagger} \psi - \psi^{\dagger} \nabla \psi) + 2\psi^{\dagger} (-e\mathbf{A} + \theta \sigma) \psi \}.$$
(A.6)

It is a simple exercise to verify that j^{μ} is conserved as a consequence of the equations of motion, equation (A.2), and its complex conjugate,

$$\partial_t j^0 - \nabla \cdot \mathbf{j} = 0. \tag{A.7}$$

The generating functional of (current) Green's functions of this two-dimensional system reads as

$$Z[A] = e^{i\Gamma[A]} := \int \mathcal{D}\psi^{\dagger} \mathcal{D}\psi \ e^{i\int dt \int d^2x \mathcal{L}(x)}, \tag{A.8}$$

where $\Gamma[A]$ is the effective action. In particular, we have

$$-\iota \frac{\delta \log Z[A]}{\delta A_{\mu}(x)} = \langle j^{\mu}(x) \rangle = J^{\mu}(x)$$
(A.9)

and the second functional derivative gives minus the current-current correlation function.

For a system with a gap in its spectrum (as is the case of our model with filled Landau levels), the effective action is a local gauge invariant functional of the electromagnetic external field [39, 40],

$$\Gamma[A] = \int dt \int d^2x \, \mathcal{L}_{\text{eff}}(A_{\mu}(x), \, \partial_{\nu}A_{\mu}(x), \, \ldots).$$
(A.10)

In general, in an asymptotic weak field and weak gradient expansion, gauge invariance determines that the effective Lagrangian \mathcal{L}_{eff} appears as an expansion in terms of the field intensities **E** and *B* and its derivatives. But in 2+1 dimensions, there is an additional possibility, since a Chern–Simons term can be the dominant one in this expansion (see, for example, [39, 40] and references therein):

$$\mathcal{L}_{\text{eff}} = \frac{K}{4\pi} \,\epsilon^{\mu\nu\lambda} A_{\mu} \partial_{\nu} A_{\lambda} - \frac{1}{2} \,E_i \rho^{ij} E_j - \frac{1}{2} \,\chi B^2 + \cdots, \qquad (A.11)$$

where $\epsilon^{\mu\nu\lambda}$ is the completely anti-symmetric tensor with $\epsilon^{012} = 1$ and the ellipses represent higher order terms.

In fact, it is the Chern–Simons term which determines the Hall conductivity since

$$J^{\mu}(x) = \frac{\delta\Gamma[A]}{\delta A_{\mu}(x)} = \frac{K}{2\pi} \epsilon^{\mu\nu\lambda} \partial_{\nu}A_{\lambda} + \ldots = \frac{K}{4\pi} \epsilon^{\mu\nu\lambda} F_{\nu\lambda} + \cdots.$$
(A.12)

In particular, for the spatial and temporal components of the current, the previous expression reduces to

$$J_i(x) = \frac{K}{2\pi} \epsilon_{0ij} E_j + \cdots, \quad J_0(x) = \frac{K}{2\pi} B + \cdots.$$
 (A.13)

The first equation in (A.13) implies that the Hall conductivity $\sigma_{xy} = \frac{K}{2\pi}$ (or $\sigma_{xy} = Ke^2/h$ in full units). Then,

$$J_0 = \sigma_{xy} \mathcal{B}. \tag{A.14}$$

It can be shown that K takes integer values if the (many-body) ground state is not degenerate and has a finite energy gap [39, 40, 47]. In this case, K has a topological character: it can be related to a Berry phase which, in turn, takes the value 2π times an (integer) Chern number.

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