Effective Magnetic Fields in Graphene Superlattices

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We demonstrate that the electronic spectrum of graphene in a one-dimensional periodic potential will develop a Landau level spectrum when the potential magnitude varies slowly in space. The effect is related to extra Dirac points generated by the potential whose positions are sensitive to its magnitude. We develop an effective theory that exploits a chiral symmetry in the Dirac Hamiltonian description with a superlattice potential, to show that the low energy theory contains an effective magnetic field. Numerical diagonalization of the Dirac equation confirms the presence of Landau levels. Possible consequences for transport are discussed.

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Graphene is a honeycomb lattice of carbon atoms with a host of remarkable electronic properties [1, 2]. Many of these are a result of the low energy states, which are governed by a massless Dirac equation. When undoped the Fermi surface of the system is a set of discrete points, with resulting properties that compromise between metallic and insulating behavior. For example, an ideal graphene system can conduct diffusively even in the absence of impurities to scatter the electrons [3, 4]. The possibility of a universal conductivity even in the presence of impurities remains an active subject of discussion [5–13]. Beyond this, graphene exhibits signatures of “effective time-reversal symmetry breaking,” effects most naturally described in terms of effective magnetic flux, due to a Berry’s phase intrinsic to graphene [14–16], or to strain in the lattice system [17–19], even in the absence of a real applied magnetic field.

In this paper we discuss a completely different way to induce an effective magnetic field in graphene, of sufficient uniformity that Landau levels in principle should be apparent in the electronic spectrum. This involves immersing the graphene system in a unidirectional periodic (i.e., superlattice) potential, for example by subjecting a sample with periodic ripples [20] to a perpendicular electric field. Recently it has been shown [21, 22] that such a potential can induce new Dirac points from the original one, whose positions are determined by the amplitude $V_0$ of the periodic potential, and by its period. We demonstrate that by promoting $V_0$ to a slowly varying function of position, the effective theory for states near an induced Dirac point contains a gauge field, such that there are states in the spectrum corresponding to Landau levels.

The possibility of generating such a description turns out to be intimately related to a chiral symmetry which is present in the Dirac equation even with a periodic potential, as we show below. The symmetry requires that any state with energy $\varepsilon$ have a chiral partner at energy $-\varepsilon$. These are degenerate for $\varepsilon = 0$. Eigenstates near zero energy can be approximately constructed from these chiral partners, and they have a number of properties guaranteeing that the resulting effective theory has the form of a Dirac equation with a vector potential, representing a magnetic field with strength proportional to the rate at which the position of the induced Dirac point varies with position. We demonstrate the effect explicitly by numerically solving the Dirac equation with a potential $V_0(x) \cos G_0 x$. Results for the energy spectrum are illustrated in Fig. 1.

![Energy bands for Dirac equation in a modulated superlattice potential. Inset: Ratio of first to second excited state energies near $k_F a/2\pi = 1.3$, demonstrating the $\sqrt{2}$ ratio (dotted line) expected of Landau levels.](https://example.com/figure1.png)

Within this spectrum we find level
spacings that are essentially identical to those of Landau levels (see inset of Fig. 1), and wavefunctions contained in envelopes whose center positions—guiding centers—are fixed by a wavevector, and have the forms of harmonic oscillator states. We find that states that straddle regions in which \( dV_0/dx \) changes sign are strongly dispersive, acting much like edge states of graphene islands in a magnetic field [23]. The presence of such edge states should have very interesting consequences for transport in this system.

Superlattice Potential and Chiral Symmetry—Low energy states of graphene in a single valley may be modeled by a Dirac equation [24]. When an external potential is added, the eigenvalue equation to be solved has the form

\[
H\tilde{\psi} = \varepsilon\tilde{\psi}, \quad H = v_F(\sigma_x p_x + \sigma_y p_y) + V_0(x)\mathbb{L},
\]

where \( \tilde{\psi} \) is a two-component spinor whose entries are the wavefunction amplitudes for the two sublattices, \( v_F \) is the speed of electrons with momenta near the Dirac point, \( \sigma_x, \sigma_y \) are Pauli matrices, \( p_x = -i\partial/\partial x, V_0 \) is the scale of an external potential with functional form \( v(x) \), and \( \mathbb{L} \) is a unit \( 2 \times 2 \) matrix. For \( V_0 = 0 \), \( H \) is known to have a chiral symmetry, which can be expressed as \( \{ \sigma_z, H \} = 0 \), with \( \sigma_z \) the third Pauli matrix. This anticommutator relation implies that for any eigenstate \( \tilde{\psi} \) with energy \( \varepsilon \), there is a particle-hole partner \( \sigma_z \tilde{\psi} \) which is also an eigenstate of \( H \), with energy \(-\varepsilon\). Moreover one may use this property to show that the state at zero energy is doubly degenerate.

When there is a non-vanishing periodic superlattice potential satisfying \( v(x + a/2) = -v(x) \) with energy scale \( V_0 \), it was shown recently [21, 22] that for sufficiently large \( V_0a \) one obtains new Dirac cones in the spectrum. These appear as pairs which emerge from \( k = 0 \) along the \( k_y \) axis, one moving up and the other down as \( V_0a \) increases, so that there are new zero energy states at \( k = (0, \pm k_y^0) \). Further increasing \( V_0a \) generates yet more pairs of new Dirac cones. For simplicity, in what follows we will only consider the situation where a single new pair has emerged.

The Hamiltonian with this type of periodic potential also has chiral symmetry, although the associated operator is slightly more complicated. Defining a shift operation \( S\tilde{\psi}(x, y) = \tilde{\psi}(x + a/2, y) \), it takes the form \( T = \sigma_z S \); one may easily demonstrate \( \{ H, T \} = 0 \) even in the presence of the periodic potential. Because there is a chiral operator, one may show that a zero energy state \( \tilde{\phi}_0 \) at \( (0, k_y^0) \) has a chiral partner \( T\tilde{\phi}_0 \) which is also at zero energy at the same wavevector. For what follows it is convenient to define eigenstates of \( T \), \( \tilde{\phi}_\pm = \tilde{\phi}_0 \pm T\tilde{\phi}_0/N_\pm \), where \( N_\pm \) are normalization integrals. These states have a number of interesting properties which will be useful to us:

\[
<\phi_+|\phi_-> = 0, \quad (1)
\]
\[
<\phi_+|\sigma_{x,y}\phi_\pm> = 0, \quad (2)
\]
\[
<\phi_+|\sigma_x\phi_-> = iu_x, \quad (3)
\]
\[
<\phi_+|\sigma_y\phi_-> = u_y, \quad (4)
\]

where \( u_{x,y} \) are purely real. Eq. (1) is easily shown by observing that \( \tilde{\phi}_\pm \) are eigenstates of \( T \) with different eigenvalues. Eq. (2) follows from \( \{\sigma_{x,y}, \sigma_z\} = 0 \) and \( <\phi_\pm|T^2|\phi_\pm> = 1 \). Eqs. (3) and (4) follow from a specific form which \( \tilde{\phi}_0 \) may take, \( \tilde{\phi}_0(r) = \phi(x)|i\sin \theta(x), \cos \theta(x)|e^{ik_yy} \), where \( \phi \) and \( \theta \) are real periodic functions [21]. The quantities \( u_{x,y} \) may be understood as the velocities of the electron states in the \((x, y)\) directions of a Dirac cone, in units of \( v_F \), as we shall see below.

Modulated Superlattice—We now consider the situation \( V_0 \rightarrow V_0(x) \), where \( V_0 \) varies slowly on the scale of the superlattice parameter \( a \). The eigenfunctions generated from \( \tilde{\phi}_0 \) have an extra spatial dependence from their implicit dependence on \( V_0 \): \( \tilde{\phi}_0(x, y; V_0) \rightarrow \tilde{\phi}_0(x, y; V_0(x)) \).

In what follows it will be convenient to consider functions where the explicit spatial dependence and that inherited from \( V_0(x) \) are independent, so that we promote the wavefunctions to functions of two variables: \( \tilde{\phi}_0(r'; V_0(r)) \rightarrow \tilde{\phi}_0(r', r; V_0(x)) \). Note that in this construction, if \( r \) is held constant then \( \tilde{\phi}_0(x' \rightarrow a, y'; r) \rightarrow \tilde{\phi}_0(x', y'; r) \).

We can then define matrix elements of the form

\[
(\phi_1|A|\phi_2)_r = \int_{x-a/2}^{x+a/2} \int_{-\infty}^{\infty} dy' g(y-y')\tilde{\phi}_1(r', r)\varepsilon\tilde{\phi}_2(r', r),
\]

where \( g(y-y') \) is a peaked function centered at 0 and of width much greater than 2\( \pi/k_y^0 \) [23] which integrates to 1, and \( A \) is an operator. Identifying \( \tilde{\phi}_\pm \equiv \tilde{\phi}_\pm(r'; V_0(x)) \), Eqs. (1)-(4) remain true if we replace the standard matrix elements with those above. We now consider a generalized eigenvalue problem \( H(r', r)\tilde{\psi}(r', r) = \varepsilon\tilde{\psi}(r', r) \), where

\[
H(r', r) = v_F[-i\sigma_x(\partial_{r'}+\partial_x)-i\sigma_y(\partial_{r'}+\partial_y)]+V_0(x)v(x')\mathbb{L},
\]

and we have set \( \hbar = 1 \). In the limit \( r' \rightarrow r \), this becomes the eigenvalue equation one needs to solve for the physical Hamiltonian; the advantage of the generalized problem is that it retains the chiral symmetry \( \{H, T\} = 0 \), with the shift operation \( (S) \) contained in \( T \) acting only \( r' \). To approximately solve this problem, we consider solutions of the form [24]

\[
\tilde{\psi}(r', r) = \psi_+(r)\tilde{\phi}_+(r', r) + \psi_-(r)\tilde{\phi}_-(r', r).
\]

Acting on this wavefunction with \( H \), and then forming matrix elements \( (\phi_1|H-\varepsilon|\psi_1) \), we seek choices of the coefficients \( \psi_{\pm} \) for which these matrix elements vanish. After some algebra [25], with the use of Eqs. (1)-(4), one finds that they must obey the equations

\[
\begin{pmatrix}
0 & v_x \partial_x - iv_y \partial_y + a_+ (r) \\
-v_x \partial_x - iv_y \partial_y + a_- (r) & 0
\end{pmatrix}
\begin{pmatrix}
\psi_+ \\
\psi_-
\end{pmatrix}
= \varepsilon
\begin{pmatrix}
\psi_+ \\
\psi_-
\end{pmatrix}.
\]

Here \( v_{x,y}(x) = v_F u_{x,y}[V_0(x)] \), and

\[
a_{\pm, \mp} (r) = -iv_F(\phi_0|\vec{\sigma} \cdot \vec{V}|\phi_\pm) r.
\]
It is important to note because the gradient operator in \( a_{±} \) acts on \( r \) and not \( r' \), \( a_{±} \neq a_{±}^\dagger \). Thus the matrix appearing in Eq. 17 is not Hermitian. However, one may show that \( \text{Im} a_{±} = -\text{Im} a_{±} = v_y \phi \partial_{x} k_y^0 \equiv v_y A_x \) is purely antisymmetric, while the real antisymmetric component of \( a_{±} a_{±}^\dagger \equiv \text{Re}(a_{±} - a_{±})/2 \), can be removed by a similarity transformation. Rewriting Eq. 17 in terms of \( \hat{\psi}_± \) we obtain a Hermitian eigenvalue equation \( \hat{H} \psi = \varepsilon \psi \), with

\[
\hat{H} = \sigma_y v_x [-i \partial_x - A_x] + \sigma_x v_y [-i \partial_y - A_y],
\]

where \( A_y = -\text{Re}(a_{±} + a_{±})/2v_y \). Eq. 6 has the form of a Dirac Hamiltonian with a vector potential, albeit with the roles of the \( x \) and \( y \) momenta interchanged, and the effective velocities \( v_{x,y}(x) \) being direction and position dependent. Several remarks are in order. (i) For \( V_0 \) constant, \( v_{x,y} \) lose their position dependence, and we expect them to give the velocities of the induced Dirac points for an unmodulated superlattice potential. Fig. 2 illustrates by comparison with velocities found numerically around an induced Dirac point of a graphene superlattice \([21,22]\) that this is indeed the case. (ii) Chiral symmetry plays a central role in giving Eq. 6 its Dirac form, through the properties of the \( \phi_{±} \) states expressed in Eqs. 11. In particular they guarantee that diagonal elements do not appear in the Hamiltonian, and that the off-diagonal terms have a form that can naturally be grouped together into canonical momenta with vector potential contributions. (iii) The effective field represented by the vector potential is \( B_{eff} = \partial_x A_y - \partial_y A_x \). In the limit of a slowly varying modulation \( (|a| \partial_x V_0)/V_0 \ll 1 \), the first term is second order in derivatives and is negligibly small, so that the effective field is \( B_{eff} = -\partial_x k_y^0 \equiv - (\partial_x k_y^0 V_0)/(\partial_x V_0(x)) \). We see that the presence of an induced magnetic field is intimately connected with the fact that the position of the Dirac point in the Brillouin zone can be changed via an external parameter. (iv) With a gauge transformation, \( \hat{H} \) can be written in a form for which there is no explicit \( y \) dependence, so that \( k_y \) is a good quantum number, as we expect because the system is translationally invariant along the \( \hat{y} \) direction. If we define a magnetic length \( \ell_0 = 1/\sqrt{\mu B_{eff}} \) then \( X = k_y \ell_0^2 \) becomes a guiding center coordinate around which the wavefunction is centered. Approximating \( v_{x,y}(x) \rightarrow v_{x,y}(X) \), \( \hat{H} \) becomes a standard Hamiltonian for Dirac electrons in a magnetic field; the approximation is valid provided \( \ell_0 \) is sufficiently small. In this case one expects eigenenergies of the form \( \varepsilon_n(X) = \text{sgn}(n) \sqrt{2v_x v_y |n|B_{eff}} \) with \( n \) an integer. Although the values of \( \varepsilon_n(X) \) depend on \( k_y \) through \( v_x v_y B_{eff} \), neighboring energy levels \( n_1, n_2 \) at fixed \( k_y \) should obey \( \varepsilon_{n_1}/\varepsilon_{n_2} = \sqrt{n_1/n_2} \). This is an unambiguous signature of Landau levels, which we next demonstrate numerically to be present in the energy spectrum of this system.

**Numerical Studies** – Returning to the original Hamiltonian, \( \hat{H} = v_F(\sigma_x p_x + \sigma_y p_y) + V_0(x) \psi(x) \mathcal{P} \), as a concrete example we take \( V_0(x) \) to have a sawtooth shape (illustrated as a dashed line in Fig. 3), acting as an envelop for a pure cosine potential \( v(x) = \cos G_0 x \), with \( G_0 = 2\pi/a \). The period of \( V_0 \) is chosen to be \( L_0 = 50a \), and its minimum/maximum values are \( \pm 0.1U_0 \), with \( U_0 = 0.12\epsilon_F \) and \( \epsilon_F = \hbar v_F/a \). We expand \( \hat{H} \) in plane wave states, for fixed values of \( k_x \) (the Bloch wavevector) and \( k_y \), and diagonalize the resulting matrix. A typical example of a few of the eigenvalues as a function of \( k_y \), for \( k_x = 0 \), are illustrated in Fig. 3. Near \( k_y = 0 \) the Dirac point which would also be present for \( V_0 = 0 \) is apparent, but at larger \( |k_y| \) one finds bands that vary nearly linearly with \( k_y \), in the vicinity of the emergent Dirac points one would find if \( V_0(x) \) were replaced by its spatial average in the potential \([21,22]\). These are the Landau levels generated by the potential modulation. The zero energy states one expects of Dirac Landau bands are very apparent. The inset of Fig. 1 shows the ratio of energies for the two lowest positive energy states; for Landau levels these should scale as \( \sqrt{n_2/n_1} = \sqrt{2} \), which is indeed the case as illustrated in the inset. The levels shown are each very nearly doubly degenerate. This is because our chosen \( V_0(x) \) satisfies \( dV_0(x)/dx = -dV_0(x + L_0/2)/dx \), so that the effective magnetic field has regions of equal magnitude but opposite directions. We find that the spectra in the vicinity of the Landau levels are quantitatively nearly identical for any choice of \( k_x \).

Fig. 3 illustrates wavefunction overlaps with \( \phi_{±} \) for the lowest positive energy states for a single choice of \( k_y \). Their similarity to a lowest harmonic oscillator state for one of these, and to a first excited excited state for the other, is apparent. This is precisely what one expects for eigenstates of the Dirac equation in a uniform magnetic field \([1]\). Note that the wavefunction has support in both...
the positive and negative effective field regions, but the roles of $\phi_+$ and $\phi_-$ which multiply the lower or higher harmonic oscillator states are reversed. This reflects the fact that Landau level raising and lowering operators interchange roles when the field sign is reversed.

It is interesting to note that the degeneracy of the positive and negative field states breaks down when $k_y$ is chosen so that the wavefunction is significantly different than zero at the cusps of $V_0(x)$. In this situation the energy rises quickly as a function of $k_y$, indicating a large current. These behave much as edge states in a quantum Hall system, and should have interesting consequences for transport. In particular we expect that $\sigma_{\bar{y}y}$, the conductance along the $\bar{y}$ direction, should have steps as a function of Fermi energy $E_F$, occurring at energies approximately reflecting the $\sqrt{nB_{eff}}$ behavior of the Landau levels. An observation of this effect would constitute evidence of the effective magnetic field present in this system. It is also interesting to note that left- and right-moving states are spatially separated, existing separately at the upward and downward cusps of $V_0(x)$. We thus expect that backscattering due to disorder will be somewhat suppressed, although scattering into a time-reversed partner from the opposite valley will be present if disorder at short wavelengths is significant in the system. Studies of this and related transport phenomena are currently underway.

In summary, we have demonstrated that an effective magnetic field can be generated in a graphene superlattice by a slow modulation of the superlattice potential amplitude. A description in terms of a gauge field can be produced from zero energy states at a Dirac point of the unmodulated superlattice which are chiral partners. These may be used as a basis to find states in the vicinity of the modulated potential. For a slowly varying amplitude, it was found that the effective field is produced by the changing position of the Dirac point in real space. Numerical studies of the Dirac equation with this type of potential verified the presence of Landau levels in the spectrum, and suggest the possibility of interesting transport phenomena in this system.

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[25] See EPAPS Document No. XXX for some details of the derivation of Eq. 5.
In this supplement, we provide some details of how Eq. 5 is obtained from the generalized eigenvalue problem, 
\[ H(\mathbf{r}', \mathbf{r}) \psi(\mathbf{r}', \mathbf{r}) = \varepsilon \psi(\mathbf{r}', \mathbf{r}), \]
with
\[ H(\mathbf{r}', \mathbf{r}) = \hbar v_F \left[ -i \sigma_x (\partial_{x'} + \partial_x) - i \sigma_y (\partial_{y'} + \partial_y) \right] + V_0(x) v(x') \Theta. \]  
(7)

We consider approximate solutions of the form
\[ \psi(\mathbf{r}', \mathbf{r}) = \psi_0(\mathbf{r}) \hat{\phi}_+(\mathbf{r}', \mathbf{r}) + \psi_0(\mathbf{r}) \hat{\phi}_-(\mathbf{r}', \mathbf{r}), \]
where \( \hat{\phi}_\pm = [\hat{\phi}_0 \pm T \hat{\phi}_0] / N_\pm \), and \( H_0 \hat{\phi}_0 = 0 \), with
\[ H_0(\mathbf{r}', \mathbf{r}) = \hbar v_F \left[ -i \sigma_x \partial_{x'} - i \sigma_y \partial_{y'} \right] + V_0(x) v(x') \Theta. \]  
(8)

In this context, the operator \( T \) is defined as
\[ T \hat{\phi}(\mathbf{r}', \mathbf{r}) \equiv \sigma_z S \hat{\phi}(\mathbf{r}', \mathbf{r}) \equiv \sigma_z \hat{\phi}(\mathbf{r}' + a \hat{\mathbf{x}}/2, \mathbf{r}). \] Since \( \mathbf{r} \) is a fixed parameter, the potential in \( H_0 \) is a periodic function of \( x' \), and \( \hat{\phi}_0 \) and \( \hat{\phi}_\pm \) are zero energy Bloch states with respect to \( \mathbf{r}' \). As shown in Refs. 21 and 22, such states are present at \( k_x = 0 \) and \( k_y = k_y^* \), with \( k_y^* \) depending on \( V_0(x) \). Applying Eq. (8) to \( \hat{\phi}_\pm \), multiplying by \( g(y - y') \phi_\pm \), and integrating \( \mathbf{r}' \) over all \( y \) and a narrow interval of width \( a \) around \( x \), one obtains two equations for the coefficients \( \psi_\pm \), which may be expressed as
\[ \begin{align*}
(\phi_+ | \sigma \cdot \nabla | \phi_+)_r &= \int_{x - a/2}^{x + a/2} dx' \int_{-\infty}^{\infty} g(y - y') \phi_+^A [x'; V_0(x)] e^{ik_y^* [V_0(x)] y'} \\
&\times \left\{ \phi_+^A [x'; V_0(x)] [\partial_x + iy' (\partial_x k_y^*)] \phi_+^B [x'; V_0(x)] + \phi_+^B [x'; V_0(x)]^* [\partial_x + iy' (\partial_x k_y^*)] \phi_+^A [x'; V_0(x)] \right\}.
\end{align*} \]
(11)

The terms involving \( \partial_x k_y^* \) are proportional to \( \int_{x - a/2}^{x + a/2} dx' \phi_- | x', V_0(x) \rangle^\dagger \sigma_x \phi_- | x', V_0(x) \rangle \), which vanishes identically, as in Eq. 2 of the text. What remains of this matrix element can be cast in the form
\[ \begin{align*}
(\phi_- | \sigma \cdot \nabla | \phi_-)_r &= \int_{x - a/2}^{x + a/2} dx' \phi_- | x', y' = 0; V_0(x) \rangle^\dagger \sigma_x \partial_y \phi_- | x', y' = 0; V_0(x) \rangle \\
&= \int_{x - a/2}^{x + a/2} dx' \frac{1}{N_- (x)} \left[ \phi_0^A [x'; V_0(x)] - \sigma_z S \phi_0^A [x'; V_0(x)] \right] \\
&\times \sigma_x \partial_x \left\{ \frac{1}{N_- (x)} \left[ \phi_0^A [x'; V_0(x)] - \sigma_z S \phi_0^A [x'; V_0(x)] \right] \right\},
\end{align*} \]
(12)

where in the last expression we have dropped the explicit \( y' \) in the arguments of the wavefunctions. Note the normalization factors \( N_- \) obtain their spatial dependence from \( V_0(x) \). When the derivative operator acts on \( 1/N_- \),
the resulting contribution is again proportional to \((\phi_-\sigma_x|\phi_-)_r\) and vanishes. We now have

\[
(\phi_-|\vec{\sigma} \cdot \vec{\nabla}|\phi_-)_r = \frac{1}{N_-(x)^2} \int_{x-a/2}^{x+a/2} dx' \left[ \phi_0[x';V_0(x)] - \sigma_z S \phi_0[x';V_0(x)] \right]^{\dagger} \\
\times \sigma_x \partial_z \left\{ \phi_0[x';V_0(x)] - \sigma_z S \phi_0[x';V_0(x)] \right\} \\
= \frac{1}{N_-(x)^2} \int_{x-a/2}^{x+a/2} dx' \left\{ \phi_0^* \sigma_x \partial_z \phi_0 - [S \phi_0][\sigma_z \partial_z S \phi_0] \\
+ [S \phi_0]^\dagger \sigma_x \sigma_z \partial_z \phi_0 - \phi_0^* \sigma_x \sigma_z \partial_z \phi_0 \right\}.
\]

(13)

where we used \(\{\sigma_x,\sigma_z\} = 0\) and \(\sigma_z^2 = 1\). Finally, noting that \(S\) is a Hermitian operator in \(x\) and that \(S^2 = 1\) since the wavefunctions are periodic in \(x\), it easy to see that the four terms of the last expression mutually cancel. We thus have established \((\phi_-|\vec{\sigma} \cdot \vec{\nabla}|\phi_-)_r = 0\). A very similar manipulation leads to \((\phi_+|\vec{\sigma} \cdot \vec{\nabla}|\phi_+)_r = 0\) as well. Note that in reaching these results, the antisymmetry of \(\sigma_x\) and \(\sigma_z\) played a crucial role, so that the vanishing of this matrix element is in part a consequence of the fact that we can define a chiral operator for this system. That \((\phi_\pm|\vec{\sigma} \cdot \vec{\nabla}|\phi_\pm)_r\) vanish is significant because had they not, the matrix equation for \(\psi_\pm\) would contain diagonal elements, effectively adding a scalar potential in addition to the vector potential we find generated by the modulation of \(V_0\).

Our equations now reduce to the form

\[-i\hbar v_F(\phi_+|\sigma_x|\phi_-)_r \partial_z \psi_- - i\hbar v_F(\phi_+|\sigma_y|\phi_-)_r \partial_y \psi_- + -i\hbar v_F(\phi_+|\vec{\sigma} \cdot \vec{\nabla}|\phi_-)_r \psi_- = \varepsilon \psi_+ \]

\[-i\hbar v_F(\phi_-|\sigma_x|\phi_+)_r \partial_z \psi_+ - i\hbar v_F(\phi_-|\sigma_y|\phi_+)_r \partial_y \psi_+ + -i\hbar v_F(\phi_-|\vec{\sigma} \cdot \vec{\nabla}|\phi_+)_r \psi_+ = \varepsilon \psi_-\]

(14)

Using Eqs. 3 and 4 from the text, we can write four of the coefficients as position-dependent velocities:

\[
\hbar v_F(\phi_+|\sigma_x|\phi_-)_r \equiv iv_x(x), \quad \hbar v_F(\phi_-|\sigma_x|\phi_+)_r \equiv -iv_x(x), \\
\hbar v_F(\phi_+|\sigma_y|\phi_-)_r \equiv v_y(x), \quad \hbar v_F(\phi_-|\sigma_y|\phi_+)_r \equiv v_y(x).
\]

(15)

Note again that the position dependence of these velocities enters through the \(x\) dependence of \(V_0(x)\). With the further definition

\[
a_{\pm,+} \equiv -i\hbar v_F(\phi_\pm|\vec{\sigma} \cdot \vec{\nabla}|\phi_\pm)_r, = -i\hbar v_F(\phi_\pm|\sigma_x \partial_x|\phi_\pm)_r,
\]

(16)

where the second equality is a consequence of the fact that \(V_0\) depends on \(x\) but not on \(y\), we arrive at Eq. 5 of the text:

\[
\begin{pmatrix}
0 \\
-v_x \partial_x - iv_y \partial_y + a_{+,+}(r)
\end{pmatrix}
\begin{pmatrix}
\psi_+ \\
\psi_-
\end{pmatrix} = \varepsilon \begin{pmatrix}
\psi_+ \\
\psi_-
\end{pmatrix}.
\]