

PHYS 525
Solution to HW 3

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PROBLEM 1

a) A possible choice of anticommuting hermitian 4×4 matrices reads

$$\Gamma_1 = \tau_z \sigma_x, \quad \Gamma_2 = \tau_z \sigma_y, \quad \Gamma_3 = \tau_z \sigma_z, \quad \Gamma_4 = \tau_x \sigma_0, \quad \Gamma_5 = \tau_y \sigma_0. \quad (1)$$

Here and hereafter σ_0 is used to denote a 2×2 unit matrix and we use the tensor product notation with \otimes sign omitted for the sake of brevity. The matrices defined above are often called Dirac Γ matrices and are used to formulate the relativistic version of the Schrödinger equation.

b) The calculation of the energy spectrum follows as in the 2×2 case discussed in class. We square the Hamiltonian

$$H^2 = \sum_{i,j} \underbrace{d_i(\mathbf{k}) d_j(\mathbf{k})}_{\text{symmetric in } i \leftrightarrow j} \underbrace{\Gamma_i \Gamma_j}_{\frac{1}{2} \{\Gamma_i, \Gamma_j\}} = |\mathbf{d}(\mathbf{k})|^2 \mathbb{I}$$

Taking a square root then gives, as before

$$E_{\pm}(\mathbf{k}) = \pm |\mathbf{d}(\mathbf{k})|. \quad (2)$$

c) Extending the construction in part (a), a possible choice of 8×8 matrices reads

$$\rho_z \tau_z \sigma_x, \quad \rho_z \tau_z \sigma_y, \quad \rho_z \tau_z \sigma_z, \quad \rho_z \tau_y \sigma_0, \quad \rho_z \tau_x \sigma_0, \quad \rho_x \tau_0 \sigma_0, \quad \rho_y \tau_0 \sigma_0, \quad (3)$$

The structure is self-explanatory. In general, one can extend this scheme to construct a set of anticommuting $2^n \times 2^n$ matrices based on the knowledge of $2^{n-1} \times 2^{n-1}$ such matrices. At every step two new matrices are added which adds up to a total of $3 + 2 \times (n - 1) = 2n + 1$ matrices.

Mathematically, this corresponds to the Clifford algebra $Cl_{d-1}(\mathbb{R})$ for even d , plus the chiral matrix. The Clifford algebra always contains d matrices Γ_i with $i = 0, \dots, d - 1$ and for even d it admits the chiral matrix $\Gamma_c = i^{d/2-1} \Gamma_0 \dots \Gamma_{d-1}$ which anticommutes with all Γ_i . For $2^{d/2} \times 2^{d/2} \equiv 2^n \times 2^n$ we have $d = 2n$ and adding Γ_c one arrives at $2n + 1$ again.

PROBLEM 2

$$H = -t \sum_{\langle ij \rangle} \left(e^{i\phi_{ij}} c_i^\dagger c_j + h.c. \right) \quad (4)$$

a) In the gauge $\vec{A} = B(0, x, 0)$, the phase factors along the x-bonds remain 1 but the phase factors along the y-bonds are x-dependent. We denote $\phi = 2\pi/q$, and illustrate the situation in Fig. ??.

The smallest unit cell contains q sites for

$$B = \frac{p}{q} \frac{\Phi_0}{a^2} \quad (5)$$

The band structure thus consists of q bands for each \vec{k} in the 1st BZ, the latter being $k_x \in (-\frac{\pi}{qa}, \frac{\pi}{qa})$, $k_y \in (-\frac{\pi}{a}, \frac{\pi}{a})$.

b) For $p = 1$, $q = 2$ the unit cell has 2 sites (Fig. ??), denoted by $a_{\vec{r}}$, $b_{\vec{r}}$

$$H = -t \sum_{ij} \left[\left(a_{ij}^\dagger b_{ij} + a_{ij}^\dagger b_{i-1j} \right) + \left(a_{ij}^\dagger a_{ij+1} + b_{ij}^\dagger b_{ij+1} \right) + h.c. \right]. \quad (6)$$

Fourier transforming, we get $H = \sum_{\vec{k}} \Psi_{\vec{k}}^\dagger \mathcal{H}(\vec{k}) \Psi_{\vec{k}}$ with $\Psi_{\vec{k}} = (a_{\vec{k}}, b_{\vec{k}})^T$ and

$$\mathcal{H}(\vec{k}) = -2t \begin{pmatrix} \cos k_y & e^{-ik_x} \cos k_x \\ e^{ik_x} \cos k_x & -\cos k_y \end{pmatrix}, \quad (7)$$

$$E_{\vec{k}} = \pm 2t \sqrt{\cos^2 k_x + \cos^2 k_y}. \quad (8)$$

The Dirac points are at $\vec{k} = (\frac{\pi}{2a}, \pm \frac{\pi}{2a})$, as shown in Fig. ??.

Near the Dirac points, the low- E Hamiltonian can be written as

$$\mathcal{H}_{eff}(\vec{q}) = v_F (\tau_z \sigma_z q_y + \sigma_y q_x) \quad (9)$$

where $\tau_z = \pm 1$ labels the two "valleys".

c) Time reversal - this is a spinless Hamiltonian so \mathcal{T} is represented as

$$\mathcal{T}: \quad \Theta = K, \quad \mathcal{H}^*(\vec{k}) = \mathcal{H}(-\vec{k}) \quad (10)$$

Inversion is more complicated because the system is not inversion symmetric about the midpoint between a and b sites. However, either site a or site b can be taken as the center of inversion. This is generated by

$$\mathcal{P}: \quad P \mathcal{H}(\vec{k}) P^{-1} = \mathcal{H}(-\vec{k}), \quad \text{where } P = \begin{pmatrix} e^{-2ik} & 0 \\ 0 & 1 \end{pmatrix} \quad (11)$$

In the low- E theory, valleys get exchanged under both τ and \mathcal{P} :

$$\begin{aligned} \mathcal{T}: \quad \tau_x \mathcal{H}_{eff}^*(\vec{q}) \tau_x &= \mathcal{H}_{eff}(-\vec{q}) & (12) \\ \mathcal{P}: \quad \tau_x \sigma_z \mathcal{H}_{eff}(\vec{q}) \sigma_z \tau_x &= \mathcal{H}_{eff}(-\vec{q}) & (13) \end{aligned}$$

\mathcal{T} -invariance of H with $q = 2$ is special. It holds because the smallest loop that an electron can traverse (one square plaquette) contains a flux $(1/2)\Phi_0$ and thus gives rise to a phase factor $e^{i\pi} = -1$. This is real and the electron cannot differentiate between clockwise and counter-clockwise hopping.

d) Possible mass terms are: $\sigma_x \tau_z$, σ_x , $\sigma_z \tau_x$, and $\sigma_z \tau_y$ (these are the only terms that anticommute with $\mathcal{H}_{eff}(\vec{q})$).

Mass term	\mathcal{P}	\mathcal{T}
$\sigma_x \tau_z$	✓	✗
$\sigma_z \tau_x$	✓	✓
$\sigma_z \tau_y$	✗	✓
σ_x	✗	✓

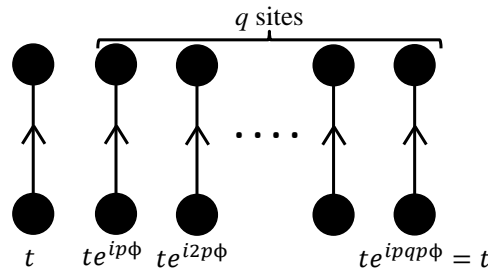
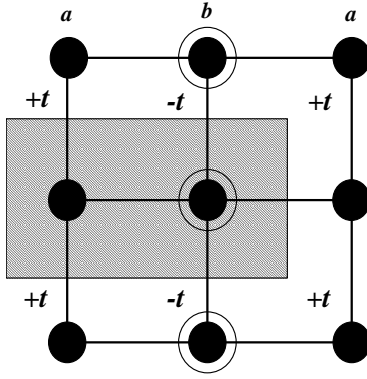


FIG. 1. Hopping parameters along y direction.

FIG. 2. Unit cell for $p = 1, q = 2$

The \mathcal{T} -breaking mass $\sigma_x \tau_z$ describes second-neighbor hopping with imaginary amplitude. The \mathcal{P} -breaking mass σ_x can be realized by dimerizing the lattice along the x-direction, i.e. alternating $t + \delta t$ and $t - \delta t$.

e) With the \mathcal{T} -breaking mass we have

$$\mathcal{H}_{eff}(\vec{q}) = v_F (\tau_z \sigma_z q_y + \sigma_y q_x) + m \sigma_x \tau_z \quad (14)$$

We can calculate the Chern number for the two valleys in analogy with graphene

$$C_1 = \frac{1}{2} \text{sgn} [(+v_F)(+v_F)(+m)] = \frac{1}{2} \text{sgn}(m) \quad (\tau_z = +1) \quad (15)$$

$$C_2 = \frac{1}{2} \text{sgn} [(-v_F)(+v_F)(-m)] = \frac{1}{2} \text{sgn}(m) \quad (\tau_z = -1) \quad (16)$$

Thus, $C = C_1 + C_2 = \text{sgn}(m) = \pm 1$. The system is a Chern insulator.

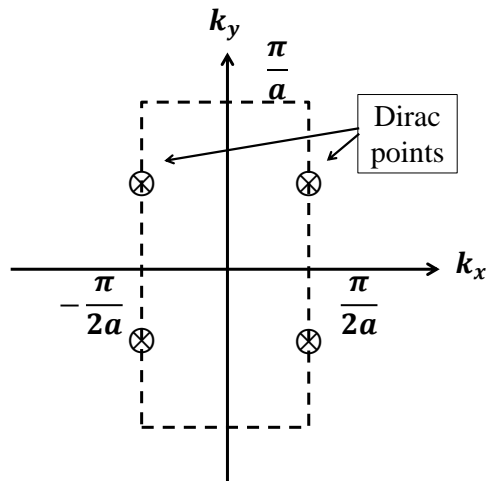


FIG. 3. Brillouin Zone and Dirac points.