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Polar Kerr Effect and Time Reversal Symmetry Breaking in Bilayer Graphene

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The unique sensitivity of optical response to different types of symmetry breaking can be used to detect and identify spontaneously ordered many-body states in bilayer graphene. We predict a strong response at optical frequencies, sensitive to electronic phenomena at low energies, which arises because of nonzero interband matrix elements of the electric current operator. In particular, the polar Kerr rotation and reflection anisotropy provide fingerprints of the quantum anomalous Hall state and the nematic state, characterized by spontaneously broken time-reversal symmetry and lattice rotation symmetry, respectively. These optical signatures, which undergo a resonant enhancement in the near-infrared regime, lie well within reach of existing experimental techniques.

Optical experiments have been successfully used to probe diverse electronic phenomena in graphene [1]. For bilayer graphene (BLG), physical properties such as the gate tunable band gap [2,3] and the electron phonon coupling [4,5] were investigated with the help of infrared and optical spectroscopy. These techniques have also been used to probe interaction effects such as band renormalization [6,7] and exciton formation [8,9]. However, optical methods have not yet been employed to investigate strongly correlated states, which are expected to form in BLG at low energies [10–18]. This can be partly due to the low characteristic energy scales for these symmetry breaking states, estimated to be of order 1 meV [11], which lie far outside the range of characteristic energies probed in optical experiments. In this Letter, we point out that the problem of energy scales is offset by the unique sensitivity of optical response to broken symmetries, making these methods ideally suited to the investigation of the interacting ground state of BLG.

A large number of possible interacting phases have been proposed for BLG [10–18]. Recent compressibility and transport experiments on charge neutral, suspended, double gated bilayer graphene [19–21] appear to confirm the prediction of a nontrivial interacting ground state. The experimental data were argued [21] to be consistent with only two of the proposed phases: the quantum anomalous Hall phase (QAH) predicted in [13,14], and the nematic phase predicted in [15–17]. Both these phases are uniquely interesting phases. The QAH phase spontaneously breaks time-reversal symmetry (TRS) and exhibits quantum Hall effect at zero magnetic field, while the nematic state involves a distortion of the Dirac band structure that spontaneously breaks the exact rotational symmetry of the lattice. If either of these phases is confirmed in BLG, it would fulfill a long quest for an experimental realization of a QAH instability [22] (QAH phase) or a Pomeranchuk instability [23,24] (nematic phase).

The possible broken symmetries are expected to manifest themselves through characteristic transport properties such as a nonzero Hall response or anisotropy in longitudinal conductance [10–18]. Detecting these effects in transport experiments requires fabrication of samples of BLG with at least four contacts, which proves challenging in suspended BLG currently used in these experiments. However, optical experiments allow us to measure the ac conductivity in a contact-free manner. Furthermore, the ac conductivity shows distinctive signatures of broken symmetry. The polar Kerr effect, wherein linearly polarized light has its polarization axis rotated upon reflection, is a well known optical probe of the Hall conductivity. It has been used to probe quantum Hall states [25], and more recently has been applied to topological insulator thin films in the vicinity of a ferromagnet [26], and to $p + ip$ superconductors [27]. The Kerr effect is closely related to the Faraday effect, which has been measured for graphene in the quantum Hall regime [28]. However, unlike the Faraday effect, which requires breaking of either TRS or inversion symmetry, the polar Kerr effect can arise only if TRS is broken [29], hence offering a direct test of the QAH scenario for BLG.

As we show below, the QAH state exhibits an ac Hall conductivity that undergoes a resonant enhancement in the optical and near-infrared regime (see Fig. 1). The enhancement occurs because the microscopic current operator has interband matrix elements (Fig. 1 inset) corresponding to transitions from the low-energy bands 1, 2 to the high-energy bands $1', 2'$. As a result, the Kerr rotation is many orders of magnitude larger than that observed in $p$-wave superconducting materials [27], and lies well within reach of existing experimental techniques.

Optical methods can be used to probe domain formation expected to occur in the TRS breaking QAH phase. Since different domains will produce a Kerr rotation of opposite sign, the spatial domain structure can be directly imaged—a significant advantage over transport experiments, which
can only measure the net effect of all domains. For a nonfocused optical experiment, the effect of random domains will be to reduce the total Kerr angle by a factor $\frac{m}{C_{11}}$, where $m$ is the number of domains.

While the Kerr rotation allows us to test for TRS breaking, anisotropy in reflection allows us to test for rotation symmetry breaking. As we discuss below, this leads to a characteristic dependence of the reflection amplitude on the polarization angle of incident light which offers a way to test the nematic scenario for BLG [15,17].

Finally, we note that spontaneous symmetry breaking is only expected to occur below a critical temperature, estimated to be of order 1–10 K [11,21]. The optical signatures of interacting states will thus show a strong temperature dependence, and will vanish entirely above a critical temperature. This provides a way to distinguish spontaneously broken symmetries from explicitly symmetry breaking effects (e.g., magnetic impurities), which will not show any comparable temperature dependence.

Electron properties of a clean BLG are governed by a four-band Hamiltonian written for the four component wave function $\psi = (\psi_1, \psi_2, \psi_3, \psi_4)$, describing electron wave function on the sublattices $A, B$ and $A', B'$ on the two layers:

$$H(p) = \begin{bmatrix} 0 & t_p & 0 & 0 \\ t_p^* & 0 & E_0 & 0 \\ 0 & E_0 & 0 & t_p \\ 0 & 0 & t_p^* & 0 \end{bmatrix}. \quad \text{with } t_p = e^{-i\Delta_0 / \hbar}$$

with $t_p = t_0(1 + e^{-i\Delta / \hbar} + e^{-i\Delta / \hbar})$, where $t_0 = 3.1$ eV is the hopping amplitude, and $E_0$ is band gap parameter for the upper and lower bands. The quantity $t_p$ vanishes at the $K$ and $K'$ points, behaving as $vp_\perp$ near point $K$ and as $-vp_\perp$ near point $K'$, where $p_\perp = p_x \pm ip_y$.

The Hamiltonian (1) features four bands with energies

$$\epsilon^2(p) = |t_p|^2 + \frac{1}{2} E_0^2 \pm \frac{1}{2} E_0^2 \sqrt{1 + 4|t_p|^2/E_0^2}. \quad (2)$$

Near the points $K$ and $K'$, this gives two massless Dirac bands $\epsilon_{1,2}(p)$ that cross quadratically at zero energy, and two high-energy bands $\epsilon_{1',2'}(p) = \pm E_0$. The dispersion near $K$ and $K'$ can be obtained by expanding in small $t_p/E_0$, giving $\epsilon_{1,2} = \pm |t_p|^2/E_0 = \pm v^2 p^2/E_0$, $\epsilon_{1',2'} = \pm (E_0 + v^2 p^2/E_0)$.

We now consider the effect of interactions. Interactions can open a bulk band gap between bands 1 and 2 [10–14], resulting in a band structure of the form Fig. 1 (inset). One particularly interesting gapped state is the QAH state, [13,14], the mean field Hamiltonian of which we present below. To exhibit more clearly the block structure we reorder basis vectors by interchanging the components $\psi_2$ and $\psi_4$. In this representation, we obtain

$$H_K(p, \Delta) = \begin{bmatrix} \Delta & 0 & v p_+ & 0 \\ 0 & -\Delta & 0 & v p_- \\ v p_- & 0 & 0 & E_0 \\ 0 & v p_+ & E_0 & 0 \end{bmatrix}, \quad (3)$$

where $\Delta$ is the order parameter describing gap opening at the $K$ and $K'$ points, where $H_K(p, \Delta) = H_{K'}(-p, -\Delta)$. The other possible gapped states [10–12] have a similar mean field Hamiltonian, but the sign of $\Delta$ is distributed differently among the spins and valleys. We note that under time reversal, $H_K(\Delta) \otimes H_{K'}(-\Delta) \rightarrow H_{K'}(-\Delta) \otimes H_K(\Delta)$, so this phase breaks TRS. In consequence, the QAH state can exhibit a nonvanishing Hall conductance at zero magnetic field. However, the gap preserves the isotropy of the band structure. Thus, the QAH state must exhibit isotropic longitudinal conductivity.

Next, we discuss the relation between the Hall response in the QAH phase and the Kerr rotation. We consider an experimental setup where light is incident normally on a BLG sheet that is placed on a substrate. If the BLG sheet has a nonvanishing Hall conductance, then incident linearly polarized light will be reflected as elliptically polarized light, with the major axis of the ellipse rotated with respect to the incident polarization by the Kerr angle $\theta_K$. The standard formula relating the Kerr angle to the Hall conductivity and Kerr angle must be calculated afresh. This was accomplished using topological field theoretic methods in [26]. More

FIG. 1 (color online). Kerr angle (in units of fine structure constant $\alpha = e^2/\hbar c$) as a function of photon energy for BLG in the QAH phase. Note the resonant enhancement near $E_0 = 0.4$ eV, arising from direct transitions to the higher BLG bands, Eq. (12).
straightforwardly, the relationship may be obtained by solving the Maxwell equations on two sides of the BLG sheet and matching solutions at the boundary [30]. This leads to an expression in agreement with [26], which takes the form

$$\theta_K = \text{Re} \frac{-(8\pi/e)\sigma_{xx}}{1 - (\eta + \frac{\pi}{e}(\sigma_{xx} + i\sigma_{xy}))^2} = \frac{8\pi \text{Re} \sigma_{xx}}{c(n^2 - 1)},$$  \hspace{1cm} (4)$$

where $n$ is the refractive index of the substrate, which is taken to be real (complex case considered in [30]).

We now calculate the magnitude of the Kerr rotation, by evaluating the conductivity. The ac conductivity can be written using the Kubo formula as

$$\sigma_{xy} = \frac{e^2}{i\omega} \sum_{i,j,p} \langle i, p | V_{i,j} | j, p \rangle \langle j, p | V_{j,i} | i, p \rangle + \gamma (n_{i,p} - n_{j,p}),$$  \hspace{1cm} (5)$$

where $i$ and $j$ are band indices, $n_{i,p} = n(e_{i,p})$ is a Fermi function, and the sum over momenta $p$ stands for an integral. The velocity operators $V_{i}$ are defined as $V_{i} = \partial H(p) / \partial p_i$, and $\gamma$ describes the excited state lifetime.

We focus on the contributions which correspond to optical interband transitions between the massless low-energy bands ($i = 1, 2$), and the high-energy bands ($i = 1', 2'$), which are separated from the low-energy bands by the energy $E_0$. We focus on these transitions because they are of resonant character at a frequency close to the band separation energy $E_0$, and hence dominate the optical response. We now note that

$$\langle i, p | V_{i,j} | j, p \rangle \langle j, p | V_{j,i} | i, p \rangle = \text{Tr}(V_{i} \Pi_{i} V_{j} \Pi_{j}),$$

where $V_{i} \Pi_{i}$ projects onto the state in band $i$ with momentum $p$. Assuming we are at a temperature $T < E_0/k_B = 400K$, we obtain

$$\sigma_{xy}(\omega) = \frac{e^2}{i\omega} \int d^2 p \left[ \text{Tr}(V_{i} \Pi_{i} V_{j} \Pi_{j}) + \frac{\pi}{e} \frac{(p - e_{i,p} + i\gamma)}{\omega - e_{1'}(p) + e_{1}(p) + i\gamma} \right],$$  \hspace{1cm} (6)$$

where we used the relation $e_{x'} - e_{x} = e_{x'} - e_{x} = e_{x'} - e_{x}$. That follows from particle and hole symmetry of the Hamiltonian (3).

We evaluate the expression (6) for $p$ near point $K$ with the help of the projectors

$$\Pi_{1,2} = \frac{1}{2} \left( 1 \pm \left( \frac{h(p)}{\|h(p)\|} \right) \right), \quad \Pi_{1',2'} = \frac{1}{2} \left( \pm \hat{\sigma}_{x} \right),$$  \hspace{1cm} (7)$$

Here $\hat{\sigma}_{x}$ acts on the $B$ and $A'$ sublattices [lower right corner of the Hamiltonian in (3)] and $\Pi_{1',2'}$ projects on this subspace. Meanwhile, $\Pi_{1,2}$ project on the $A1$ and $B2$ sublattices [upper left corner of Eq. (3)], and $h(p)$ is the effective two band Hamiltonian for the massless Dirac states, which has eigenvalues $E(p) = \pm \|h(p)\|$. The trace over projectors takes the form

$$g_{a\beta}^{1'} = \text{Tr}(V_{a} \Pi_{1} V_{\beta} \Pi_{2}) = \langle 1' | V_{\beta} \Pi_{2} V_{a} | 1' \rangle$$

$$= \frac{1}{4} \left[ \nabla_{\beta t'}(p) \right]^{\dagger} \left( 1 - \frac{h(p)}{\|h(p)\|} \right) \left[ \nabla_{a t'}(p) \right] g_{a\beta}^{1'}$$

$$= \text{Tr}(V_{a} \Pi_{1} V_{\beta} \Pi_{2}) = \langle 2' | V_{\beta} \Pi_{1} V_{a} | 2' \rangle$$

$$= \frac{1}{4} \left[ \nabla_{a t'}(p) \right]^{\dagger} \left( 1 + \frac{h(p)}{\|h(p)\|} \right) \left[ \nabla_{\beta t'}(p) \right].$$  \hspace{1cm} (8)$$

Here $\nabla_a$ denotes $\partial / \partial p_{a}$. We now compute $h(p)$ by using second order perturbation theory in $v_F / E_0$, and obtain

$$h_{K}(p) = \left[ \begin{array}{cc} \Delta & v_F p_z^2 / E_0 \\ -\Delta & -\Delta \end{array} \right].$$  \hspace{1cm} (9)$$

$h_{K}(\Delta) = h_{K}(-\Delta)$. This result agrees with [31]. We substitute these two band Hamiltonian into Eq. (8) and obtain

$$g_{a\beta}^{1'} = g_{a\beta}^{2'} = \frac{1}{4} \left[ 1 - \frac{\Delta}{\|h(p)\|} \right] \nabla_{a t'}(p) \nabla_{\beta t'}(p)$$

$$+ \frac{1}{4} \left[ 1 + \frac{\Delta}{\|h(p)\|} \right] \nabla_{\beta t'}(p) \nabla_{a t'}(p),$$

where we suppressed the terms arising from off-diagonal parts of $h(p)$—these terms give zero upon integration over $d^2 p$. Hence, we find $g_{a\beta}^{1'} = -g_{a\beta}^{2'} = \frac{1}{2} iv_F^2 \Delta / \|h(p)\|$. We substitute these results into Eqs. (6) and (8), to obtain

$$\sigma_{xy}(\omega) = \int d^2 p \left[ \frac{A \Delta}{\omega + i\gamma - \Omega_p} + (\Omega_p - \Omega_p) \right],$$  \hspace{1cm} (11)$$

where $A = Ne^2 v_F^2 / ((2\pi)^2 \omega)$, where $N = 4$ is the number of spin/valley flavors, and $\Omega_p = e_2(p) - e_1(p)$.

We now specialize to optical frequencies $\omega \gg \Delta$, and also assume $\gamma \gg \Delta$. We approximate by taking $\|h(p)\| \approx v_F p^2 / E_0$ and $\Omega_p = E_0 + 2v_F^2 p^2 / E_0$, and perform the momentum integral in polar coordinates. The log divergence near $p^2 = 0$ is cut by $|\Delta|$, but there is no need for any high-energy cutoff. In this manner, we obtain the Hall conductivity

$$\sigma_{xy}(\omega) = \frac{Ne^2}{\hbar} \left[ \frac{E_0 \omega}{\omega + E_0 + i\gamma} \ln \left( \frac{E_0 + \omega + i\gamma}{2|\Delta|} \right) \right]$$

$$+ \frac{E_0}{\omega - E_0 + i\gamma} \ln \left( \frac{E_0 - \omega - i\gamma}{2|\Delta|} \right),$$  \hspace{1cm} (12)$$

There is also a contribution from $1 \rightarrow 2$ transitions, which may be evaluated in the two band model [18,32]. This contribution, extrapolated to optical frequencies $\omega \approx E_0 \gg \Delta$ is of order $(\Delta / |\Delta|)^3 v_F e^2 / \hbar$. This is smaller than the contribution (12) by a large factor

$$\frac{E_0}{\Delta} \ln \frac{E_0}{\Delta} \approx 1.$$  \hspace{1cm} (13)$$

Thus, the Hall conductivity at optical frequencies is dominated by transitions to the higher bands, necessitating our four-band analysis.
From the result Eq. (12) and the expression Eq. (4) we can extract the Kerr angle \( \theta_K \). We take \( n = 1.5 \), which describes SiO2 substrate, and take \( \Delta = 10^{-3} \) eV [11]. To estimate \( \gamma \), we use the electron lifetime \( \tau \approx 0.02 \) ps from [33], which corresponds to \( \gamma \approx 0.01 \) eV. In Fig. 1 we plot the resulting Kerr angle as a function of frequency using value \( \gamma = 0.05 \) eV (a conservative choice) and find a Kerr angle of order \( 10^{-3} \) radians. In optical experiments on cuprate materials, Kerr angles as small as \( 10^{-9} \) radians have been measured [27]. The 6 orders of magnitude larger Kerr rotation in the QAH phase should thus be comfortably within reach of experiments.

The nematic state [15–17] is another interesting ordered state proposed to explain the experiments [20,21]. This state is time-reversal invariant, featuring no Kerr rotation. Instead, it breaks rotation symmetry of graphene crystal lattice. The Hamiltonian for this state is

\[
H_K(p, \Delta) = \begin{bmatrix}
0 & \Delta & v p_+ & 0 \\
\Delta^* & 0 & 0 & v p_-
\end{bmatrix}
\begin{bmatrix}
0 \\
v p_- \\
v p_+ \\
E_0
\end{bmatrix}
\] (14)

After reduction to the two low-energy bands, it becomes

\[
h_K(p, \Delta) = \begin{bmatrix}
0 & \frac{v p_+^2}{E_0} + \Delta^* \\
\frac{v p_-^2}{E_0} + \Delta^* & 0
\end{bmatrix}
\] (15)

where \( H_K(p, \Delta) = H_K^*(-p, \Delta) \). This Hamiltonian describes splitting of the quadratic band crossing into two linear band crossings. The argument of the nematic order parameter \( \Delta \) specifies the orientation of the nematic axis, which is defined as the line joining the two linear band crossings. The nematic axis makes an angle \( \varphi = -\pi/2 + \arg(\Delta)/2 \) with respect to the \( p_z \) axis. The nematic state manifestly breaks the approximate rotation invariance of the low-energy band structure, which manifests itself in an anisotropic longitudinal conductivity. Writing \( \sigma(\theta) = \sigma_0 + 6\sigma(\theta) \), where \( \theta \) is the angle with respect to the \( x \) axis, we obtain an expression for the reflection amplitude \( r(\theta) \),

\[
r(\theta) = 1 - \frac{n}{n + 1} - \frac{8\pi}{c(n + 1)^2} \Delta \sigma(\theta)
\] (16)

For high frequencies \( \omega \gg \Delta \), we calculate using the formalism introduced above that

\[
\Delta \sigma(\theta) \sim \frac{i e^2}{h} \left| \frac{\Delta}{\omega} \right| \ln \frac{E_0 - \omega - i\gamma}{\Delta} \cos[2(\theta - \varphi)]
\] (17)

Again, this exceeds the anisotropy calculated in the two band model [18] by the large factor Eq. (13). We note that trigonal warping of the BLG band structure arising from higher neighbor hopping can also lead to a reflection anisotropy. However, these effects reflect the threefold rotation symmetry of the lattice. In contrast, the anisotropy resulting from formation of a nematic state exhibits a twofold rotation symmetry. The breaking of the exact lattice rotation symmetry can serve as diagnostic of the nematic state.

To conclude, optical experiments can be used to probe broken symmetries in BLG by measuring the conductivity in a contact-free manner. The polar Kerr effect, by providing a means for measuring Hall conductivity, can be used to detect the QAH phase. TRS breaking gapped states that do not display a Hall conductance [14] can also be probed using the Kerr effect, although for these states the Kerr angle will be smaller than that for the QAH state by the small parameter \( d/\lambda \), where \( d = 3 \) Å is the BLG interlayer spacing and \( \lambda \) is the wavelength of the light used in the experiment [34]. Nevertheless, this much weaker Kerr rotation will still be much larger than that measured in [27], and will be within reach of experiments. Meanwhile, the nematic scenario for BLG may be probed by looking for an angle dependence of the reflection amplitude, which provides a direct test of broken rotational symmetry.

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