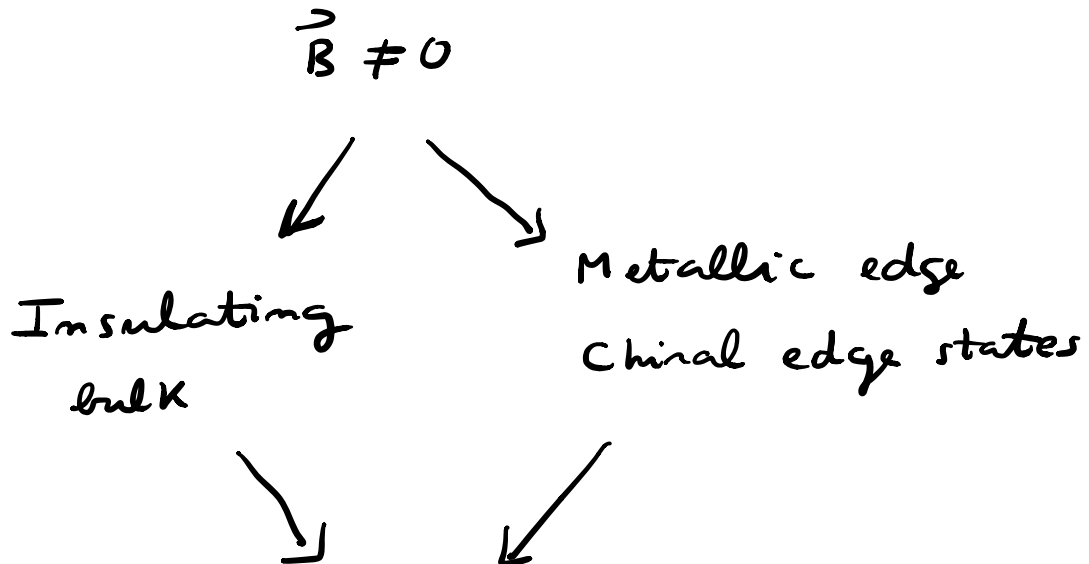


② Quantum Hall insulator



$$\sigma_{xy} = \nu \frac{e^2}{h}$$

$$\nu = \sum_{m \in \text{occ}} c_m$$

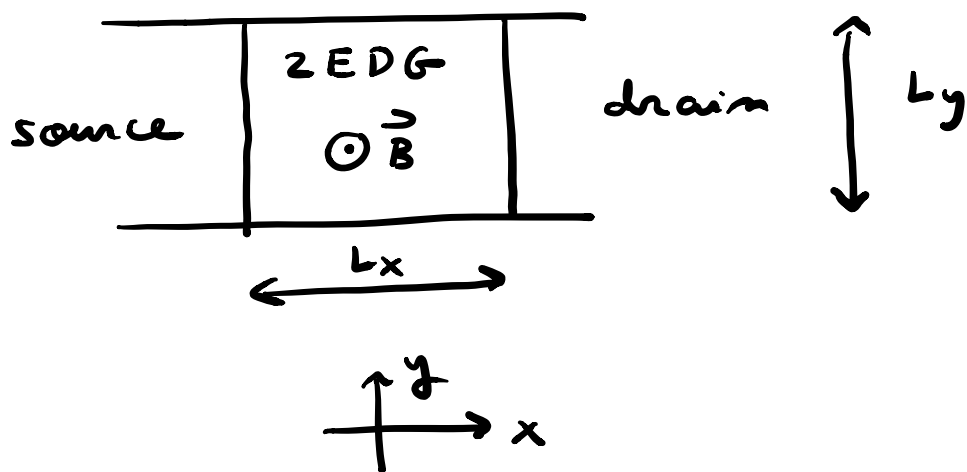
= # of occupied LLs in bulk

= # of chiral edge states

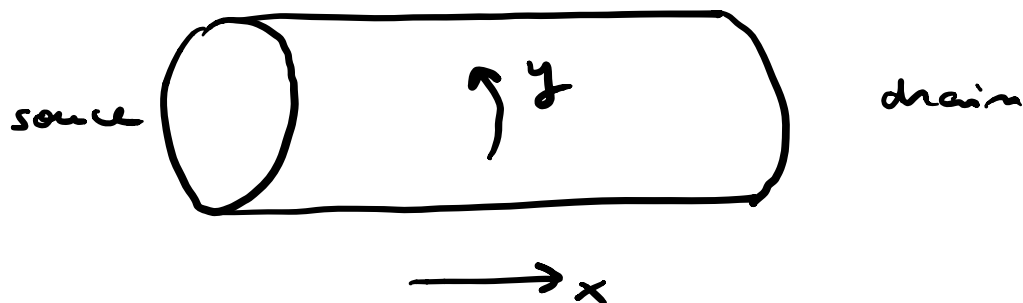
"Bulk-edge correspondence"

2.4 Laughlin's argument for Hall quantitation

[Laughlin, PRB 23, 5632 (1981)]



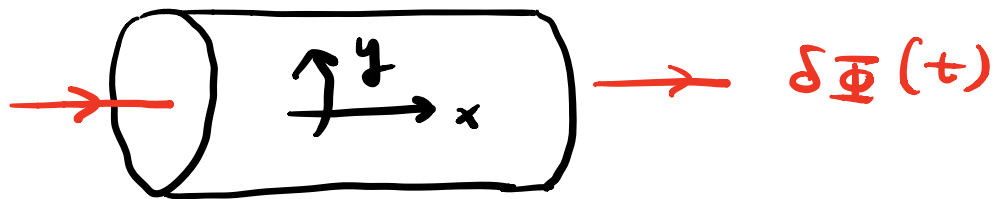
Assume periodic boundary conditions along y .



$\vec{B}$ is radial w.r.t. cylinder axis.

2 DEG: surface of cylinder.

Insert a magnetic flux
(adiabatically) through hole
of cylinder.



vector potential generated by
flux:

$$\vec{\delta A}(t) = \frac{\delta\Phi(t)}{L_y} \hat{y}$$

check:

$$\delta \Phi(t) = \oint \delta \vec{A} \cdot d\vec{\ell}$$


loop around
cylinder

$$= \oint dy \frac{\delta \Phi}{L_y} = \delta \Phi \quad \checkmark$$

The flux insertion produces
an electric field:

$$\delta \vec{E}(t) = - \frac{\partial \delta \vec{A}}{\partial t}$$
$$= - \hat{y} \frac{1}{L_y} \frac{\partial}{\partial t} \delta \Phi \parallel \hat{y}$$

Source-drain charge current density induced by $\delta \vec{E}$:

$$\delta j_x(t) = \sigma_{xy} \delta E(t)$$

$\uparrow$
A/m

$\uparrow$
intrinsic
(independent of δE)

Charge transferred from source to drain, between $t=0$ and $t=T$:

$$\delta Q = \oint dy \int_0^T dt \delta j_x$$

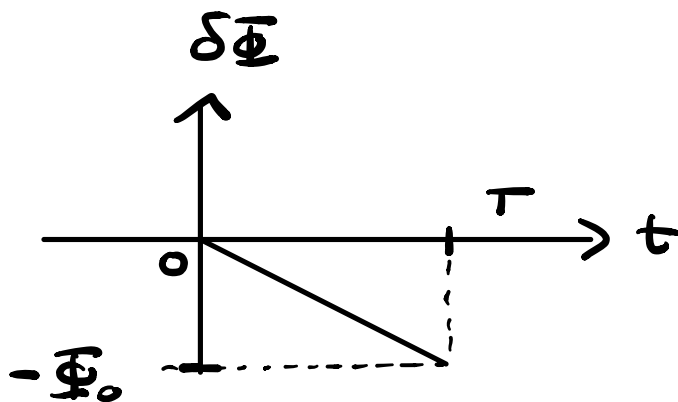
$\uparrow$
C

$$= -\sigma_{xy} \int_0^{\tau} \frac{\partial}{\partial t} \delta\Phi(t)$$

$$= -\sigma_{xy} [\delta\Phi(\tau) - \delta\Phi(0)]$$

↑
independent
of $\delta\Phi$

Let's choose to insert a
single quantum of flux



$$\Phi_0 = h/e$$

$$\Rightarrow \delta Q = \sigma_{xy} \Phi_0$$

Gauge invariance : when
inserting a flux quantum,
one must have

$$\delta Q = n e, \text{ where}$$

$$n \in \mathbb{Z}.$$

$$\Rightarrow n e = \sigma_{xy} \frac{h}{e}$$

$$\Rightarrow \sigma_{xy} = n \frac{e^2}{h}$$

* "Proof" # 1 for $\delta a = ne$

Total vector potential:

$$\vec{A} = \underbrace{B \times \hat{y}}_{\substack{\uparrow \\ \text{due to} \\ \text{radial} \\ \text{field}}} + \underbrace{\delta \vec{A}}_{\substack{\uparrow \\ \text{due to} \\ \text{flux insertion} \\ \text{along cylinder}}}$$

$$= \left(B \times + \frac{\delta \Phi}{L_y} \right) \hat{y}$$

$$\mathcal{H} = \frac{1}{2m} \left(\vec{p} - e \vec{A} \right)^2$$

$$\Psi(x, y) = \frac{e^{ik_y y}}{\sqrt{L_y}} \phi_k(x)$$

$$\hbar(k) \phi_k(x) = E_k \phi_k(x)$$

where

$$\hbar(k) = \frac{1}{2m} \left[p_x^2 + (\hbar k - eA)^2 \right]$$

$$= \frac{p_x^2}{2m} + \frac{1}{2} m \omega_c^2 (x - x_k)^2$$

where

x_k = guiding center

$$= \frac{\hbar k}{eB} + \frac{\delta \Phi}{Ly}$$

Now recall

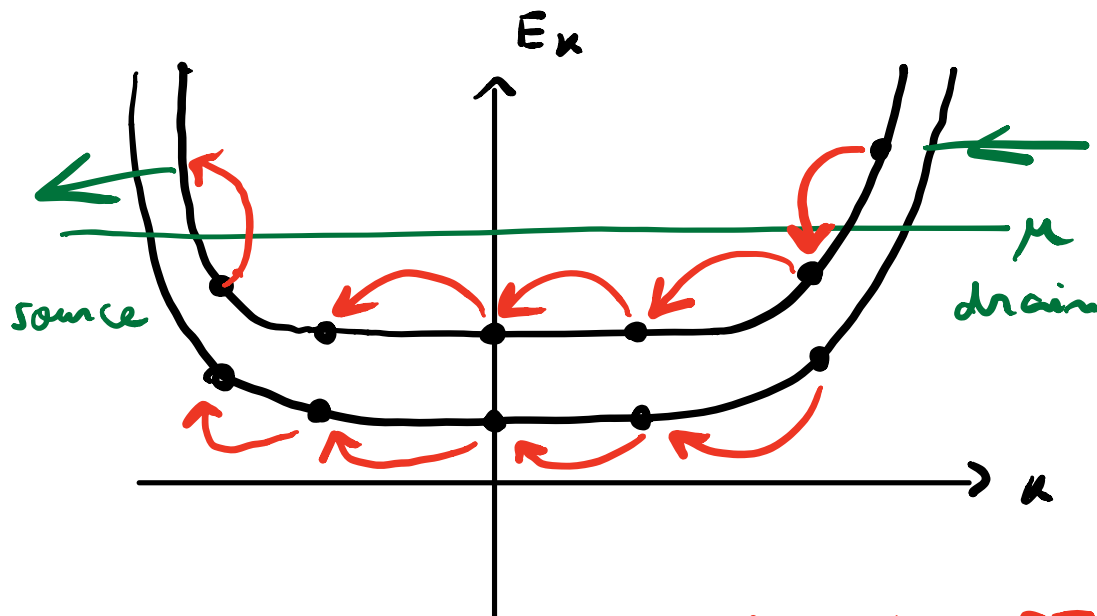
$$k = \frac{2\pi}{L_y} l, \quad l \in \mathbb{Z}$$

$$\Rightarrow \underbrace{x_l}_{\uparrow} = \frac{h}{e B L_y} \left(l - \frac{\delta \Phi}{\Phi_0} \right)$$

center of
wave function.

Having $\delta \Phi = \Phi_0$ is

equivalent to $l \rightarrow l - 1$



$\curvearrowright = \text{flow of states when } \delta \Phi = \Phi_0$

For every value of x ,

$$\# \text{ of } e^-s = \# \text{ of} \\ \text{occupied Ls} = \nu$$

Transferred $\#$ of e^-s upon
insertion of Φ_0 : ν

This proves $\delta Q = e \nu$

$$\Rightarrow \sigma_{xy} = \nu \frac{e^2}{h}$$

* Proof # 2 for $\delta Q = ne$:

System must be periodic
in Φ , w/ periodicity Φ_0
(AB effect).

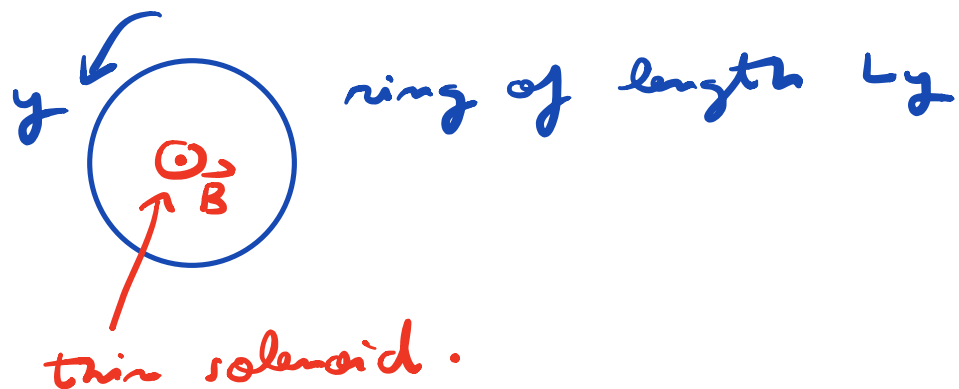
In other words, inserting Φ_0 gradually is like completing a full turn to Archimedes' screw.

Therefore: charge pumped in a cycle is equal to integer $\times e$.

$$\text{integer} = \sum_{n \in \mathbb{Z}} C_n.$$

In a QH insulator, this is precisely ν .

* Why the system is periodic
in Φ with periodicity
 Φ_0 ?



Take $\vec{A} = \underbrace{A}_{\substack{\uparrow \\ \text{independent of } y}} \hat{y}$

$$\mathcal{H} = \frac{1}{2m} (p_y - eA)^2 + v(y)$$

$$V(y) = V(y + Ly)$$

$$\mathcal{H}\psi = E\psi$$

$$\text{If } A=0, \quad \psi(y) = \psi(y + Ly)$$

$$\text{Let} \quad i \frac{2\pi}{\Phi_0} \int_0^y A dy'$$

$$\tilde{\psi}(y) \equiv \psi(y) e$$

$$= \psi(y) e^{i \frac{2\pi}{\Phi_0} A y}$$

Replace this in Schrödinger equation:

$$\left[\frac{\hbar^2}{2m} \frac{d^2}{dy^2} + V(y) \right] \tilde{\psi} = E \tilde{\psi}$$

A disappeared? No.

It went to the boundary condition.

$$\tilde{\psi}(y+L_y) = e^{i \frac{2\pi}{\Phi_0} A L_y}$$

$$\underbrace{e^{i \frac{2\pi}{\Phi_0} A y} \underbrace{\psi(y+L_y)}_{\psi(y)}}_{\tilde{\psi}(y)}$$

$$\tilde{\psi}(y+L_y) = \tilde{\psi}(y) e^{i \frac{2\pi}{\Phi_0} A L_y}$$

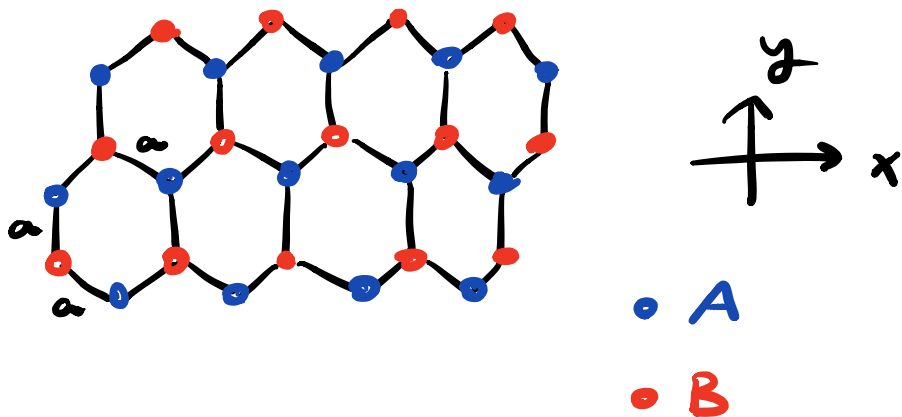
$$= \tilde{\psi}(y) e^{i \frac{2\pi \Phi}{\Phi_0}}$$

Φ alters the solutions of the Schrödinger eq. unless

$\Phi = n \Phi_0$. In the latter case, the flux completely disappears from the problem.

③ Graphene

* Hexagonal lattice of C atoms



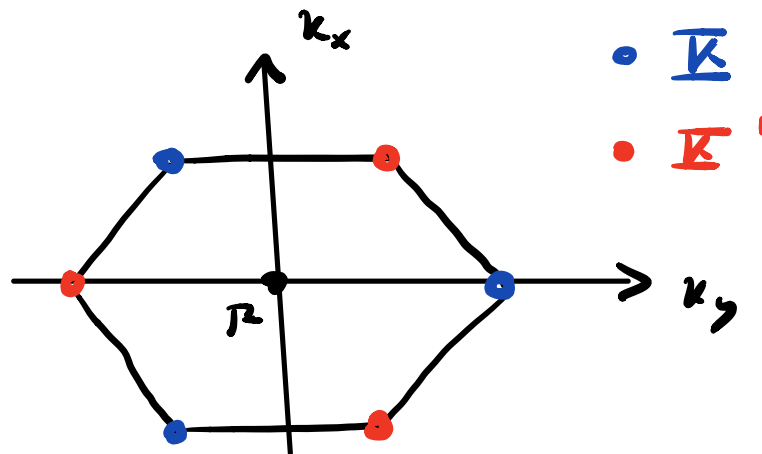
Two atoms per unit cell.

In graphene, A and B are carbon atoms.

BN : A = Boron

B = Nitrogen.

* 1 BZ



3.1 Low-energy electron states

* C : $1s^2 2s^2 2p^2$



π orbital is partially filled.

$\Rightarrow$ low-energy physics is associated to π orbital.

* Minimal tight-binding model (one orbital per site):

$$\mathcal{H}_0 = -t \sum_{\langle ij \rangle} (c_i^\dagger c_j + \text{h.c.})$$

$\langle \rangle$: nearest neighbor.

For now, neglect the spin
of e^- -s.

Fourier transform:

$$H_0 = \sum_{\vec{k}} (c_{\vec{k}A}^{\dagger}, c_{\vec{k}B}^{\dagger}) \\ h_0(\vec{k}) \begin{pmatrix} c_{\vec{k}A} \\ c_{\vec{k}B} \end{pmatrix}$$

where

$$h_0(\vec{k}) = \vec{d}(\vec{k}) \cdot \vec{\sigma}$$

$\vec{\sigma}$: sublattice pseudospin

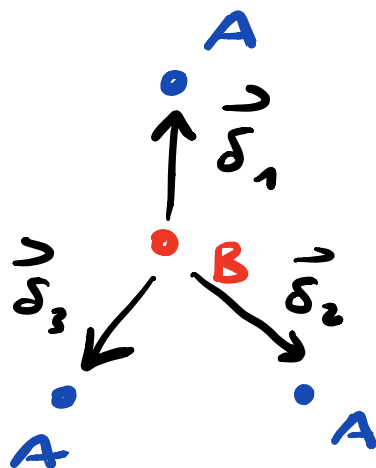
$\sigma_z \uparrow$: A

$\sigma_z \downarrow$: B

$$d_x(\vec{k}) = -t \sum_{j=1}^3 \cos(\vec{k} \cdot \vec{\delta}_j)$$

$$d_y(\vec{k}) = -t \sum_{j=1}^3 \sin(\vec{k} \cdot \vec{\delta}_j)$$

$$d_z(\vec{k}) = 0.$$



Nearest
-neighbor
vectors.

$$\vec{\delta}_1 = a(0, 1)$$

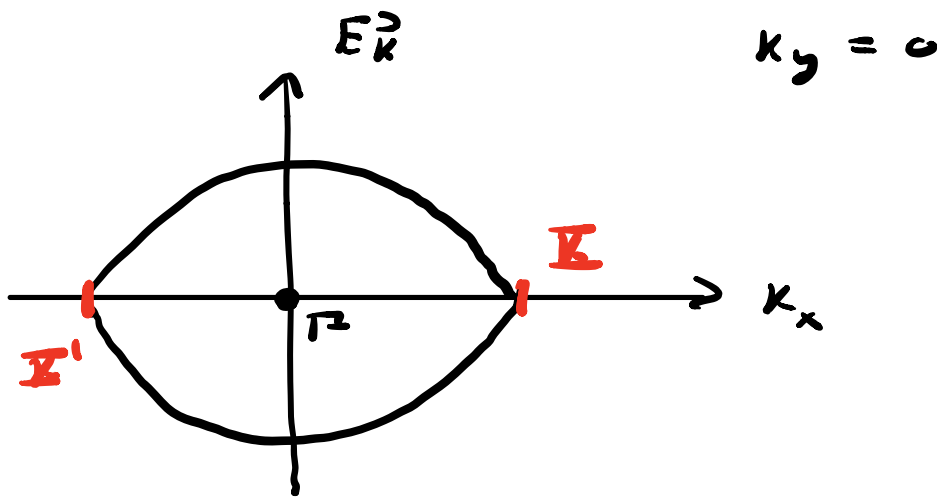
$$\vec{\delta}_2 = a\left(\frac{\sqrt{3}}{2}, -\frac{1}{2}\right)$$

$$\vec{\delta}_3 = a\left(-\frac{\sqrt{3}}{2}, -\frac{1}{2}\right)$$

Energy eigenvalues:

$$E_{\vec{k} \pm} = \pm |\vec{d}(\vec{k})|$$

$$= \pm \sqrt{d_x(\vec{k})^2 + d_y(\vec{k})^2}$$



Bands touch at Γ and Γ'

$\Rightarrow$ at half-filling (one e^- per unit cell), graphene is a semimetal.

More precisely, it's a Dirac semimetal. Why Dirac?

Expand $\vec{k}$ near $\vec{K}$ and $\vec{K}'$:

$$(i) \quad \vec{k} = \vec{K} + \vec{q},$$

where $qa \ll 1$.

Then,

$$d_x \approx -\frac{3}{2}atq_x \equiv vq_x$$

$$d_y \approx -\frac{3}{2}atq_y \equiv vq_y$$

$$\Rightarrow \left[\hbar_0^{(\vec{K})} \left(\vec{q} \right) = v \left(\sigma^x q_x + \sigma^y q_y \right) \right] \quad (1)$$



momentum

wrt to $\vec{K}$

Massless 2D Dirac fermion,

with Dirac velocity v .

$$(ii) \quad \vec{k} = \vec{k}' + \vec{q}$$

$$qa \ll 1 \quad (2)$$

$$h_0^{(\vec{k}')}(\vec{q}) \approx v (-\sigma^x q_x + \sigma^y q_y)$$

Massless 2D Dirac fermion.

Combine (1) and (2) into
a 4×4 Hamiltonian:

$$\overline{h_{eff}(\vec{q})} = \begin{pmatrix} \overline{h_0^{(\vec{k})}} & 0 \\ 0 & \overline{h_0^{(\vec{k}')}} \end{pmatrix}$$

$$= v (\tau^z \sigma^x q_x + \sigma^y q_y)$$

where $\vec{\tau}$ is the valley pseudospin.

$$\tau^z \uparrow : \underline{K}$$

$$\tau^z \downarrow : \underline{K}'$$

$h_{\text{eff}}(\vec{\tau})$ is the low-energy effective theory of graphene.

Dirac fermions may be

gapped (they can acquire a nonzero mass) by breaking either space inversion symmetry or time-reversal symmetry.

Some gapped phases are topologically nontrivial.