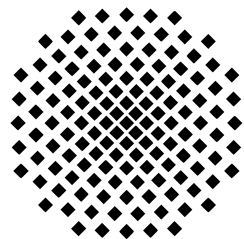
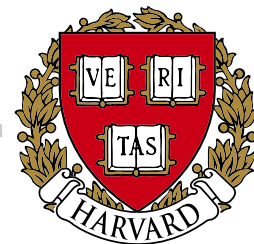


Topological flat bands with Chern number $C = 2$
by dipolar exchange interactions

David Peter — Palaiseau — November 6th, 2014



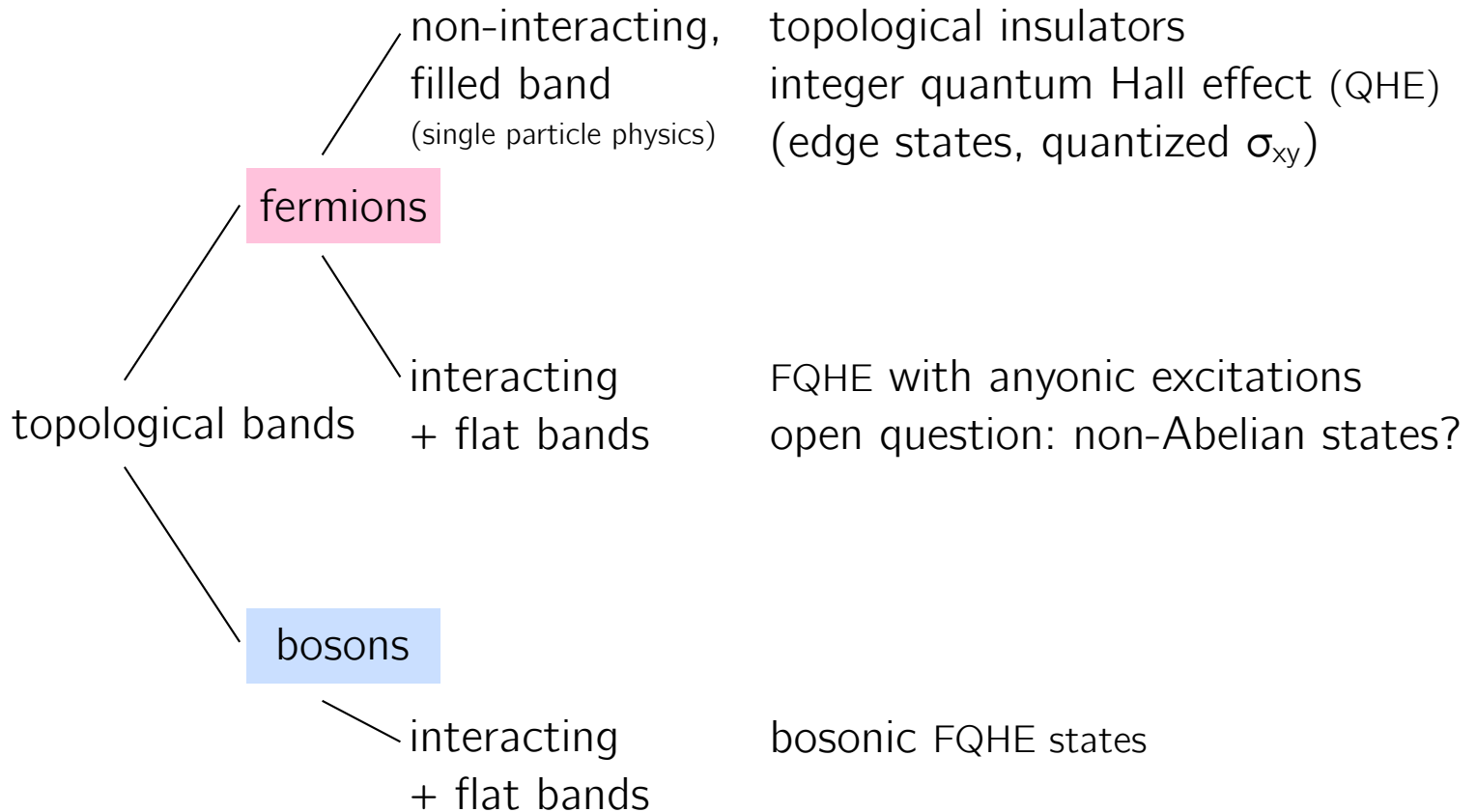
Hans Peter Büchler
Nicolai Lang

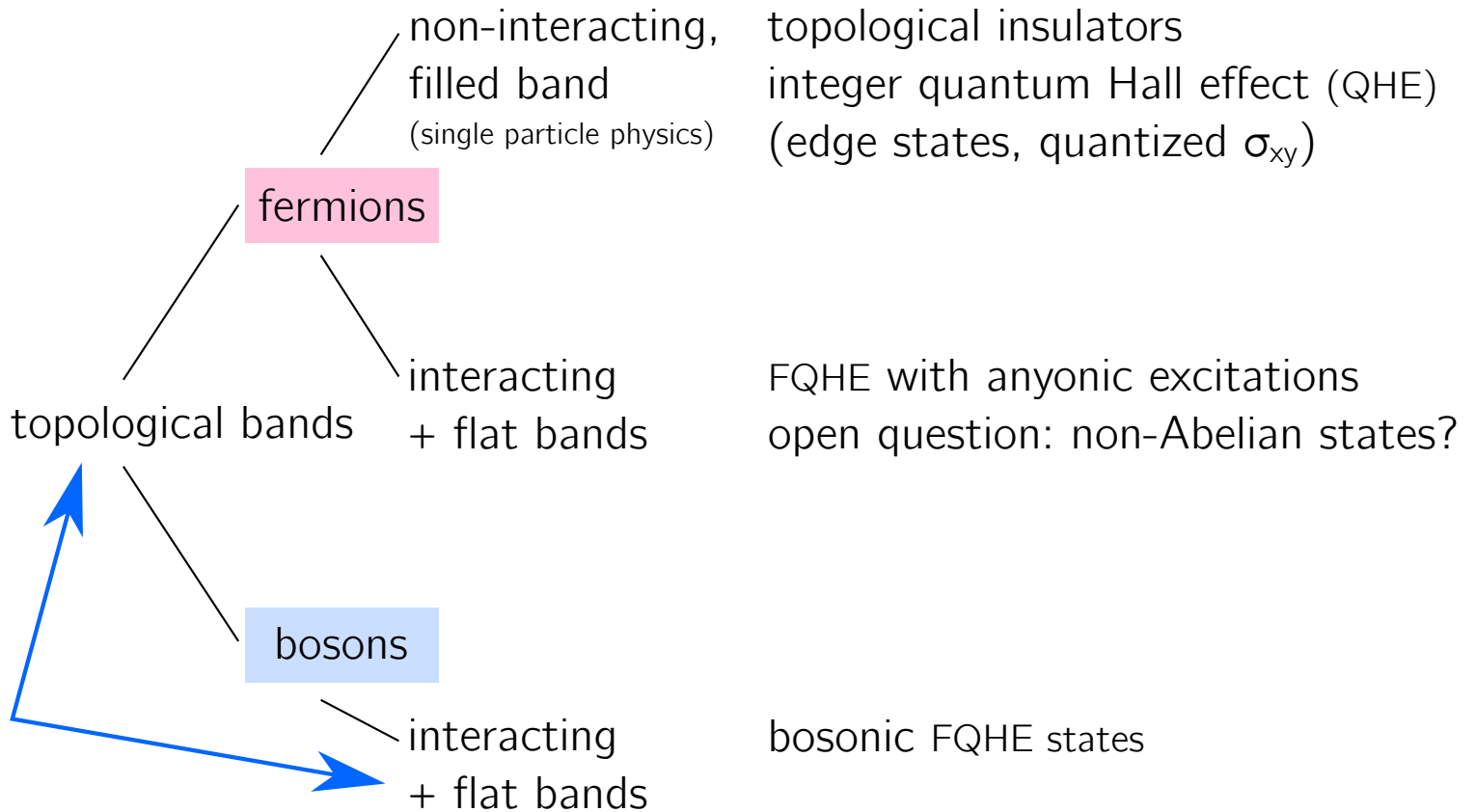


Mikhail Lukin
Norman Yao



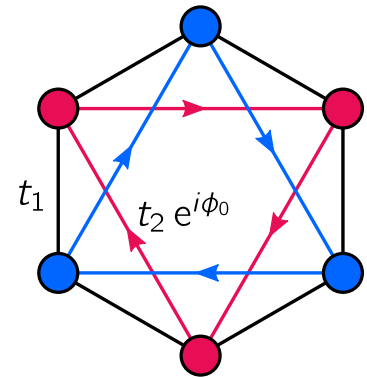
Sebastian Huber



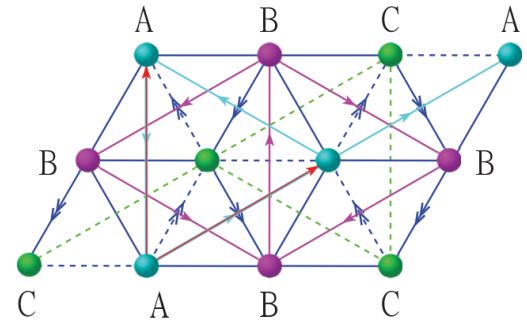


topological bands

tunneling phases for neutral particles?



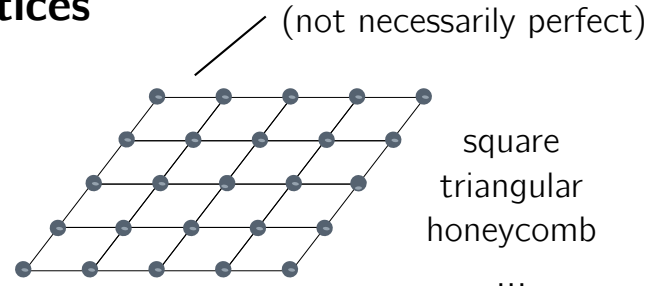
F. D. M. Haldane, Phys. Rev. Lett. **61**, 2015 (1988)



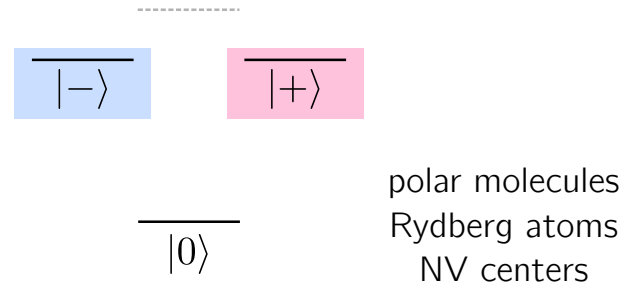
Y. Wang. Phys. Rev. B **86**, 201101 (2012)

"natural" realization with dipoles on lattices

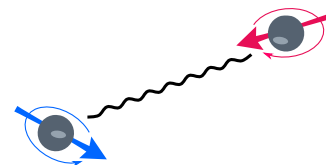
▷ Two-dimensional lattice

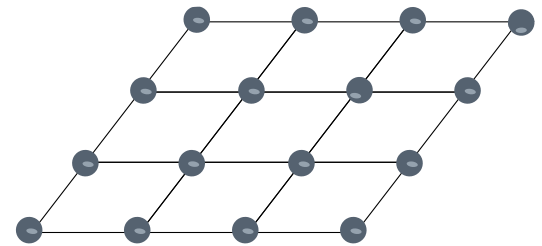


▷ Three-level dipole



▷ Strong dipole-dipole interactions





one molecule
pinned at each lattice site

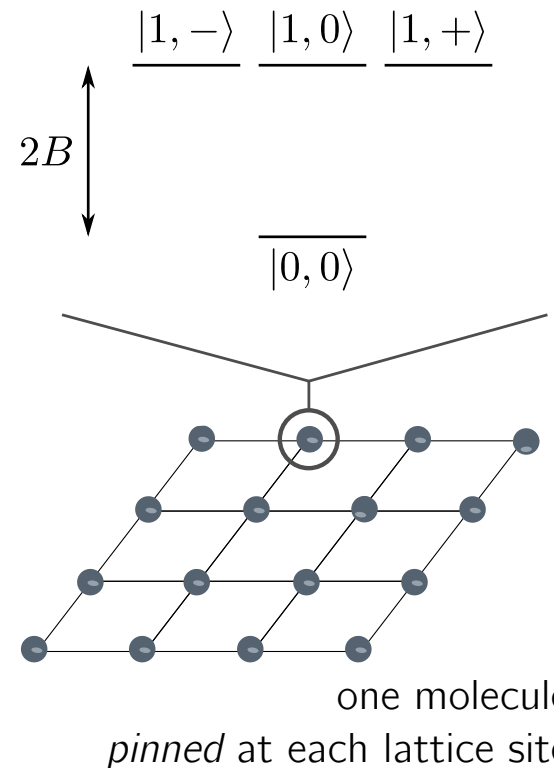
Three level dipole from rotational structure:

$$H_i^{\text{rot}} = B J_i^2 - \mathbf{d}_i \cdot \mathbf{E}$$

rotational constant angular momentum dipole moment external fields

eigenstates without electric field:

$$|J, m\rangle$$



see also:

N. Yao, Phys. Rev. Lett. **109**, 266804 (2012)

N. Yao, Phys. Rev. Lett. **110**, 185302 (2013)

S. Syzranov, arXiv:1406.0570

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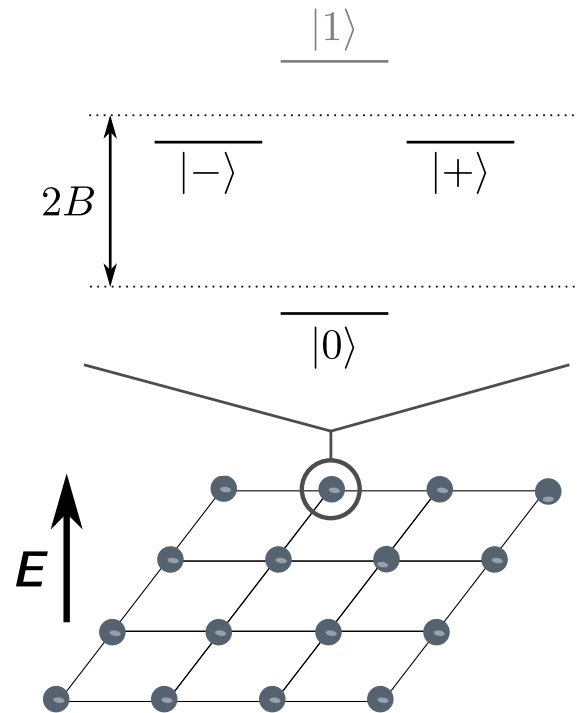
rotational constant angular momentum dipole moment external fields

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eigenstates with electric field:

m is still a good quantum number



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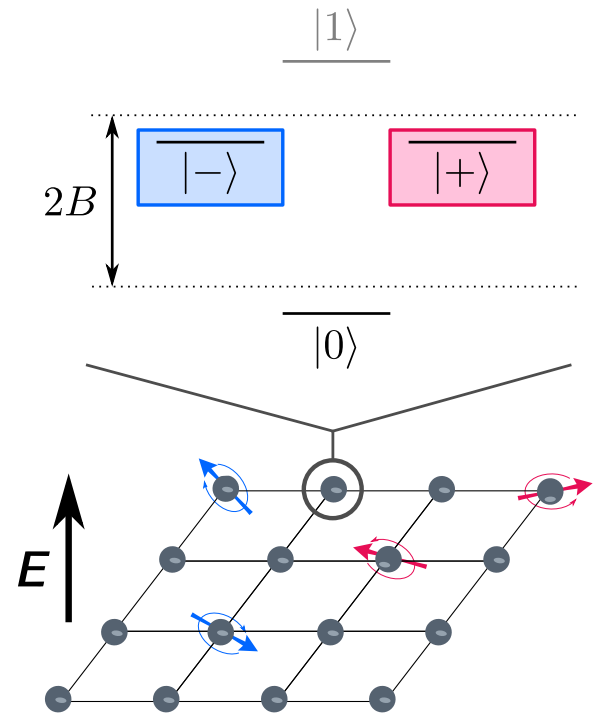
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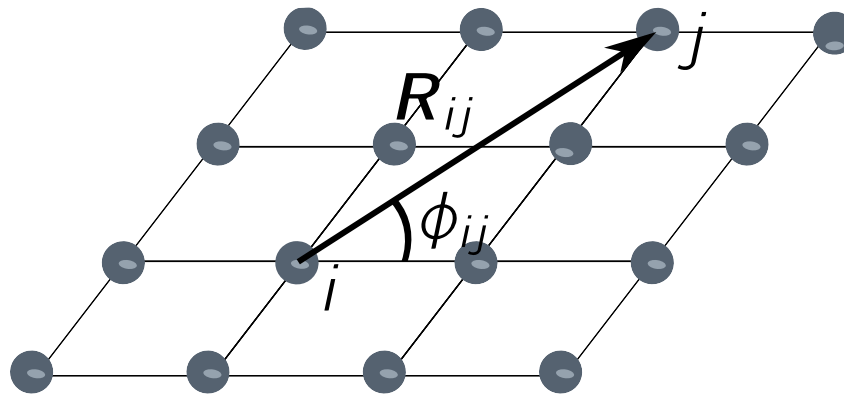
N. Yao, Phys. Rev. Lett. **109**, 266804 (2012)

N. Yao, Phys. Rev. Lett. **110**, 185302 (2013)

S. Syzranov, arXiv:1406.0570

dipole-dipole interactions in the plane

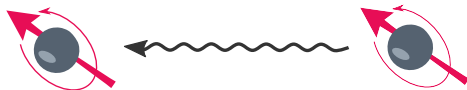
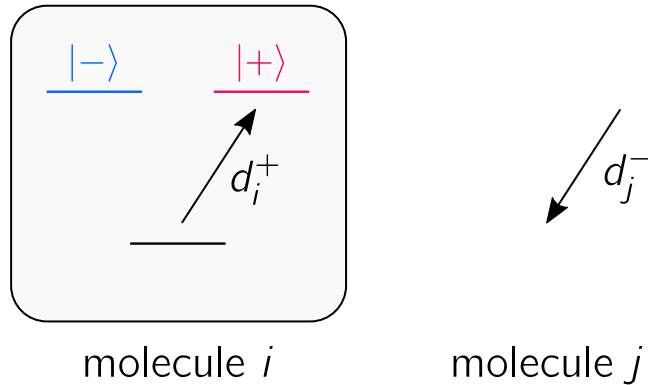
$$H_{ij}^{\text{dd}} = \frac{1}{|\mathbf{R}_{ij}|^3} \left[d_i^0 d_j^0 + \frac{1}{2} (d_i^- d_j^+ + d_i^+ d_j^-) - \frac{3}{2} (d_i^- d_j^- e^{2i\phi_{ij}} + d_i^+ d_j^+ e^{-2i\phi_{ij}}) \right]$$



dipole-dipole interactions in the plane

$$H_{ij}^{\text{dd}} = \frac{1}{|\mathbf{R}_{ij}|^3} \left[d_i^0 d_j^0 + \frac{1}{2} (d_i^- d_j^+ + d_i^+ d_j^-) - \frac{3}{2} (d_i^- d_j^- e^{2i\phi_{ij}} + d_i^+ d_j^+ e^{-2i\phi_{ij}}) \right]$$

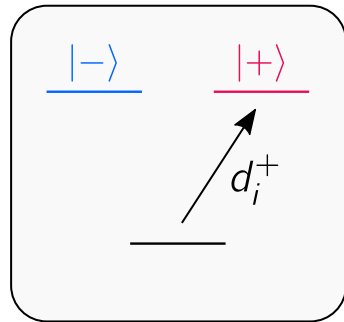
energy conserving processes:



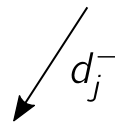
dipole-dipole interactions in the plane

$$H_{ij}^{\text{dd}} = \frac{1}{|\mathbf{R}_{ij}|^3} \left[d_i^0 d_j^0 + \frac{1}{2} (d_i^- d_j^+ + d_i^+ d_j^-) - \frac{3}{2} (d_i^- d_j^- e^{2i\phi_{ij}} + d_i^+ d_j^+ e^{-2i\phi_{ij}}) \right]$$

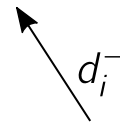
energy conserving processes:



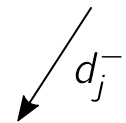
molecule *i*



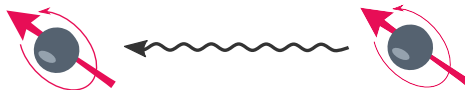
molecule *j*

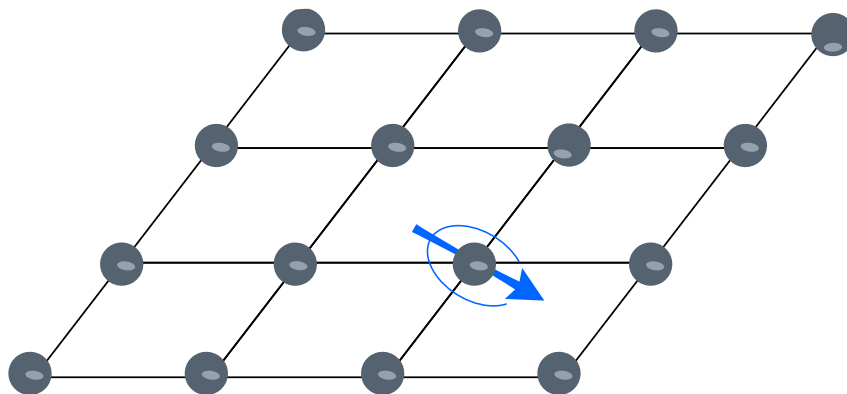


molecule *i*



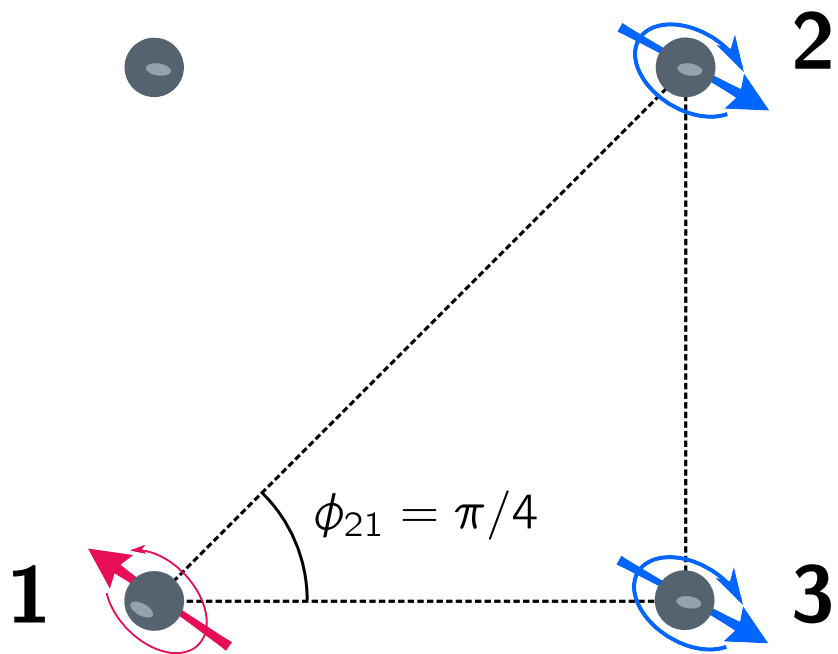
molecule *j*

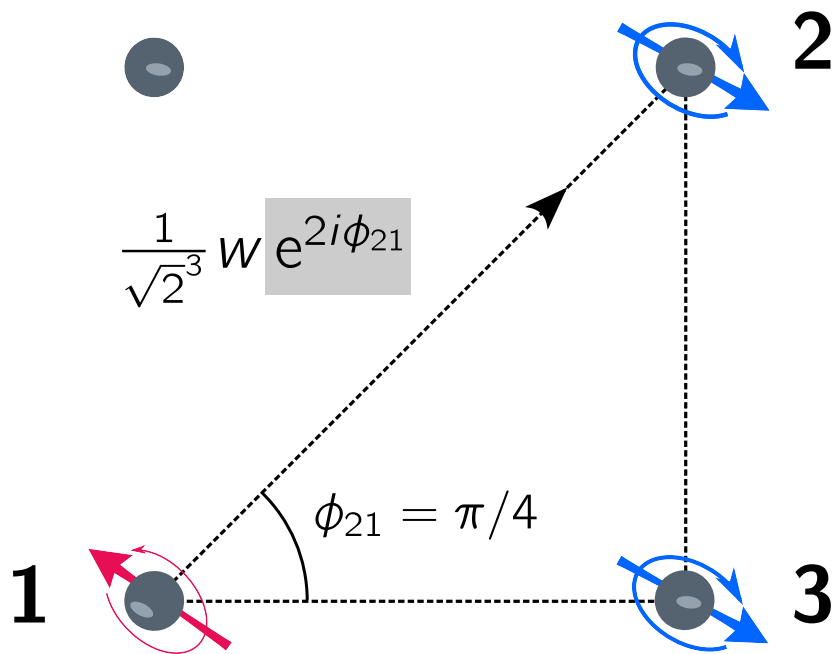


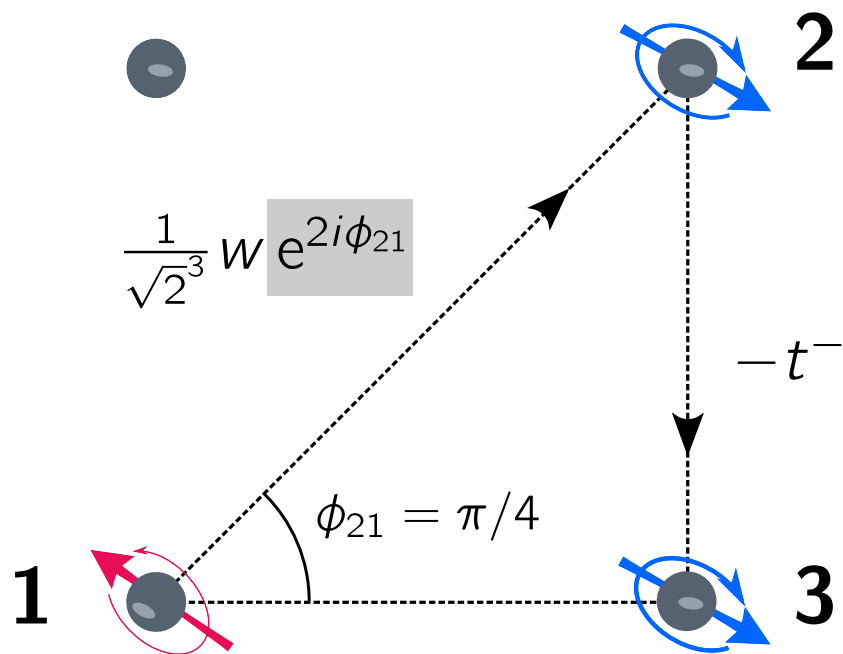


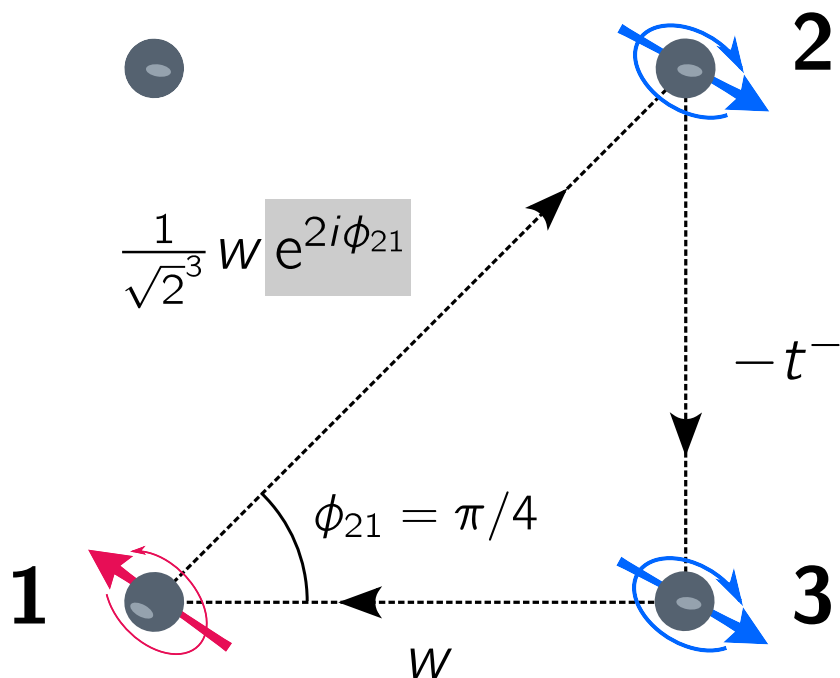
$$H_{ij}^{\text{dd}} = \frac{1}{|\mathbf{R}_{ij}|^3} \left[d_i^0 d_j^0 + \frac{1}{2} (d_i^- d_j^+ + d_i^+ d_j^-) - \frac{3}{2} (d_i^- d_j^- e^{2i\phi_{ij}} + d_i^+ d_j^+ e^{-2i\phi_{ij}}) \right]$$

interaction term, not relevant
for single excitation dynamics









$$\varphi_{\text{flux}} = \arg \left(\prod_{\triangle} t_{ij} \right) = -\pi/2$$

Describe excitations as effective particles (hardcore bosons)

$$b_{i,+}^\dagger = |+\rangle_i \langle 0|_i$$

$$b_{i,-}^\dagger = |-\rangle_i \langle 0|_i$$

$$\psi_i^\dagger = \begin{pmatrix} b_{i,+}^\dagger \\ b_{i,-}^\dagger \end{pmatrix}$$

Tunneling Hamiltonian

$$H = \sum_{i \neq j} \frac{a^3}{|\mathbf{R}_{ij}|^3} \psi_i^\dagger \begin{pmatrix} -t^+ & w e^{-2i\phi_{ij}} \\ w e^{2i\phi_{ij}} & -t^- \end{pmatrix} \psi_j$$

long-range tunneling

spin-orbit coupling

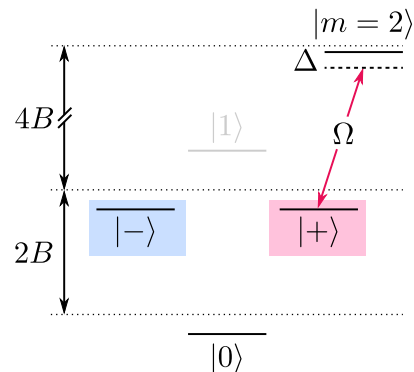
\mathcal{T} -symmetry: $t^+ = t^-$

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Tunneling Hamiltonian

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$$+ \sum_i \psi_i^\dagger \begin{pmatrix} \mu & 0 \\ 0 & -\mu \end{pmatrix} \psi_i$$

~~\mathcal{T} -symmetry: $t^+ = t^-$~~

$$H = \sum_k \psi_k^\dagger \begin{pmatrix} -t^+ \epsilon_k^0 + \mu & W \epsilon_k^{-2} \\ W \epsilon_k^2 & -t^- \epsilon_k^0 - \mu \end{pmatrix} \psi_k$$

$$\psi_k = \frac{1}{\sqrt{N_s}} \sum_j \psi_j e^{ikR_j}$$
$$\epsilon_k^m = \sum_{j \neq 0} \frac{a^3}{|R_j|^3} e^{ikR_j + im\phi_j}$$

$$H = \sum_k \psi_k^\dagger \begin{pmatrix} -t^+ \epsilon_k^0 + \mu & w \epsilon_k^{-2} \\ w \epsilon_k^2 & -t^- \epsilon_k^0 - \mu \end{pmatrix} \psi_k$$

$$= \sum_k \psi_k^\dagger (n_k^0 \mathbb{1} + \mathbf{n}_k \cdot \boldsymbol{\sigma}) \psi_k$$

$$n_k^0 = -\bar{t} \epsilon_k^0$$

spin-orbit coupling with: $\mathbf{n}_k = \begin{pmatrix} w \operatorname{Re} \epsilon_k^2 \\ w \operatorname{Im} \epsilon_k^2 \\ \mu + t \epsilon_k^0 \end{pmatrix}$

$$\bar{t} = (t^- + t^+) / 2$$

$$t = (t^- - t^+) / 2$$

$$\psi_k = \frac{1}{\sqrt{N_s}} \sum_j \psi_j e^{ikR_j}$$

$$\epsilon_k^m = \sum_{j \neq 0} \frac{a^3}{|R_j|^3} e^{ikR_j + im\phi_j}$$

$$H = \sum_{\mathbf{k}} \psi_{\mathbf{k}}^{\dagger} \begin{pmatrix} -t^{+} \epsilon_{\mathbf{k}}^0 + \mu & w \epsilon_{\mathbf{k}}^{-2} \\ w \epsilon_{\mathbf{k}}^2 & -t^{-} \epsilon_{\mathbf{k}}^0 - \mu \end{pmatrix} \psi_{\mathbf{k}}$$

$$= \sum_{\mathbf{k}} \psi_{\mathbf{k}}^{\dagger} (n_{\mathbf{k}}^0 \mathbb{1} + \mathbf{n}_{\mathbf{k}} \cdot \boldsymbol{\sigma}) \psi_{\mathbf{k}}$$

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spin-orbit coupling with: $\mathbf{n}_{\mathbf{k}} = \begin{pmatrix} w \operatorname{Re} \epsilon_{\mathbf{k}}^2 \\ w \operatorname{Im} \epsilon_{\mathbf{k}}^2 \\ \mu + t \epsilon_{\mathbf{k}}^0 \end{pmatrix}$

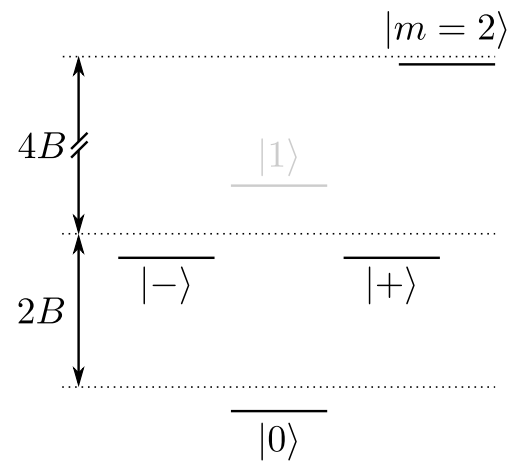
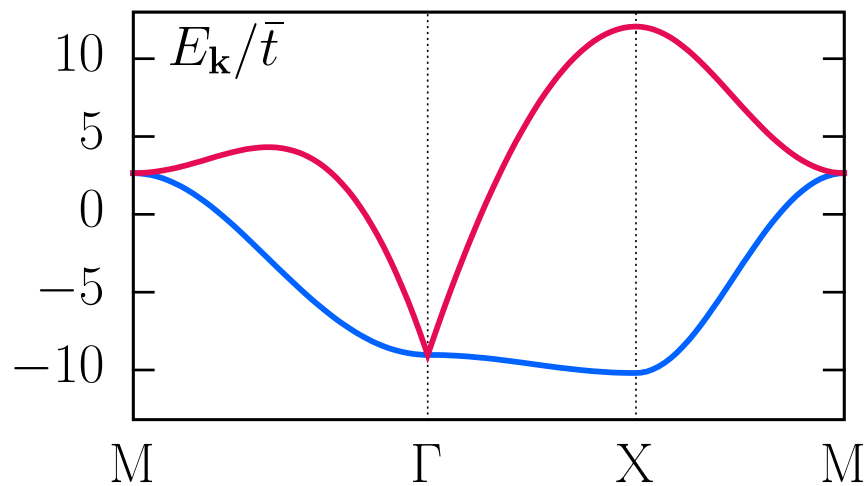
$$\bar{t} = (t^{-} + t^{+}) / 2$$

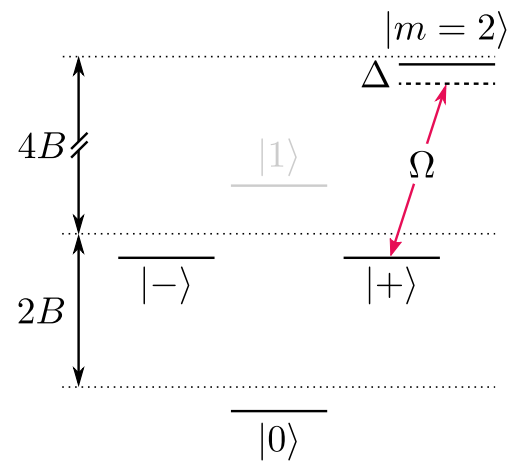
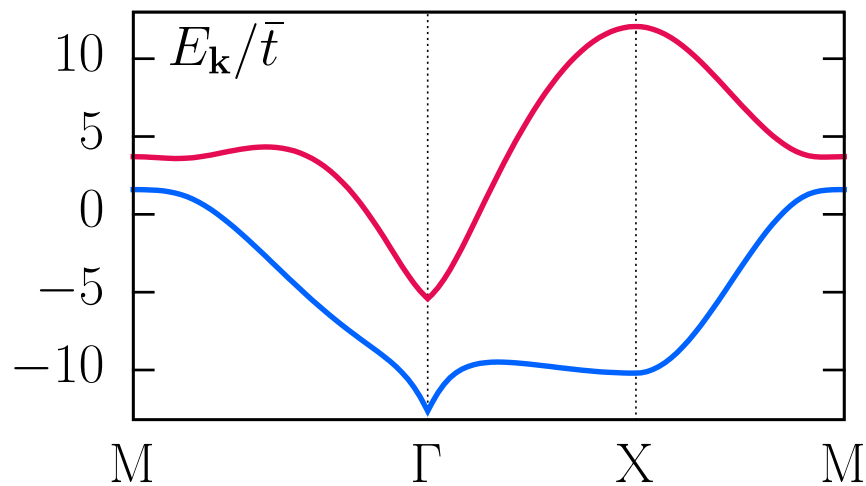
$$t = (t^{-} - t^{+}) / 2$$

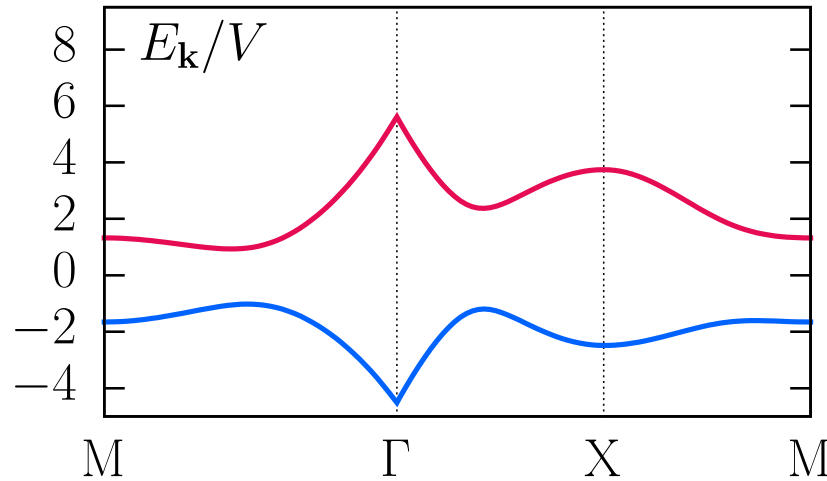
$$E_{\pm}(\mathbf{k}) = n_{\mathbf{k}}^0 \pm |\mathbf{n}_{\mathbf{k}}| = -\bar{t} \epsilon_{\mathbf{k}}^0 \pm \sqrt{w^2 |\epsilon_{\mathbf{k}}^2|^2 + (\mu + t \epsilon_{\mathbf{k}}^0)^2}$$

$$\psi_{\mathbf{k}} = \frac{1}{\sqrt{N_s}} \sum_j \psi_j e^{i\mathbf{k} \cdot \mathbf{R}_j}$$

$$\epsilon_{\mathbf{k}}^m = \sum_{j \neq 0} \frac{a^3}{|\mathbf{R}_j|^3} e^{i\mathbf{k} \cdot \mathbf{R}_j + i m \phi_j}$$







← after tuning parameters

are these bands topological?

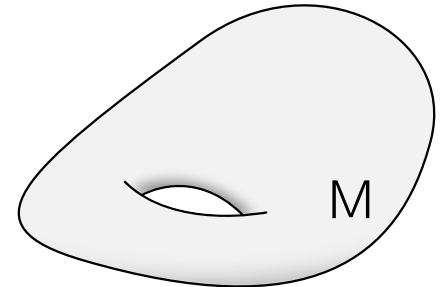
Gauss-Bonnet-Theorem

For a two-dimensional closed manifold M :

$$\int_M dS K = 4\pi(1 - g)$$

locally defined quantity
(Gaussian curvature)

topological index
(genus $g \in \mathbb{N}_0$)



Review on topological insulators:

M. Hasan and C. Kane, Rev. Mod. Phys. **82**, 3045 (2010)

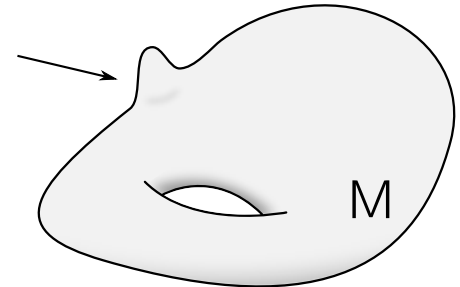
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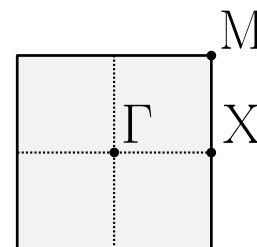
M. Hasan and C. Kane, Rev. Mod. Phys. **82**, 3045 (2010)

Chern number

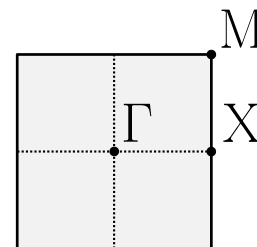
$$\frac{1}{2\pi} \int_{\text{BZ}} d^2\mathbf{k} \mathcal{B}_\nu(\mathbf{k}) = C_\nu$$

locally defined quantity
(Berry curvature)

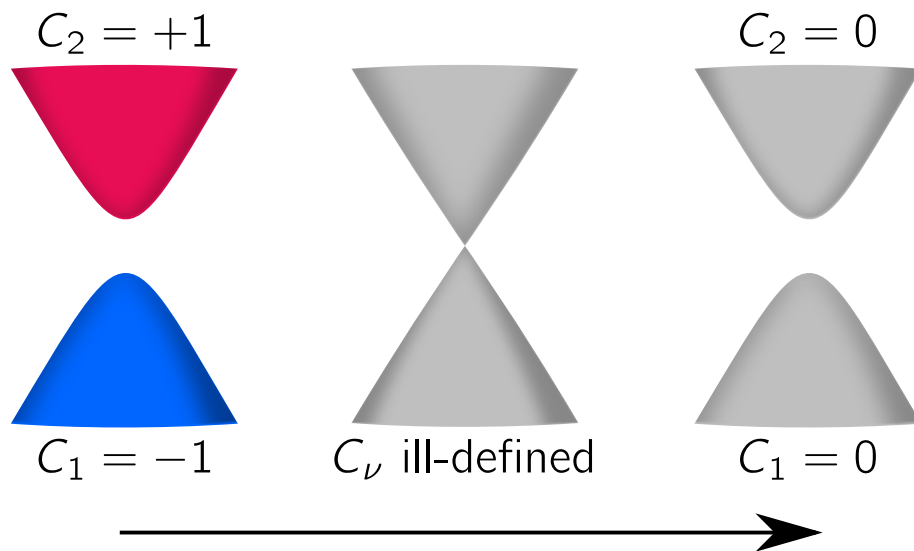
topological index
(Chern number $C_\nu \in \mathbb{Z}$)



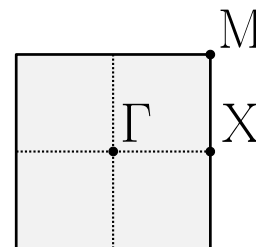
Chern number



$$\frac{1}{2\pi} \int_{\text{BZ}} d^2\mathbf{k} \mathcal{B}_\nu(\mathbf{k}) = C_\nu$$



Chern number



$$\frac{1}{2\pi} \int_{\text{BZ}} d^2\mathbf{k} \mathcal{B}_\nu(\mathbf{k}) = C_\nu$$

$$\mathcal{B}_\nu(\mathbf{k}) = \partial_{k_x} \mathcal{A}_\nu^y(\mathbf{k}) - \partial_{k_y} \mathcal{A}_\nu^x(\mathbf{k})$$

$$\mathcal{A}_\nu^j(\mathbf{k}) = i \langle u_\nu(\mathbf{k}) | \partial_{k_j} | u_\nu(\mathbf{k}) \rangle$$

Berry vector potential
(gauge dependent)

Bloch functions
of the ν -th band

Gapped system: $\mathbf{n}_k = \begin{pmatrix} w \operatorname{Re} \epsilon_k^2 \\ w \operatorname{Im} \epsilon_k^2 \\ \mu + t \epsilon_k^0 \end{pmatrix} \in \mathbb{R}^3 \setminus \{\mathbf{0}\}$

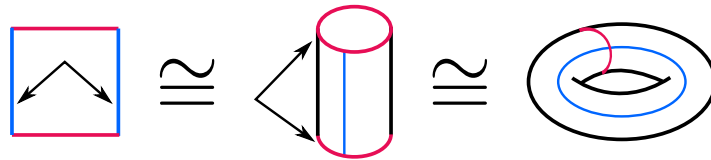
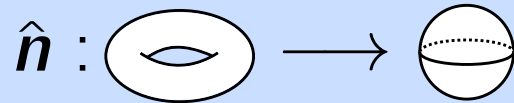
$$E_{\pm}(\mathbf{k}) = n_k^0 \pm |\mathbf{n}_k|$$

Gapped system: $\mathbf{n}_k = \begin{pmatrix} w \operatorname{Re} \epsilon_k^2 \\ w \operatorname{Im} \epsilon_k^2 \\ \mu + t \epsilon_k^0 \end{pmatrix} \in \mathbb{R}^3 \setminus \{\mathbf{0}\}$

$$E_{\pm}(k) = n_k^0 \pm |\mathbf{n}_k|$$

normalized vector!

$$\hat{\mathbf{n}} : T^2 \longrightarrow S^2$$

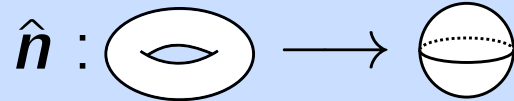


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$$E_{\pm}(\mathbf{k}) = n_k^0 \pm |\mathbf{n}_k|$$

normalized vector!

$$\hat{\mathbf{n}} : T^2 \longrightarrow S^2$$



How can we classify such mappings $\hat{\mathbf{n}}$?

▷ winding number:

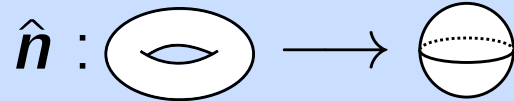
how many times does $\hat{\mathbf{n}}$ wrap the sphere (as we integrate over the BZ)?

Gapped system: $\mathbf{n}_k = \begin{pmatrix} w \operatorname{Re} \epsilon_k^2 \\ w \operatorname{Im} \epsilon_k^2 \\ \mu + t \epsilon_k^0 \end{pmatrix} \in \mathbb{R}^3 \setminus \{\mathbf{0}\}$

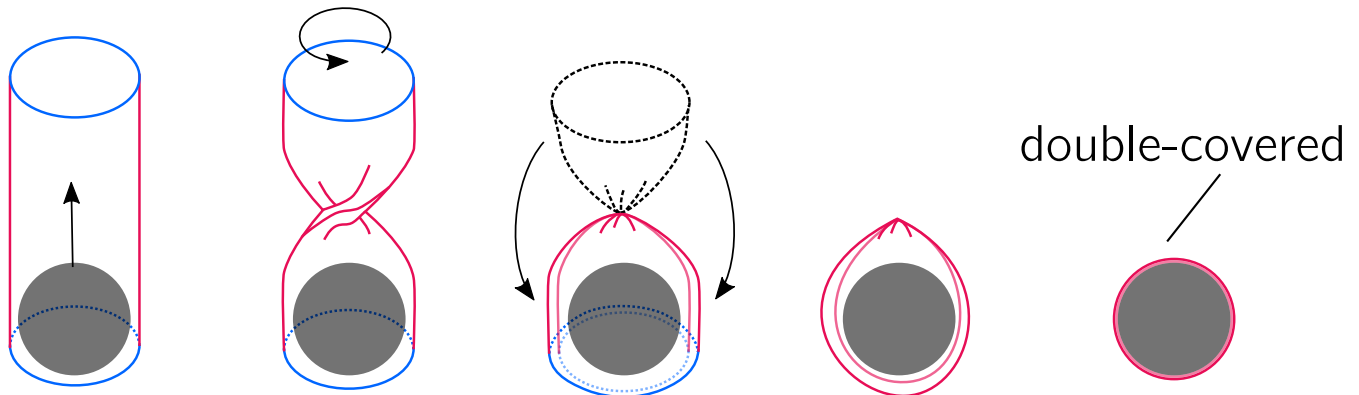
$$E_{\pm}(k) = n_k^0 \pm |\mathbf{n}_k|$$

normalized vector!

$$\hat{\mathbf{n}} : T^2 \longrightarrow S^2$$



How does a nontrivial mapping with $C = 2$ look like?

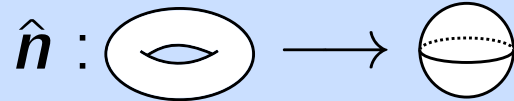


Gapped system: $\mathbf{n}_k = \begin{pmatrix} w \operatorname{Re} \epsilon_k^2 \\ w \operatorname{Im} \epsilon_k^2 \\ \mu + t \epsilon_k^0 \end{pmatrix} \in \mathbb{R}^3 \setminus \{\mathbf{0}\}$

$$E_{\pm}(\mathbf{k}) = n_k^0 \pm |\mathbf{n}_k|$$

normalized vector!

$$\hat{\mathbf{n}} : T^2 \longrightarrow S^2$$



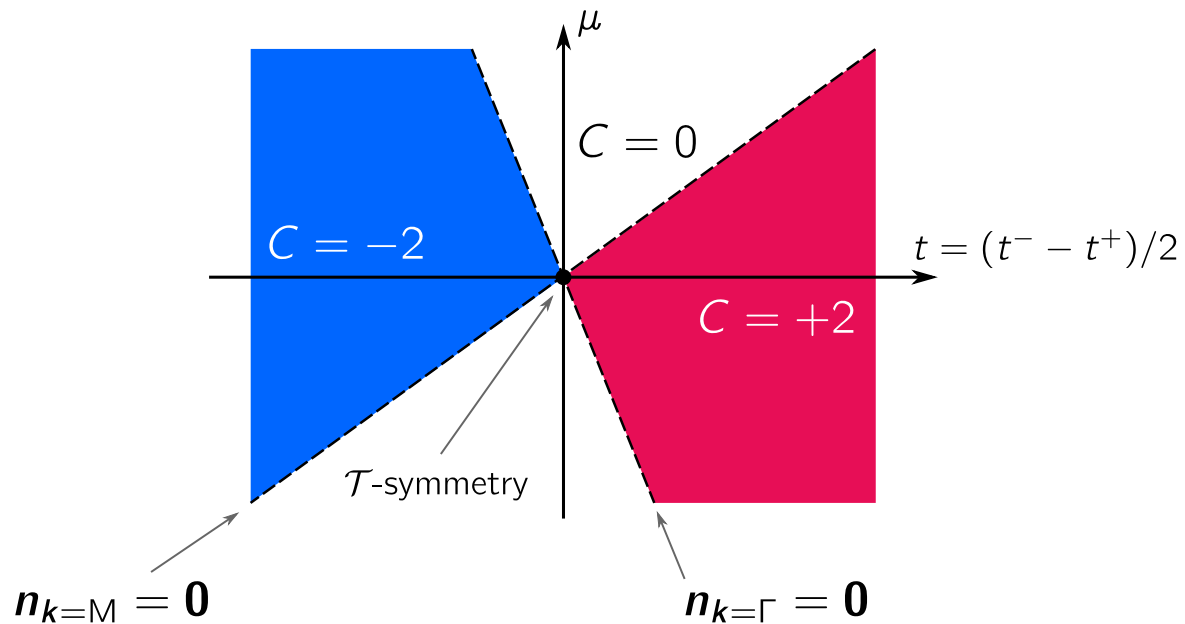
▷ second homotopy group of two-sphere:

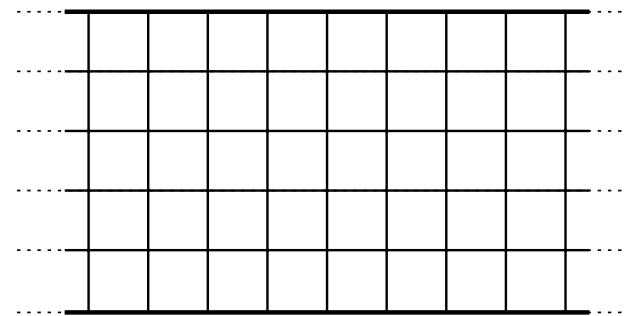
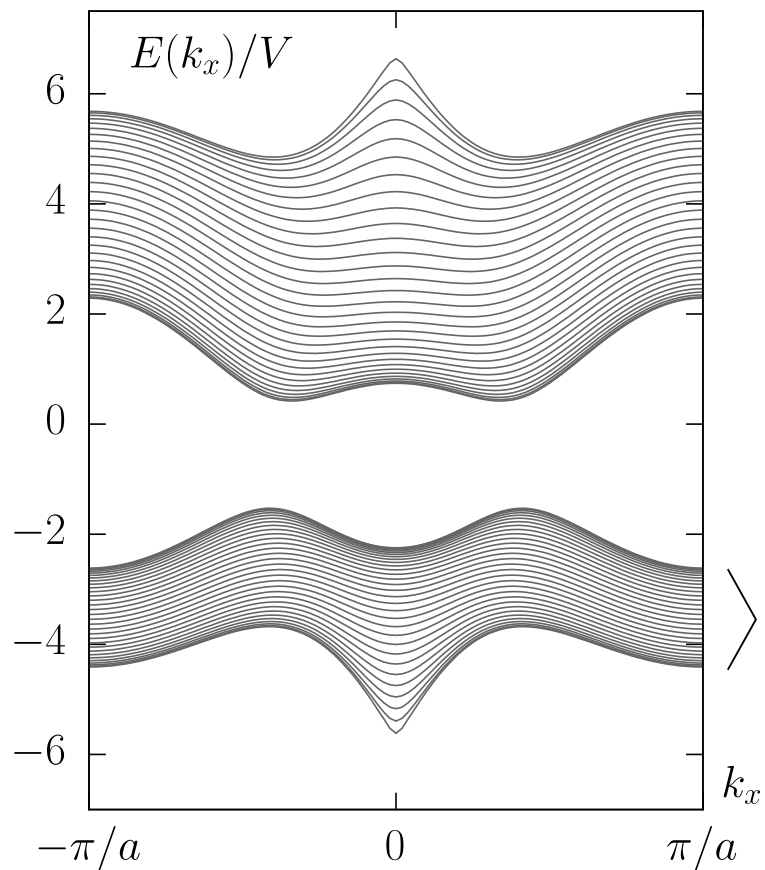
$$\pi_2(S^2) = \mathbb{Z}$$

The Chern number or the winding number of the vector $\hat{\mathbf{n}}$

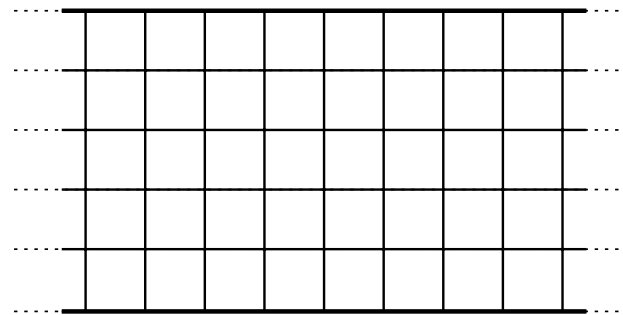
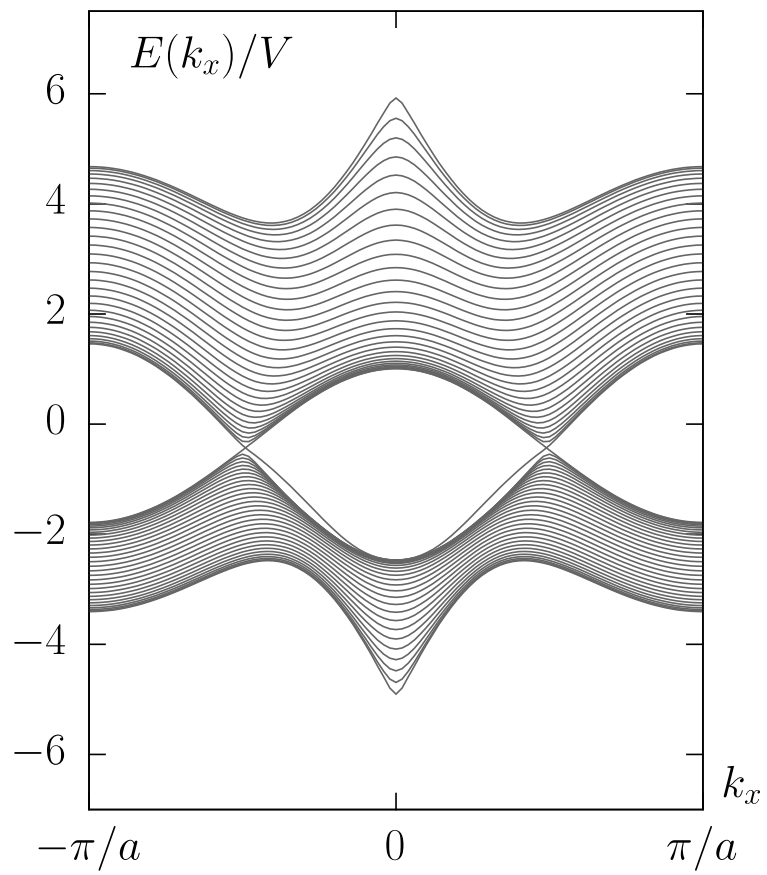
C as winding number

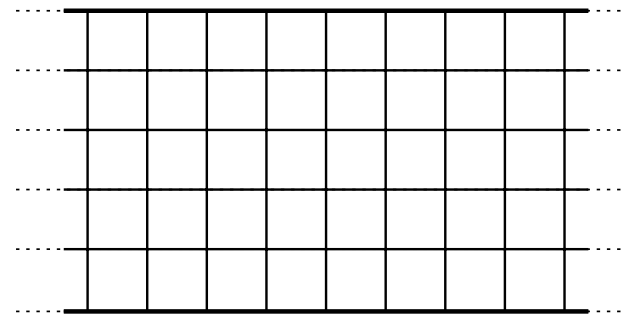
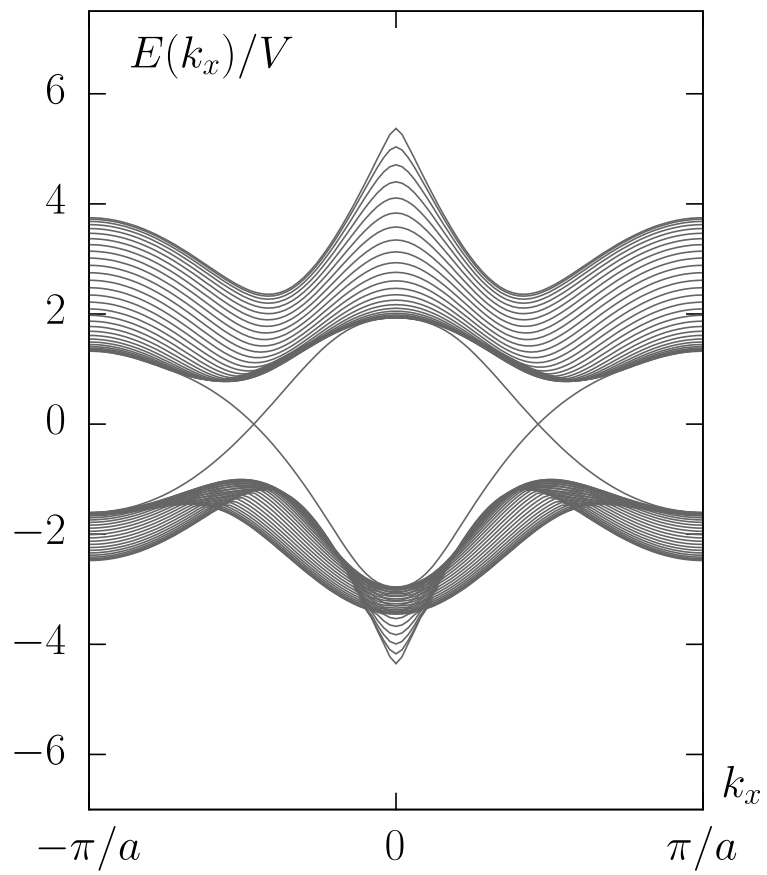
$$C = \frac{1}{4\pi} \int_{\text{BZ}} d^2\mathbf{k} (\partial_{k_x} \hat{\mathbf{n}}_k \times \partial_{k_y} \hat{\mathbf{n}}_k) \cdot \hat{\mathbf{n}}_k$$

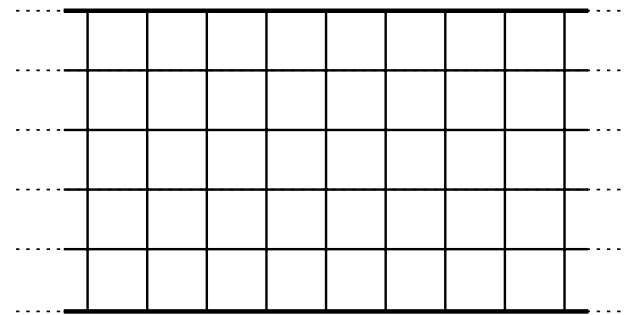
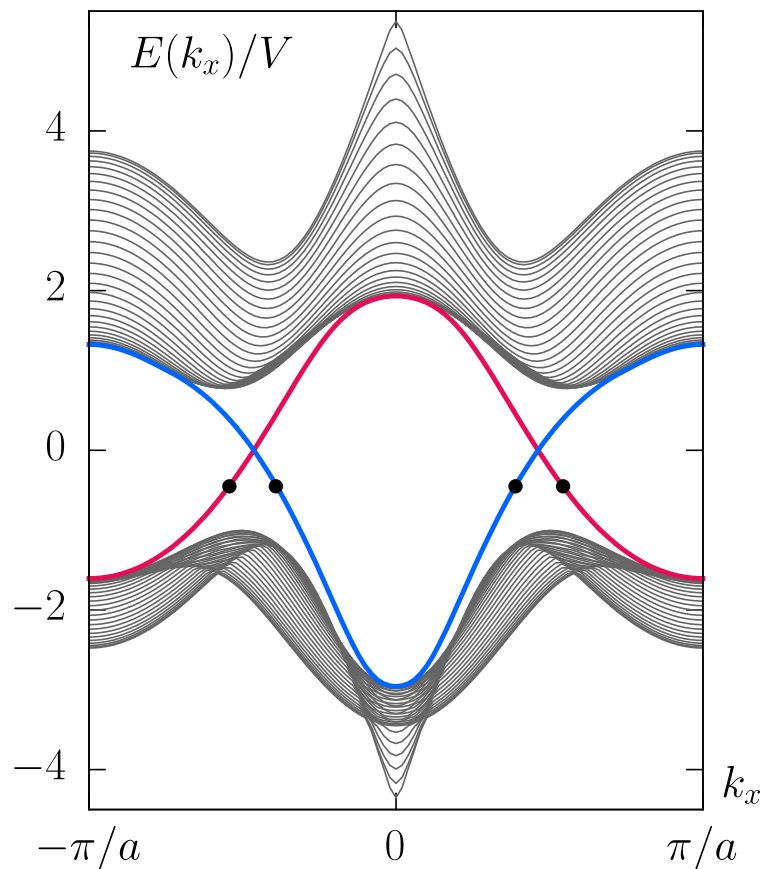




number of bands = N_y







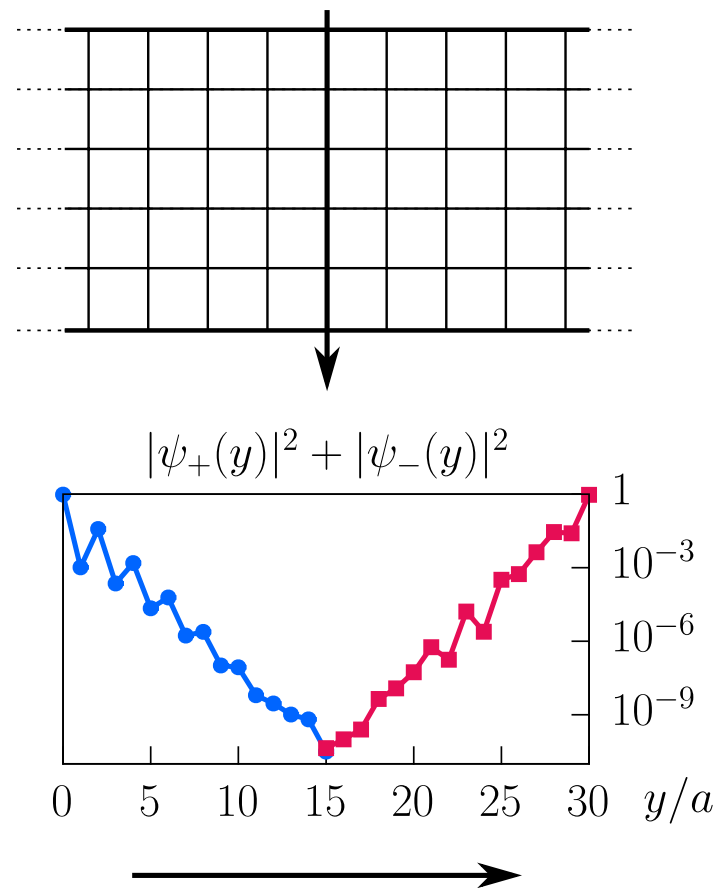
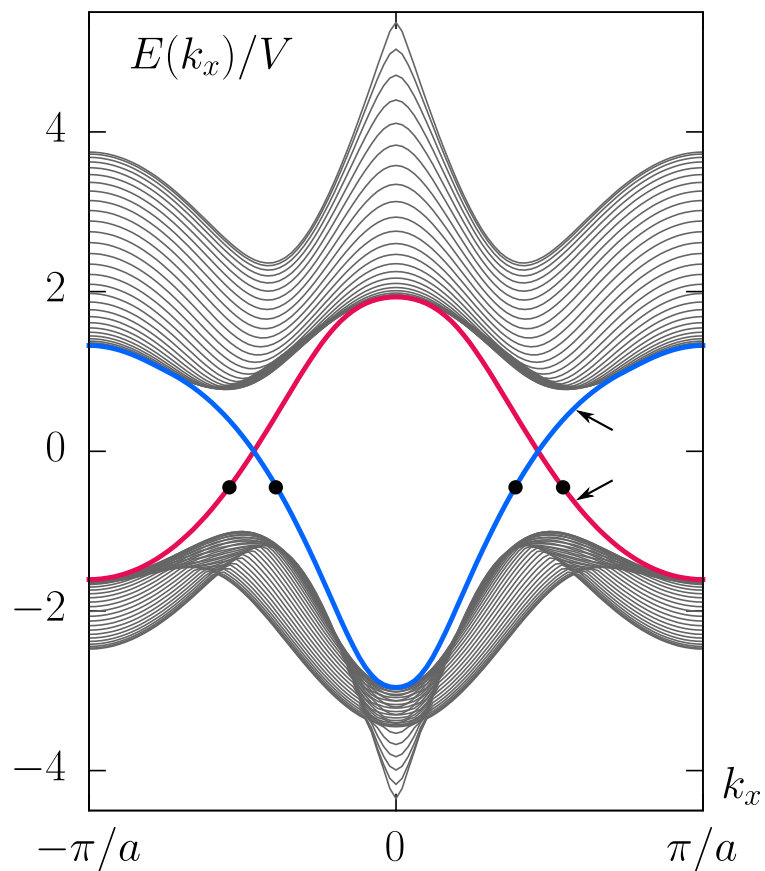
bulk–edge correspondence

Chern number

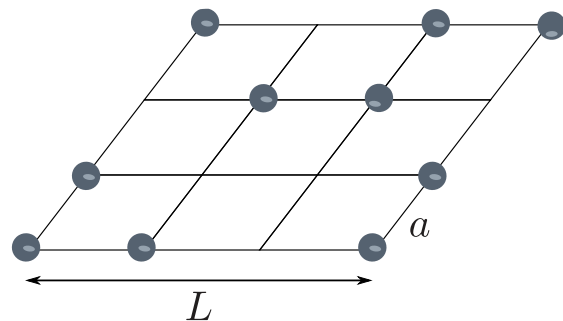
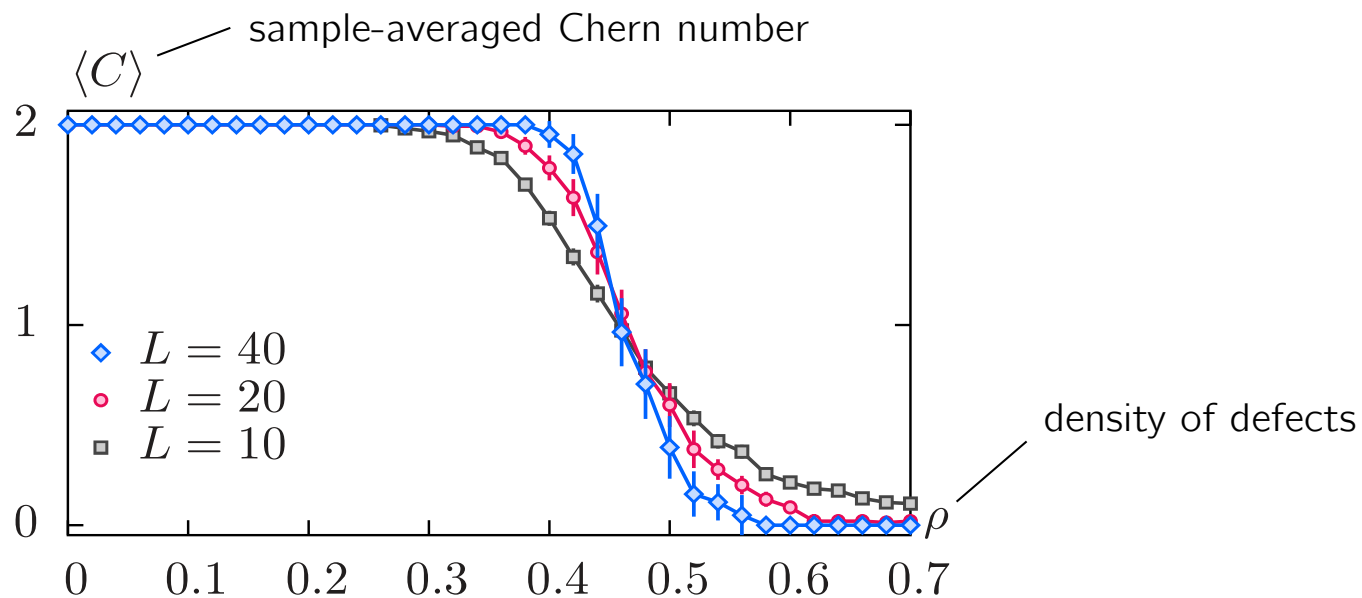
\leftrightarrow

number of edge modes

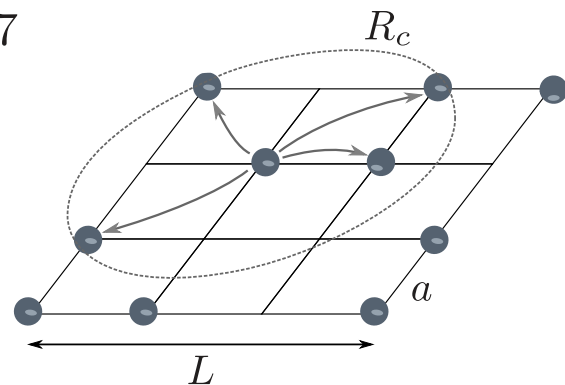
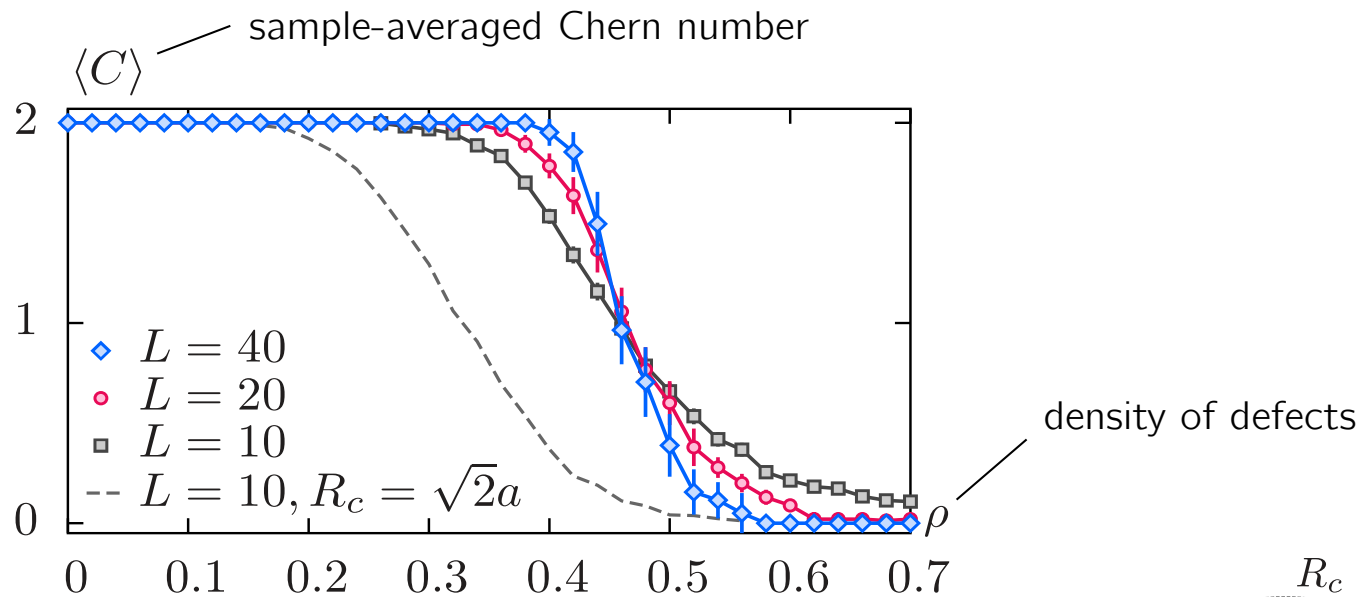
Hatsugai, Phys. Rev. Lett. **71**, 3697 (1993)

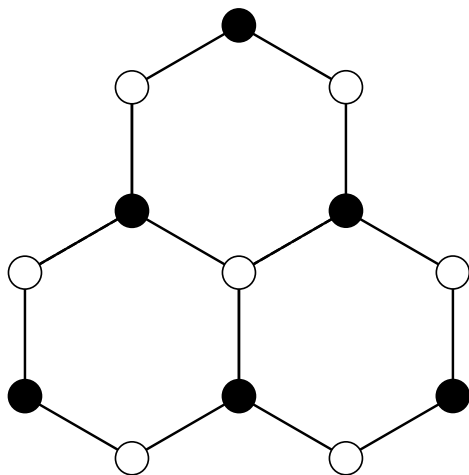


What if the lattice is not perfect?

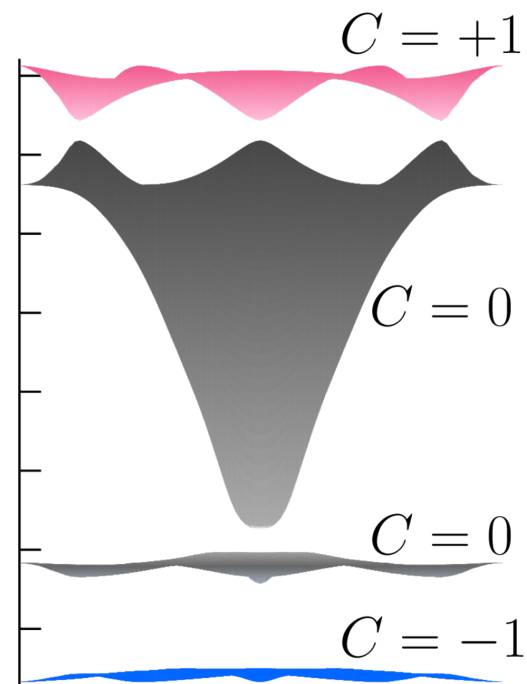


What if the lattice is not perfect?





2 sites per unit cell \times 2 orbitals
 \triangleright 4 bands



flatness = bandwidth/bandgap ≈ 6.4

- ▷ dipolar exchange interactions & broken time-reversal symmetry naturally lead to topological bands
- ▷ Chern number depends on the underlying lattice
square lattice: $C = 2$
honeycomb: $C = 1 - 4$
- ▷ Robust against missing dipoles (lattice sites)
- ▷ Numerical candidate for the interacting $C = 2$ system^{1,2}:

bosonic fractional Chern insulator state at $\nu = 2/3$: Halperin (221) state?

Thank you!

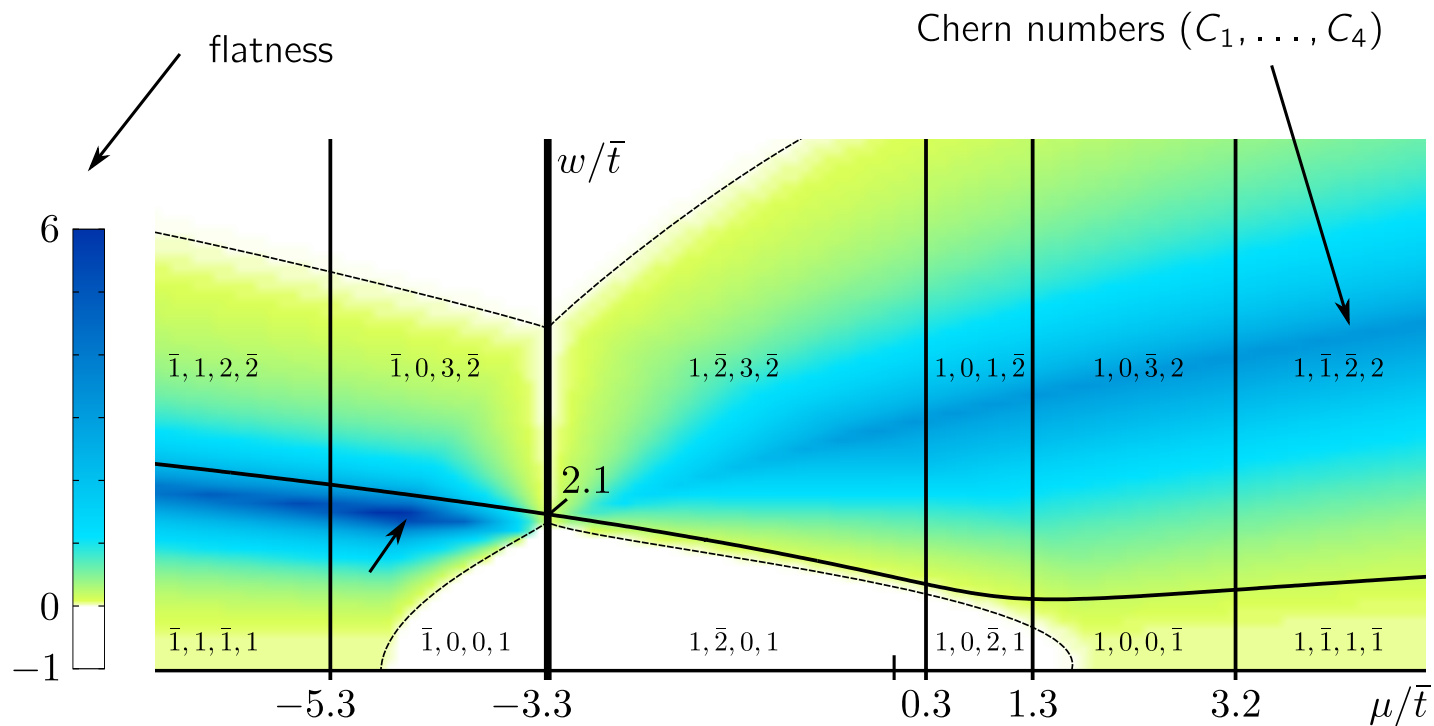
¹Y. Wang, Phys. Rev. B **86**, 201101 (2012)

²G. Möller, Phys. Rev. Lett. **103**, 105303 (2009)

[arXiv:1410.5667](https://arxiv.org/abs/1410.5667)

$$\Psi_{(l;m;n)} = \prod_{i \neq j} (z_i^\downarrow - z_j^\downarrow)^l \prod_{i \neq j} (z_i^\uparrow - z_j^\uparrow)^m \prod_{i,j} (z_i^\uparrow - z_j^\downarrow)^n e^{-\frac{1}{4} \sum_{j,\alpha} |z_j^\alpha|^2}$$

honeycomb lattice: topological phase diagram



How to calculate the Chern number in the disordered system?

impose twisted boundary conditions

$$\psi(x + L, y) = e^{i\theta_x} \psi(x, y)$$

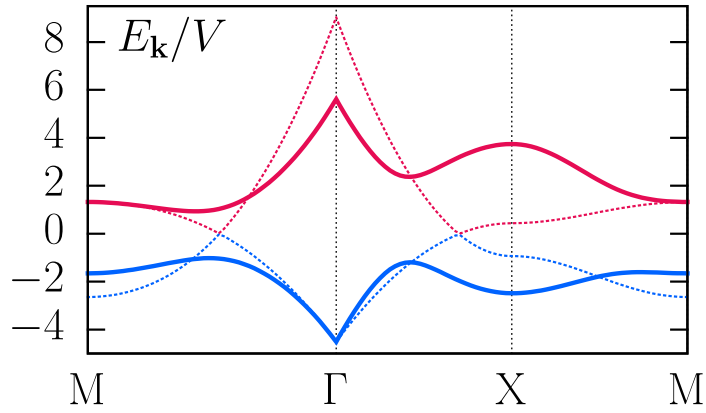
$$\psi(x, y + L) = e^{i\theta_y} \psi(x, y)$$

$$C = \frac{1}{2\pi} \iint d\theta_x d\theta_y F(\theta_x, \theta_y)$$

many-body Berry curvature

Slater-determinant of lower "band" $\Psi(\theta_x, \theta_y)$

$$F(\theta_x, \theta_y) = \text{Im} \left(\left\langle \frac{\partial \Psi}{\partial \theta_y} \middle| \frac{\partial \Psi}{\partial \theta_x} \right\rangle - \left\langle \frac{\partial \Psi}{\partial \theta_x} \middle| \frac{\partial \Psi}{\partial \theta_y} \right\rangle \right)$$



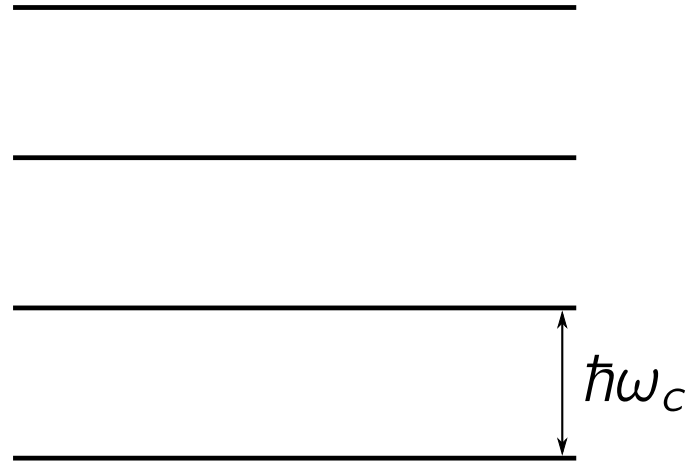
Analogy with Landau levels:

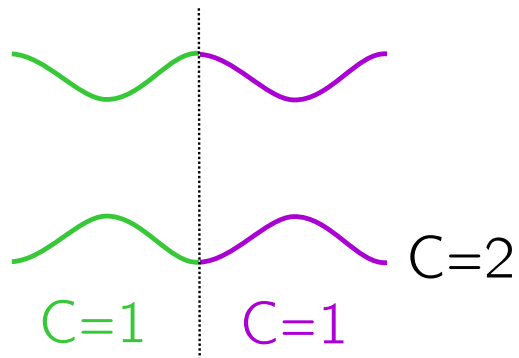
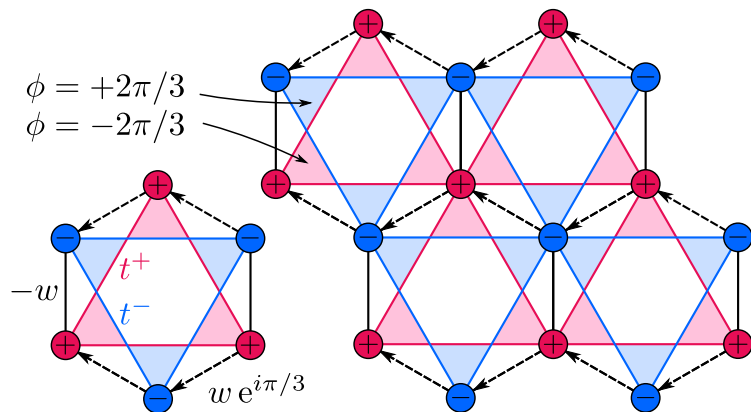
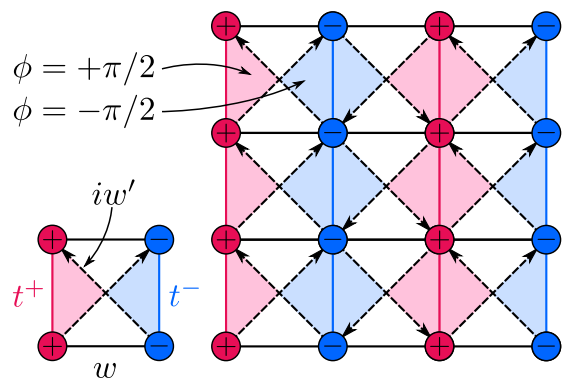
$$H = \hbar\omega_c \left(n + \frac{1}{2} \right)$$

with $\omega_c = eB/m$

▷ perfectly flat band(s) with
Chern number $C = 1$

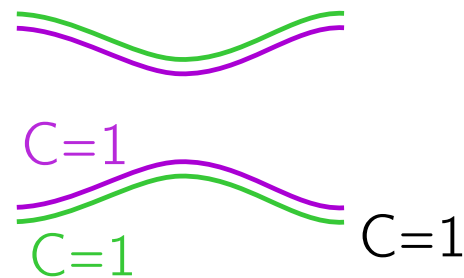
Hall conductivity: $\sigma_{xy} = \frac{e^2}{h} \times C$





full BZ

$$1 + 1 = 2$$



full BZ

$$1 + 1 = 2 \times 1$$

Setup:

- ▷ Polar molecules
- ▷ Dipolar exchange interactions

Two-band model

- ▷ Mapping to bosons
- ▷ Introduction to topological bands

Square lattice:

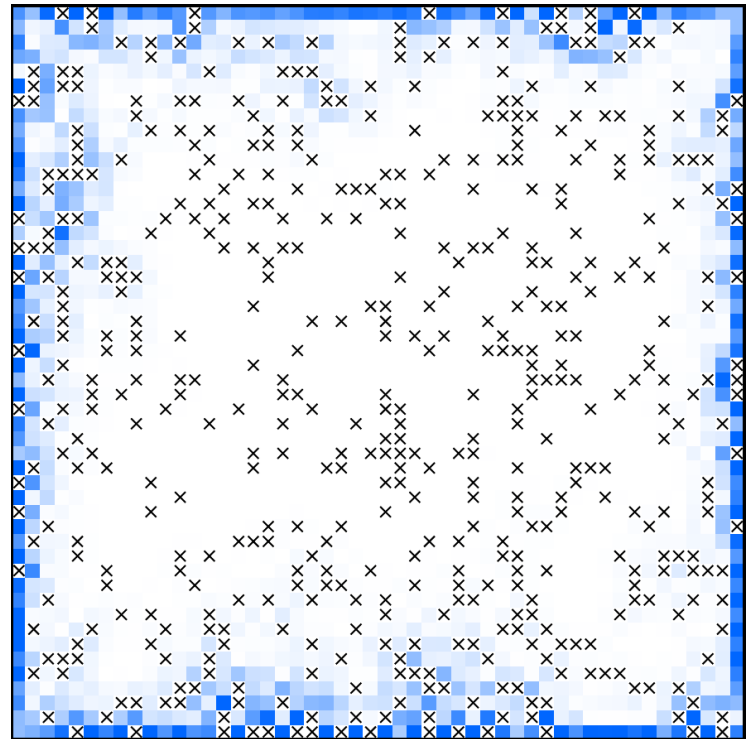
- ▷ Edge states
- ▷ Robustness against disorder

Honeycomb lattice:

- ▷ flat bands
- ▷ phase diagram

What if the lattice is not perfect?

edge states are robust
against disorder:



20% of the molecules randomly removed