

# Chern Number in a Band Structure

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Words	Equations
Berry connection / vec pot	$\vec{A}(k) = i \left\langle n(k) \left  \frac{\partial}{\partial \vec{k}} \right  n(k) \right\rangle$
Berry Phase	$\gamma = i \int d\vec{k} \cdot \vec{A}_n(k)$
Berry Curvature	$F_n = \left\langle \vec{\nabla}_k n \left  \times \right  \vec{\nabla}_k n \right\rangle = \vec{\nabla}_k \times \vec{A}(k)$
Chern Number	$\nu_n = \frac{1}{2\pi} \int_{BZ} d^2k F_n \in \mathbb{Z}$
TRS with real spin	$T = -i\sigma_y K = \begin{pmatrix} 0 & 1 \\ -1 & 0 \end{pmatrix} K$

## Outline of Talk

1. Describe what Chern number means and how it is calculated.
2. Model calculation with a single Dirac cone

- (a) Explain qualitatively why this is only half-quantized.
- (b) Explain what this means for when a gap closes and re-opens.
- (c) On hand:
  - i. Calculation for state existing at edge

### 3. Haldane Model

- (a) Introduce Hamiltonian
  - i. Full tight-binding first, without gapping terms.
  - ii. Sublattice asymmetry: show it breaks inversion but not TRS
  - iii. Second-nearest neighbor terms
- (b) Add periodic magnetic field
  - i. Show it breaks TRS, but not inversion (necessarily)
  - ii. Explain the complex hopping term qualitatively
- (c) Topological Phase Diagram
  - i. Plot the gap at the different K-points as a function of  $M \rightarrow$  show why chern # changes from 0 to 1 or -1
  - ii. Draw the entire phase diagram
- (d) On hand:
  - i. Quantitative explanation for complex hopping term

### 4. QAHE in magnetic TIs

- (a) Explain system:
  - i. Two surface states with finite tunneling between them (trivial mass gap)
  - ii. Write down Hamiltonian
  - iii. Introduce magnetic term, explain what it means
- (b) Show via same methods why it can produce finite chern # like in Haldane Model
- (c) On hand:
  - i. Van Vleck Paramagnetism
  - ii. Mean field theory standard
  - iii. Mean field theory TI?

### 5. On hand for calculating quantized conductance

- (a) Anomalous velocity?
- (b) From anomalous velocity up to Hall conductivity
- (c) Influence of Dissipative Channel

## 1 Single Dirac Cone

$$H(k) = \sum_i d_k(k) \sigma_i \tag{1}$$

$$\sigma_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \sigma_2 = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \sigma_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

If  $d_3 = M \neq 0$ , then there will be a gap. Showing TRS is broken with a mass gap:

$$THT^{-1} = -d_x\sigma_x - d_y\sigma_y - M\sigma_z \neq H(-k) \quad (2)$$

$$H(-k) = -d_x\sigma_x - d_y\sigma_y + M\sigma_z \quad (3)$$

Generally we need  $d_i(-k) = -d_i(k)$  to preserve TRS.

The normalized, orthogonal wave-functions of this two-level system are the following, including a helicity parameter  $\alpha = \pm 1$ :

$$\psi_+ = \frac{1}{\sqrt{2d(d+d_3)}} \begin{pmatrix} d_3 + d \\ (d_1 - \alpha id_2) \end{pmatrix} \quad (4)$$

$$\psi_- = \frac{1}{\sqrt{2d(d-d_3)}} \begin{pmatrix} d_3 - d \\ (d_1 - \alpha id_2) \end{pmatrix} \quad (5)$$

$$\psi_- = \frac{1}{\sqrt{2\sqrt{m^2+k^2}(\sqrt{m^2+k^2}-m)}} \begin{pmatrix} m - \sqrt{m^2+k^2} \\ k_x - \alpha ik_y \end{pmatrix} \quad (6)$$

$$= \frac{1}{\sqrt{B_-(k)}} \begin{pmatrix} m - \sqrt{m^2+k^2} \\ ke^{i\alpha\theta_k} \end{pmatrix} \quad (7)$$

The Berry connection is

$$\vec{A}_-(k) = i \langle \psi_-(k) | \vec{\nabla}_k | \psi_-(k) \rangle \quad (8)$$

$$= \frac{-\alpha}{2d(d-d_3)} (d_2 \vec{\nabla}_k d_1 - d_1 \vec{\nabla}_k d_2) \quad (9)$$

$$A_x = \frac{-\alpha k_y}{2\sqrt{k^2+m^2}(\sqrt{k^2+m^2}-m)} \quad (10)$$

$$A_y = \frac{\alpha k_x}{2\sqrt{k^2+m^2}(\sqrt{k^2+m^2}-m)} \quad (11)$$

$$F_z = \frac{1}{2(k^2+m^2)^{\frac{3}{2}}} \alpha m \quad (12)$$

$$\nu = \frac{\alpha}{2\pi} \int_{-\infty}^0 F_{xy} d^2k = \frac{\text{sign}(m\alpha)}{2} \quad (13)$$

This is for ‘infinite’ band width of a Dirac system. However, in a real situation, finite band-width gives two gaps, each of which should roughly contribute  $\pm \frac{1}{2}$  to the Chern number.

## 2 QAHE in TI

### 2.1 SS Hamiltonian in thick limit

Each surface state has a Hamiltonian

$$H_{sf}^L = v_F \vec{k} \times \vec{\sigma} = \text{Liv}_F \begin{pmatrix} 0 & k_- \\ -k_+ & 0 \end{pmatrix} \quad (14)$$

where prefactor  $L = \pm 1$  is for the top surface and bottom surface, and  $k_{\pm} = k_x \pm ik_y$ . Dispersion is

$$E_{sf} = v_F |k| \quad (15)$$

## 2.2 SS Hamiltonians with coupling

Coupling the surfaces results in a mass term.

$$H_{two} = \begin{pmatrix} 0 & iv_F k_- & m_k^* & 0 \\ -iv_F k_+ & 0 & 0 & m_k^* \\ m_k & 0 & 0 & -iv_F k_- \\ 0 & m_k & iv_F k_+ & 0 \end{pmatrix}, \quad \psi = \begin{pmatrix} t \uparrow \\ t \downarrow \\ b \uparrow \\ b \downarrow \end{pmatrix} \quad (16)$$

$$\text{TRS } T = \begin{pmatrix} 0 & -1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & -1 & 0 \end{pmatrix} K \text{ shows that } THT^{-1} = H.$$

In the limit that  $v_F k \ll m_k$ , then  $m_k \approx m_0 + m_1 k^2$  and one can rewrite the Hamiltonian in a new basis

$$H_{two} = \begin{pmatrix} m_k & iv_F k_- & 0 & 0 \\ -iv_F k_+ & -m_k & 0 & 0 \\ 0 & 0 & m_k & -iv_F k_+ \\ 0 & 0 & iv_F k_- & -m_k \end{pmatrix}, \quad \psi = \begin{pmatrix} + \uparrow \\ - \downarrow \\ + \downarrow \\ - \uparrow \end{pmatrix} \quad (17)$$

which is block diagonal yet again, with each block being an equal superposition of top and bottom surfaces, and they are both gapped:

$$E = \pm \sqrt{m_k^2 + v_F^2 k^2} \quad (18)$$

Note that this Hamiltonian is still maintains TRS, since flipping spin and complex conjugating (flipping momentum) does not change it.

## 2.3 Add magnetic term

A magnetic exchange term to a “bath” of localized spins inserts itself as a sort of Zeeman coupling for each spin. Therefore

$$H_M = \begin{pmatrix} gM & 0 & 0 & 0 \\ 0 & -gM & 0 & 0 \\ 0 & 0 & -gM & 0 \\ 0 & 0 & 0 & gM \end{pmatrix}, \quad \psi = \begin{pmatrix} \uparrow \\ \downarrow \\ \downarrow \\ \uparrow \end{pmatrix} \quad (19)$$

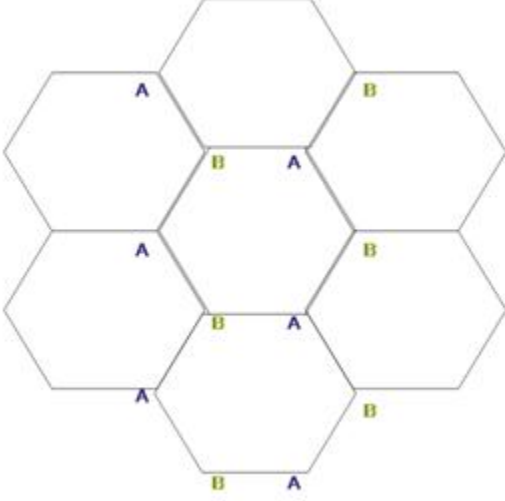
which DOES break TRS, since  $THT^{-1} = -H$ .

One can re-write the block-diagonal Hamiltonian as

$$H = \begin{pmatrix} h_k + (m_k + gM) \sigma_z & 0 \\ 0 & h_k^* + (m_k - gM) \sigma_z \end{pmatrix}, \quad \psi = \begin{pmatrix} + \uparrow \\ - \downarrow \\ + \downarrow \\ - \uparrow \end{pmatrix} \quad (20)$$

which shows two independent sections. When  $g|M| \equiv 0$ , then the system has TRS so  $\nu \equiv 0$ . As the exchange field  $|M|$  is increased, ONE of the gaps closes when  $g|M| = m_0$  and then reopens with an opposite-sign mass parameter. This means that the Chern number has changed by 1, so  $\nu = \pm 1$ , with the sign determined by the sign of  $M$ .

### 3 Haldane Model



#### 3.1 Basic Graphene

Tight binding for Graphene. First, lattice vectors, assume A and B are horizontal from each other:  $\vec{r} = a\hat{x}$

$$\vec{r}_{a \rightarrow b} = a\hat{x} \quad (21)$$

$$= -\frac{1}{2}a\hat{x} \pm \frac{\sqrt{3}}{2}a\hat{y} \quad (22)$$

$$\vec{r}_{b \rightarrow a} = -a\hat{x} \quad (23)$$

$$= \frac{1}{2}a\hat{x} \pm \frac{\sqrt{3}}{2}a\hat{y} \quad (24)$$

But either way:

$$H = t_1 \sum_{\langle i,j \rangle} e^{i\vec{k} \cdot \vec{r}_{ij}} c_i^\dagger c_j + \sum_{i=a,b} \epsilon_i c_i^\dagger c_i \quad (25)$$

Find the zero-energy spots when  $m = 0$ . We know it is at the  $K$  point, which we can find by setting  $k_x = 0$ :

$$1 + 4 \cos^2\left(\frac{\sqrt{3}}{2}ak_y\right) + 4 \cos\left(\frac{\sqrt{3}}{2}ak_y\right) \cos\left(\frac{1}{2}ak_x\right) = 0 \quad (26)$$

$$1 + 4 \cos^2\left(\frac{\sqrt{3}}{2}ak_y\right) + 4 \cos\left(\frac{\sqrt{3}}{2}ak_y\right) = 0 \quad (27)$$

$$\left(1 + 2 \cos\left(\frac{\sqrt{3}}{2}ak_y\right)\right)^2 = 0 \quad (28)$$

$$\cos\left(\frac{\sqrt{3}}{2}ak_y\right) = -\frac{1}{2} \quad (29)$$

$$\frac{\sqrt{3}}{2}ak_y = \pm \frac{2\pi}{3} \quad (30)$$

$$K_{\pm} = \pm \frac{4\pi}{3\sqrt{3}a} \hat{k}_y \quad (31)$$

Now simplify the Hamiltonian:

$$H = t_1 \begin{pmatrix} 0 & e^{-iak_x} + e^{i\frac{1}{2}ak_x} \left( e^{i\frac{\sqrt{3}}{2}ak_y} + e^{-i\frac{\sqrt{3}}{2}ak_y} \right) \\ e^{iak_x} + e^{-i\frac{1}{2}ak_x} \left( e^{i\frac{\sqrt{3}}{2}ak_y} + e^{-i\frac{\sqrt{3}}{2}ak_y} \right) & 0 \end{pmatrix} \quad (32)$$

$$= t_1 \begin{pmatrix} 0 & e^{-ia\delta_x} + e^{i\frac{1}{2}a\delta_x} \left( e^{\pm\frac{2\pi}{3}i + \frac{\sqrt{3}}{2}ia\delta_y} + e^{\mp\frac{2\pi}{3}i - \frac{\sqrt{3}}{2}ia\delta_y} \right) \\ e^{ia\delta_x} + e^{-i\frac{1}{2}a\delta_x} \left( e^{\pm\frac{2\pi}{3}i + \frac{\sqrt{3}}{2}ia\delta_y} + e^{\mp\frac{2\pi}{3}i - \frac{\sqrt{3}}{2}ia\delta_y} \right) & 0 \end{pmatrix} \quad (33)$$

$$= t_1 \begin{pmatrix} 0 & e^{-ia\delta_x} + e^{i\frac{1}{2}a\delta_x} \left( e^{\pm\frac{2\pi}{3}i} e^{\frac{\sqrt{3}}{2}ia\delta_y} + e^{\mp\frac{2\pi}{3}i} e^{-\frac{\sqrt{3}}{2}ia\delta_y} \right) \\ e^{ia\delta_x} + e^{-i\frac{1}{2}a\delta_x} \left( e^{\pm\frac{2\pi}{3}i} e^{\frac{\sqrt{3}}{2}ia\delta_y} + e^{\mp\frac{2\pi}{3}i} e^{-\frac{\sqrt{3}}{2}ia\delta_y} \right) & 0 \end{pmatrix} \quad (34)$$

$$\approx t_1 \begin{pmatrix} 0 & 1 - ia\delta_x + \left( 1 + i\frac{a}{2}\delta_x \right) \left( -1 + \alpha\frac{3}{2}a\delta_y \right) \\ 1 + ia\delta_x + \left( 1 - i\frac{a}{2}\delta_x \right) \left( -1 + \alpha\frac{3}{2}a\delta_y \right) & 0 \end{pmatrix} \quad (35)$$

$$\approx t_1 \begin{pmatrix} 0 & 1 - ia\delta_x + \left( -1 + \alpha\frac{3}{2}a\delta_y - i\frac{a}{2}\delta_x \right) \\ 1 + ia\delta_x + \left( -1 + \alpha\frac{3}{2}a\delta_y + i\frac{a}{2}\delta_x \right) & 0 \end{pmatrix} \quad (36)$$

$$\approx \frac{3}{2}at_1 \begin{pmatrix} 0 & \alpha\delta_y - i\delta_x \\ \alpha\delta_y + i\delta_x & 0 \end{pmatrix} \quad (37)$$

where  $\frac{3}{2}at_1$  is the fermi velocity  $v_F$  up to a factor of  $\hbar$ . Note that  $\alpha = \pm 1$  for  $K/K'$ . This leads to the familiar dispersion relation

$$E = \pm \hbar v_F |\delta| \quad (38)$$

## 3.2 Symmetries

### 3.2.1 Inversion

$\vec{r} \rightarrow -\vec{r}$ . This swaps the sublattices. It clearly does not affect the n.n. off-diagonal components, but it does affect the on-site energies. Breaking inversion opens a band gap.

### 3.2.2 TRS

$\hat{T} = iK\sigma_y$ , so  $T^2 = -1$ . The Hamiltonian, as is, has TRS, i.e.

$$THT^{-1} = H \quad (39)$$

A term in the Hamiltonian that breaks TRS may open up a band gap.

## 3.3 Second Nearest Neighbor

First, need the vectors:

$$\vec{r}_{a \rightarrow a} = \pm \sqrt{3}a\hat{y} \quad (40)$$

$$= \pm \frac{3}{2}a\hat{x} \pm \frac{\sqrt{3}}{2}a\hat{y} \quad (41)$$

$$\vec{r}_{b \rightarrow b} = \pm \sqrt{3}a\hat{y} \quad (42)$$

$$= \pm \frac{3}{2}a\hat{x} \pm \frac{\sqrt{3}}{2}a\hat{y} \quad (43)$$

The perturbation to the model is

$$H_2 = t_2 \sum_{\langle\langle i,j \rangle\rangle} e^{i\vec{k}\cdot\vec{r}_{ij}} c_i^\dagger c_j \quad (44)$$

which is a diagonal term. As written, this term does not break either symmetry of the lattice, and so does not open a gap. It contributes the following (this is at  $K^+$ ):

$$H_{nnn} = 2t_2 \left( \cos(\sqrt{3}ak_y) + \cos\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y\right) + \cos\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y\right) \right) \quad (45)$$

$$= 2t_2 \left( \cos\left(\sqrt{3}a\left(\frac{4\pi}{3\sqrt{3}}\frac{1}{a} + \delta_y\right)\right) + \cos\left(\frac{3}{2}a\delta_x - \frac{\sqrt{3}}{2}\left(\frac{4\pi}{3\sqrt{3}}\frac{1}{a} + \delta_y\right)\right) + \cos\left(\frac{3}{2}a\delta_x + \frac{\sqrt{3}}{2}\left(\frac{4\pi}{3\sqrt{3}}\frac{1}{a} + \delta_y\right)\right) \right) \quad (46)$$

$$= 2t_2 \left( \cos\left(\frac{4\pi}{3} + \sqrt{3}a\delta_y\right) + \cos\left(\frac{3}{2}a\delta_x - \frac{2\pi}{3} - \frac{\sqrt{3}}{2}a\delta_y\right) + \cos\left(\frac{3}{2}a\delta_x + \frac{2\pi}{3} + \frac{\sqrt{3}}{2}a\delta_y\right) \right) \quad (47)$$

And the key is that at  $\delta_{x,y} = 0$ :

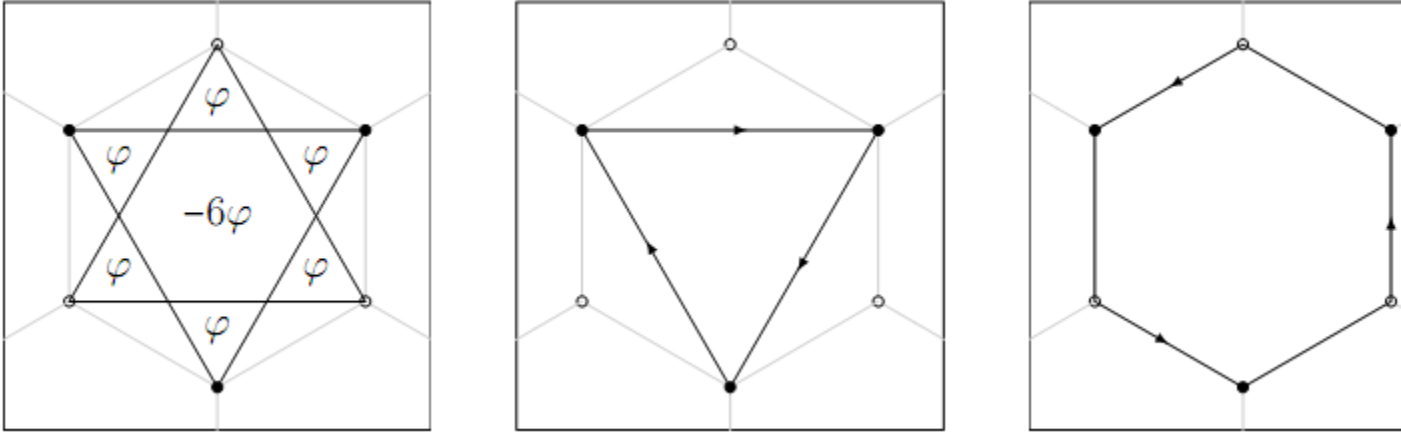
$$H_{nnn}(0,0) = 2t_2 \left( \cos\left(\frac{4\pi}{3}\right) + \cos\left(-\frac{2\pi}{3}\right) + \cos\left(\frac{2\pi}{3}\right) \right) \quad (48)$$

$$= 2t_2 \left( \cos\left(\frac{4\pi}{3}\right) + 2\cos\left(\frac{2\pi}{3}\right) \right) \quad (49)$$

$$= -3t_2 \quad (50)$$

which is just a chemical potential shift ( $K^\pm$  is the same since any  $\pm$  happens in front of a  $\delta_y$ ) and has no consequence. When  $\delta_{x,y} \neq 0$ , then this term renormalizes the Fermi velocity, and will be different for electrons and holes, but doesn't make any fundamental change in the band structure.

### 3.4 Haldane Term



Now introduce a periodic magnetic flux into the unit cell and choose a gauge so that only the nnn hoppings develop a phase.

$$H_2 = t_2 \sum_{\langle\langle i,j \rangle\rangle} e^{i\phi_{ij}} e^{i\vec{k}\cdot\vec{r}_{ij}} c_i^\dagger c_j \quad (51)$$

Be careful about minus signs!

$$H_{nnn,a} = 2t_2 \left( \cos(\sqrt{3}ak_y + \phi) + \cos\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y + \phi\right) + \cos\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y + \phi\right) \right) \quad (52)$$

$$H_{nnn,b} = 2t_2 \left( \cos(\sqrt{3}ak_y - \phi) + \cos\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y - \phi\right) + \cos\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y - \phi\right) \right) \quad (53)$$

Recall that

$$\cos(a \pm b) = \cos a \cos b \mp \sin a \sin b \quad (54)$$

$$H_{nnn,a} = 2t_2 \cos \phi \left( \cos(\sqrt{3}ak_y) + \cos\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y\right) + \cos\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y\right) \right) \quad (55)$$

$$-2t_2 \sin \phi \left( \sin(\sqrt{3}ak_y) + \sin\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y\right) + \sin\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y\right) \right) \quad (56)$$

$$H_{nnn,b} = 2t_2 \cos \phi \left( \cos(\sqrt{3}ak_y) + \cos\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y\right) + \cos\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y\right) \right) \quad (57)$$

$$+2t_2 \sin \phi \left( \sin(\sqrt{3}ak_y) + \sin\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y\right) + \sin\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y\right) \right) \quad (58)$$

So the mass-inducing terms are

$$H_{nnn}(0,0,a) = -2t_2 \sin \phi \left( \sin(\sqrt{3}ak_y) + \sin\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y\right) + \sin\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y\right) \right) \quad (59)$$

$$H_{nnn}(0,0,b) = 2t_2 \sin \phi \left( \sin(\sqrt{3}ak_y) + \sin\left(\frac{3}{2}ak_x - \frac{\sqrt{3}}{2}k_y\right) + \sin\left(\frac{3}{2}ak_x + \frac{\sqrt{3}}{2}k_y\right) \right) \quad (60)$$

So

$$H_a - H_b = 4t_2 \sin \phi \left( \sin\left(\frac{4\pi}{3}\right) + \sin\left(\frac{3}{2}a\delta_x - \frac{\sqrt{3}}{2}a\delta_y - \frac{2\pi}{3}\right) + \sin\left(\frac{3}{2}a\delta_x + \frac{\sqrt{3}}{2}a\delta_y + \frac{2\pi}{3}\right) \right) \quad (61)$$

$$= 4t_2 \sin\left(\frac{4\pi}{3}\right) \sin(\phi) \quad (62)$$

$$= -2\sqrt{3}t_2 \sin \phi \quad (63)$$

Note that the sign of this term in the Hamiltonian is **DEPENDENT** on the valley since each valley has a different helicity.

$$\textcircled{K} \rightarrow M - t_2 \sin \phi \quad (64)$$

$$\textcircled{K'} \rightarrow M + t_2 \sin \phi \quad (65)$$

$$\textcircled{K} \rightarrow \nu = \frac{\text{sign}(M - t_2 \sin \phi)}{2} \quad (66)$$

$$\textcircled{K'} \rightarrow \nu = \frac{-\text{sign}(M + t_2 \sin \phi)}{2} \quad (67)$$

See that when  $M$  goes from  $+\infty$  to  $< t_2 \sin \phi$ , the gap at  $K'$  stays open always and at  $K$  it closes and reopens, with the parameter changing sign. This means the total Chern # much change by an integer 1, and we know that at  $M \gg t_2 \sin \phi$  it is trivial due to the system appearing like an atomic system.

## 4 Supplementary Calculations

### 4.1 Edge Modes at Mass Inversion Interface

Hamiltonian of gapped Dirac system with mass that depends on the  $y$  coordinate.

$$H(y) = -i\frac{\partial}{\partial x}\sigma_x - i\frac{\partial}{\partial y}\sigma_y + m(y)\sigma_z \quad (68)$$

and look for a separable solution

$$\psi(x, y) = \phi_x(x)\phi_y(y) \quad (69)$$



Now,  $\phi_x = e^{ik_x x} \begin{pmatrix} 1 \\ -1 \end{pmatrix}$ , whereas

$$\phi_y = e^{-\int_0^y m(y') dy'} \quad (70)$$

Plugging in  $\phi_y \phi_x$ , one gets a function of  $\phi_x$  and  $m$  to find the exact spinor for  $\phi_x$ . Then you can solve for the plane-wave part which gives a chiral mode.

## 4.2 Phase From Magnetic Flux

Phase acquired from a path with a vector potential:

$$\phi = \frac{e}{\hbar} \int \vec{A} \cdot d\vec{r} \quad (71)$$

Which is Gauge dependent if not on a closed loop:

$$\vec{A}' = \vec{A} + \vec{\nabla}\chi \rightarrow \phi' \neq \phi \quad (72)$$

But if you have a closed loop, then you have a fixed quantity:

$$\varphi = \frac{e}{\hbar} \oint \vec{A} \cdot d\vec{r} = \frac{e}{\hbar} \iint \vec{B} \cdot d\vec{s} \quad (73)$$

And the complex phase around the loop is related to this phase. You can put this phase on whatever hopping term you like, due to the flexibility of gauge transformations.

## 4.3 Adiabatic Continuity

General Hamiltonian and solution to id:

$$\hat{H}(t)\psi_n(t) = E_n(t)\psi_n(t) \quad (74)$$

$$\Psi(t) = \sum_n c_n(t)\psi_n(t)e^{i\theta_n(t)} \quad (75)$$

$$\theta_n(t) = -\frac{1}{\hbar} \int_0^t E_n(t') dt' \quad (76)$$

Substitute  $\Psi$  into  $\hat{H}(t)$  and you get

$$i\hbar \sum_n \left( \dot{c}_n \psi_n + c_n \dot{\psi}_n + i c_n \psi_n \dot{\theta}_n \right) e^{i\theta_n} = \sum_n c_n \hat{H} \psi_n e^{i\theta_n} \quad (77)$$

but  $\dot{\theta}_n = -\frac{E}{\hbar}$  so

$$\sum_n \dot{c}_n \psi_n e^{i\theta_n} = - \sum_n c_n \dot{\psi}_n e^{i\theta_n} \quad (78)$$

now inner product with an arbitrary eigenfunction  $\psi_m$  and you get

$$\dot{c}_m(t) = - \sum_n c_n \langle \psi_m | \dot{\psi}_n \rangle e^{i(\theta_n - \theta_m)} \quad (79)$$

then calculating the differentiated inner product you get the exact form

$$\dot{c}_m(t) = -c_m \langle \psi_m | \dot{\psi}_m \rangle - \sum_{n \neq m} c_n \frac{\langle \psi_m | \dot{\hat{H}} | \psi_n \rangle}{E_n - E_m} e^{i(\theta_n - \theta_m)} \quad (80)$$

but if  $\frac{\partial}{\partial t} \hat{H} \ll 1$ , then the second term drops out. This means that

$$c_m(t) = c_m(0) e^{-\int_0^t \langle \psi_m | \dot{\psi}_m \rangle dt'} = c_m(0) e^{i\gamma_m(t)} \quad (81)$$

so that there is a dynamical phase factor known as the geometrical phase, but the amplitudes remain constant. Thus, a particle remains in the  $n^{\text{th}}$  eigenstate at all times if the Hamiltonian varies slowly.

## 4.4 Van Vleck Paramagnetism

Assume an insulator with no net magnetism, i.e.

$$\mu_z \cdot B = 0 \quad (82)$$

so that there is no first-order correction to the system. But then let's say that

$$\langle c|S_z|v\rangle \neq 0 \quad (83)$$

This means that with perturbation theory in a field  $\mu B \ll E_G$ , one gets a wave-function

$$\psi'_v = \psi_v + \frac{B}{E_G} \langle c|S_z|v\rangle \psi_c \quad (84)$$

And now we have a moment

$$\langle 0'|S_z|0'\rangle \approx \frac{2B}{E_G} |\langle c|S_z|v\rangle|^2 \quad (85)$$

So that in the presence of an applied field the magnetization is

$$\frac{M}{N} = \frac{2B}{E_G} |\langle c|S_z|v\rangle|^2 \quad (86)$$

And for a band insulator with more than one quantum number  $k$  then

$$\chi = 2 \sum_{(c,v),k} \frac{|\langle c|S_z|v\rangle|^2}{E_{c,k} - E_{v,k}} \quad (87)$$

In typical systems, these terms are very small. What happens in a TI? Well, there is a point in the TI band structure where the gap comes close to closing and at the same time this causes a maximization of the second order effects.

## 4.5 Mean Field Theory

### 4.5.1 Normal FM

$$H = -J \sum_{\langle i,j \rangle} s_i s_j - h \sum_i s_i \quad (88)$$

so define the mean field  $m_i = \langle s_i \rangle$  and rewrite

$$H = -J \sum_{\langle i,j \rangle} (m_i + \delta s_i)(m_j + \delta s_j) - h \sum_i s_i \quad (89)$$

$$H \approx H^{MF} \quad (90)$$

$$H^{MF} = -J \sum_{\langle i,j \rangle} (m_i m_j + m_i \delta s_j + m_j \delta s_i) - h \sum_i s_i \quad (91)$$

$$= -J \sum_{\langle i,j \rangle} (m^2 + 2m(s_i - m)) - h \sum_i s_i \quad (92)$$

and  $\sum_{\langle i,j \rangle} = \frac{1}{2} \sum_i \sum_{j \in nn(i)}$  and thus  $\sum_{j \in nn(i)} = x$  which is the number of nearest-neighbor spins. Therefore

$$H^{MF} = -\frac{1}{2} J m^2 N x - (h + m J x) \sum_i s_i \quad (93)$$

This gives us a partition function

$$Z = e^{-\frac{J m^2 N x}{2\beta}} \left[ 2 \cosh \left( \frac{h + m J x}{k_B T} \right) \right]^N \quad (94)$$

which gives a Free Energy

$$F = -k_B T \ln Z \sim N k_B T \ln \left[ 2 \cosh \left( \frac{h + m J x}{k_B T} \right) \right] \quad (95)$$

$$M = -\frac{\partial F}{\partial H} \sim \tanh \left( \frac{h + m J x}{k_B T} \right) \quad (96)$$

## 4.6 Lagrangian for wave-packet

Marder, section 16.4, page 425.

$\mathcal{L} = \langle w | i\hbar \frac{\partial}{\partial t} | w \rangle - \langle w | \vec{H} - eV | w \rangle$ , assuming no magnetic fields so  $\mathcal{H} = \frac{p^2}{2m} + U(r)$  and  $|w\rangle = \frac{1}{\sqrt{N}} \sum_k w_{kk_c} e^{-i\vec{k} \cdot \vec{r}_c} \psi_k$  and  $w_{kk_c} = e^{i(k-k_c) \cdot \vec{A}_{berry}}$  and  $\mathcal{H}\psi_k = \mathcal{E}_k \psi_k$ .

The Lagrangian after math simplifies to

$$\mathcal{L} = \hbar \vec{k}_c \cdot \dot{\vec{r}}_c + \hbar \dot{\vec{k}}_c \cdot \vec{A}_{berry} - \mathcal{E}_{k_c} - eV(\vec{r}_c) \quad (97)$$

$$\frac{\partial \mathcal{L}}{\partial \vec{r}_c} = \frac{d}{dt} \frac{\partial \mathcal{L}}{\partial \dot{\vec{r}}_c} \quad \text{and} \quad \frac{\partial \mathcal{L}}{\partial \vec{k}_c} = \frac{d}{dt} \frac{\partial \mathcal{L}}{\partial \dot{\vec{k}}_c} \quad (98)$$

$$-e\vec{E} = \frac{d}{dt} (\hbar \dot{\vec{k}}_c) \quad (99)$$

$$\hbar \dot{\vec{r}}_c + \frac{\partial \mathcal{E}_k}{\partial \vec{k}_c} = \frac{d}{dt} \frac{\partial}{\partial \dot{\vec{k}}_c} (\hbar \dot{\vec{k}}_c \cdot \vec{A}_{berry}) \quad (100)$$

and eventually with a bunch of vector identities, etc.

$$\dot{\vec{k}} = -\frac{e}{\hbar} \vec{E} \quad (101)$$

$$\dot{\vec{r}} = \frac{1}{\hbar} \frac{\partial \mathcal{E}_k}{\partial \vec{k}} - \dot{\vec{k}} \times \vec{F} \quad (102)$$

## 4.7 Anomalous Velocity

Semiclassical calculations (or Kubo formula, or other means), give the following two equations of motion for a wave-packet when the external magnetic field is zero:

$$\dot{\vec{k}} = -\frac{e}{\hbar} \vec{E} \quad (103)$$

$$\dot{\vec{r}} = \frac{1}{\hbar} \frac{\partial \mathcal{E}_k}{\partial \vec{k}} - \dot{\vec{k}} \times \vec{F} \quad (104)$$

while if the system is in a band gap then  $\int_{BZ} \frac{\partial E(k)}{\partial \vec{k}} \equiv 0$ , e.g. when integrating over completely filled bands. As such, I pre-emptively remove that term from the calculation (reader can verify for him/herself that this is ok). So, substituting, we get

$$\dot{\vec{r}} = v(k) = \frac{e}{\hbar} \vec{E} \times \vec{F} \quad (105)$$

For a 2D system,  $\vec{F} = F_z \hat{z}$ . To calculate the conductivity one needs

$$\vec{j} = \hat{\sigma} \vec{E} \quad (106)$$

so

$$j_y = \sigma_{xy} E_x \quad (107)$$

then, noting that the distribution function  $g$  is roughly the Fermi function  $g \sim f$  in this context.

$$j_y = e \iint \frac{d^2k}{4\pi^2} \vec{v}(k) g \quad (108)$$

$$= \frac{e^2}{h} \frac{1}{2\pi} \iint d^2k \vec{E} \times \vec{F} f \quad (109)$$

$$= \frac{e^2}{h} \frac{1}{2\pi} E_x \iint d^2k F_z \quad (110)$$

$$\sigma_{xy} = \frac{dj_y}{dE_x} \quad (111)$$

$$= \frac{e^2}{h} \frac{1}{2\pi} \iint d^2k F_z \quad (112)$$

$$= \frac{e^2}{h} \nu \quad (113)$$

$$\nu = \frac{1}{2\pi} \iint d^2k F_z \in \mathbb{Z} \quad (114)$$

Normally,

$$g \approx f - \tau e \frac{\partial f}{\partial \mu} \vec{v}_{\vec{k}} \cdot \vec{E} \quad (115)$$

and when you take the derivative  $\frac{dj_y}{dE_x}$  the first term drops out since there is no electric field there. However, when inside a gap and with finite Berry curvature,  $\vec{v}(k) \frac{\partial f}{\partial \mu} = 0$  and the only remaining term is the one with the Berry curvature multiplying the general Fermi function.

## 4.8 QHE and dissipative channels

$$\sigma_{xy} = \sigma_{QH} + \sigma_{dis} = \sigma_{QH} \quad (116)$$

$$\sigma_{xx} = \sigma_{QH} + \sigma_{dis} \approx \sigma_{dis} \quad (117)$$

so

$$\sigma = \begin{pmatrix} \sigma_d & \sigma_Q \\ -\sigma_Q & \sigma_d \end{pmatrix} \quad (118)$$

$$\rho = \frac{1}{\sigma_d^2 + \sigma_Q^2} \begin{pmatrix} \sigma_d & -\sigma_Q \\ \sigma_Q & \sigma_d \end{pmatrix} \quad (119)$$

so

$$\rho_{xx} = \frac{\sigma_d}{\sigma_d^2 + \sigma_Q^2} \approx \frac{h}{e^2} \frac{\xi}{\xi^2 + 1} \quad (120)$$

$$\rho_{xy} = \frac{\sigma_Q}{\sigma_d^2 + \sigma_Q^2} \approx \frac{h}{e^2} \frac{1}{\xi^2 + 1} \quad (121)$$

So as the dissipative channel's conductance goes away, the dissipating resistance goes to zero, and the Hall resistance approaches the quantized value.