Topological Floquet States in Two-Dimensional Triangle Shaking Optical Lattice

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Outline

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- Floquet Theoty and Effective Hamiltonian
- Model
- Solving and Results
- Discussion for Further Study

For a time-dependent Hamiltonian H(t), Shrödinger equation reads

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$$i\hbar\partial_t\psi=H\psi$$

Time-dependent system

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$$i\hbar\partial_t\psi=H\psi$$

For a periodically driven system,

$$H(t+T) = H(t)$$
$$T = \frac{2\pi}{\omega}$$

we have

$$\psi_{\lambda} = e^{-i(\lambda t/\hbar)} u_{\lambda}(\vec{r}, \omega t)$$

Time-dependent system

we have

$$\psi_{\lambda} = e^{-i(\lambda t/\hbar)} u_{\lambda}(\vec{r}, \omega t)$$

and

$$u_{\lambda}(\vec{r}, \omega(t+T)) = u_{\lambda}(\vec{r}, \omega t)$$

$$\hat{F} = \hat{U}(T_i + T, T_i)$$

$$= \hat{T} \exp\left(-i \int_{T_i}^{T_i + T} \hat{H}(t) dt\right)$$

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 $\{\varepsilon_n\}$: the quasi-energy bands

Effective Hamiltonian

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Effective Hamiltonian

$$\hat{F} = e^{-\mathfrak{i}H_{\mathrm{eff}}T}$$

$$\hat{H}(t) = \sum_{n = -\infty}^{\infty} \hat{H}_n(t)e^{i\omega t}$$

$$\hat{H}_{\mathrm{eff}} = \hat{H}_0 + \sum_{n=1}^{\infty} \left(\frac{[\hat{H}_n, \hat{H}_{-n}]}{n\omega} - \frac{[\hat{H}_n, \hat{H}_0]}{e^{-2\pi n i \alpha} n \omega} + \frac{[\hat{H}_{-n}, \hat{H}_0]}{e^{2\pi n i \alpha} n \omega} \right)$$

$$\mathcal{H}(t) = T + \mathcal{V}(\vec{r}, t)$$

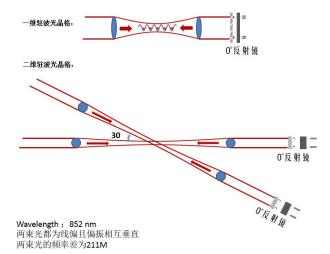
Hamiltonian

$$\mathcal{H}(t) = T + \mathcal{V}(\vec{r}, t)$$

$$\mathcal{V} = -V_1 \cos(2\vec{k_1} \cdot \vec{r_1} + f \cos(\omega t))$$
$$-V_2 \cos(2\vec{k_2} \cdot \vec{r_2} + f \cos(\omega t + \alpha))$$

Setup

Setup



Translation

Translation

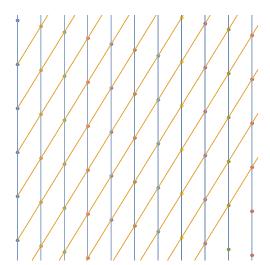
Translational transformation into the co-moving frame of reference,

$$\vec{r} \rightarrow \vec{r} + \vec{\xi}(t)$$

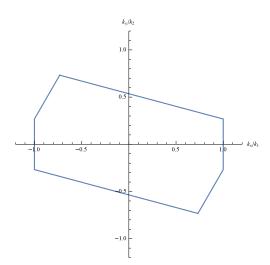
$$H(t) = \frac{p^2}{2\mu} - V_1 \cos(2\vec{k_1} \cdot \vec{r_1}) - V_2 \cos(2\vec{k_2} \cdot \vec{r_2})$$

$$+ i\frac{\hbar\omega f}{2k_0} \left[\left(\sin(\omega t) + \sin(\omega t + \alpha) \cos(30^\circ) \right) \partial_x - \sin(\omega t + \alpha) \sin(30^\circ) \partial_y \right]$$

Lattice Structure

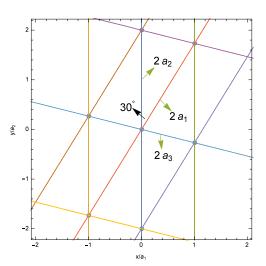


First Brillouin Zone in Reciprocal Lattice



Hopping term

Hopping term



Eigen equation reads

$$H|\psi_q^{(n)}(\vec{r})\rangle = \varepsilon^{(n)}(\vec{k})|\psi_q^{(n)}(\vec{r})\rangle$$

Eigen problem under Bloch bases

Eigen equation reads

$$H|\psi_q^{(n)}(\vec{r})\rangle = \varepsilon^{(n)}(\vec{k})|\psi_q^{(n)}(\vec{r})\rangle$$

with

$$H = \frac{p^2}{2m} + V_1 \cos(2\vec{k_1} \cdot \vec{r_1}) + V_2 \cos(2\vec{k_2} \cdot \vec{r_2})$$

Eigen equation reads

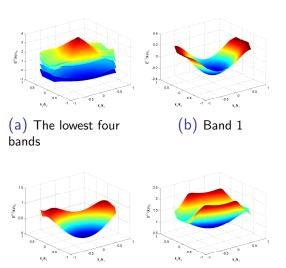
$$H|\psi_q^{(n)}(\vec{r})\rangle = \varepsilon^{(n)}(\vec{k})|\psi_q^{(n)}(\vec{r})\rangle$$

with

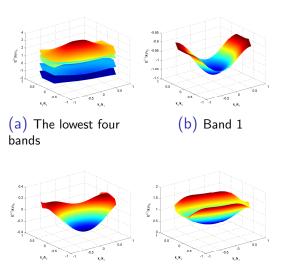
$$H = \frac{p^2}{2m} + V_1 \cos(2\vec{k_1} \cdot \vec{r_1}) + V_2 \cos(2\vec{k_2} \cdot \vec{r_2})$$

$$\psi_q(\vec{r}) = \sum_{l_1, l_2} C_{l_1, l_2}(\vec{q}) e^{i(2l_1 \vec{k_1} + 2l_2 \vec{k_2} + \vec{q}) \cdot \vec{r}}$$

Energy bands with V=1.0

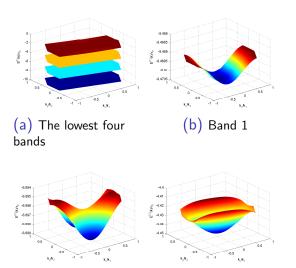


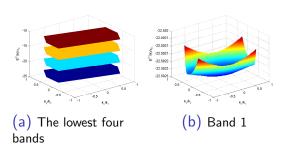
Energy bands with V=2.0

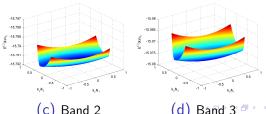


Band 2

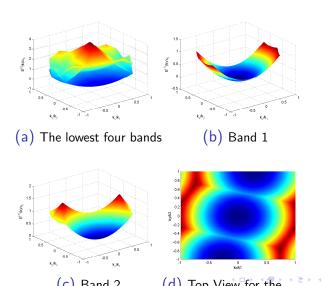
Energy bands with V=8.0







Energy bands with $V=0.05(\approx 0)$



Wannier functions

Wannier functions

$$w_n(\vec{R}_m, \vec{r}) = w_n(\vec{r} - \vec{R}_m)$$

$$= \frac{1}{\sqrt{N}} \sum_{\vec{k} \in BZ} e^{-i\vec{k} \cdot \vec{R}_m} \psi_{n\vec{k}}(\vec{r})$$

$$w_n(\vec{R}_m, \vec{r}) = w_n(\vec{r} - \vec{R}_m)$$

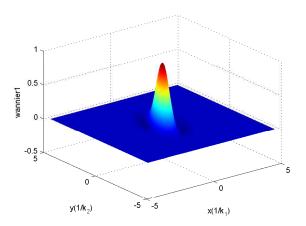
$$= \frac{1}{\sqrt{N}} \sum_{\vec{k} \in BZ} e^{-i\vec{k} \cdot \vec{R}_m} \psi_{n\vec{k}}(\vec{r})$$

$$w_{n}(\vec{r}) = w_{n}(0, \vec{r})$$

$$= \frac{1}{\sqrt{N}} \sum_{\vec{q} \in BZ} \psi_{n\vec{q}}(\vec{r})$$

$$= \frac{1}{\sqrt{N}} \sum_{\vec{q} \in BZ} \sum_{l_{1}, l_{2}} C_{l_{1}, l_{2}}(\vec{q}) e^{i(2l_{1}\vec{k}_{1} + 2l_{2}\vec{k}_{2} + \vec{q}) \cdot \vec{r}}$$

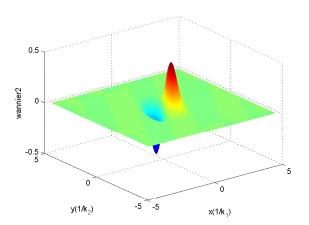
Wannier 1



(b) Wannier function for the 1st band $w_1(x,y)$



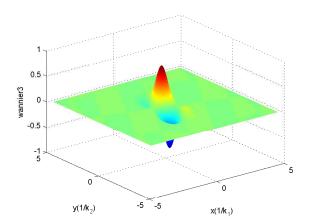
Wannier 2



(d) Wannier function for the 2^{nd} band $w_2(x,y)$



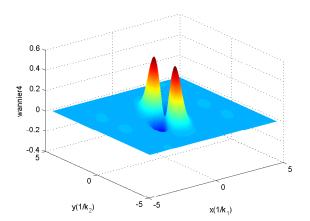
Wannier 3



(f) Wannier function for the 3^{th} band $w_3(x,y)$

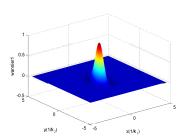


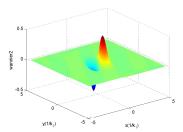
Wannier 4



(h) Wannier function for the 4^{th} band $w_4(x,y)$

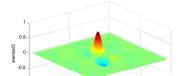


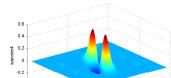




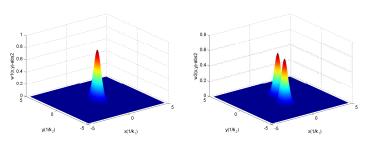
band $w_1(x,y)$

(a) Wannier function for the 1st (b) Wannier function for the 2nd band $w_2(x,y)$









(a) Square Modulus of $w_1(x,y)$ (b) Square Modulus of $w_2(x,y)$

Figure: Square modulus of the Wannier functions in Figure 6

Tight-bounding model

$$\hat{H}(t) = \sum_{\vec{k}} \left(\hat{a}_{s\vec{k}}^{\dagger} \ \hat{a}_{p\vec{k}}^{\dagger} \right) H(\vec{k},t) \left(\begin{array}{c} \hat{a}_{s\vec{k}} \\ \hat{a}_{p\vec{k}} \end{array} \right)$$

Tight-bounding model

$$\hat{H}(t) = \sum_{\vec{k}} \left(\hat{a}_{s\vec{k}}^{\dagger} \ \hat{a}_{p\vec{k}}^{\dagger} \right) H(\vec{k},t) \left(\begin{array}{c} \hat{a}_{s\vec{k}} \\ \hat{a}_{p\vec{k}} \end{array} \right)$$

$$H(\vec{k},t) = H_0(\vec{k}) + H_1(\vec{k})e^{i\omega t} + H_{-1}(\vec{k})e^{-i\omega t}$$

Tight-bounding model

$$\hat{H}(t) = \sum_{\vec{k}} \left(\hat{a}_{s\vec{k}}^{\dagger} \ \hat{a}_{p\vec{k}}^{\dagger} \right) H(\vec{k},t) \left(\begin{array}{c} \hat{a}_{s\vec{k}} \\ \hat{a}_{p\vec{k}} \end{array} \right)$$

$$H(\vec{k},t) = H_0(\vec{k}) + H_1(\vec{k})e^{i\omega t} + H_{-1}(\vec{k})e^{-i\omega t}$$

$$H_0(\vec{k}) = \begin{pmatrix} E_s & iK \\ -iK & E_p \end{pmatrix}$$

$$H_1(\vec{k}) = \begin{pmatrix} \Lambda_s & \Omega \\ -\Omega & \Lambda_p \end{pmatrix}$$

$$H_{-1}(\vec{k}) = H_1(\vec{k})^{\dagger} = \begin{pmatrix} \Lambda_s^* & -\Omega^* \\ \Omega^* & \Lambda_n^* \end{pmatrix}$$

One-photon resonance

$$\begin{split} \widetilde{H}_{\text{eff}}(\vec{k}) &= \varepsilon_0 + \vec{B}(\vec{k}) \cdot \vec{\sigma} \\ &= \varepsilon_0 + B_x(\vec{k}) \sigma_x + B_y(\vec{k}) \sigma_y + B_z(\vec{k}) \sigma_z \end{split}$$

Solving and Results

One-photon resonance

$$\begin{split} \hat{H}_{\text{eff}}(\vec{k}) &= \varepsilon_0 + \vec{B}(\vec{k}) \cdot \vec{\sigma} \\ &= \varepsilon_0 + B_x(\vec{k})\sigma_x + B_y(\vec{k})\sigma_y + B_z(\vec{k})\sigma_z \\ \varepsilon_0 &= \frac{1}{2}(E_s + \omega + E_p) \\ B_x(\vec{k}) &= \Re(\Omega) + \frac{K}{\omega} \big(\Re(\Lambda_p) - \Im(\Lambda_s) \big) \\ B_y(\vec{k}) &= -\Im(\Omega) + \frac{K}{\omega} \big(\Re(\Lambda_p) - \Re(\Lambda_s) \big) \\ B_z(\vec{k}) &= \frac{E_s + \omega - E_p}{2} - \frac{K^2}{\omega} - \frac{|\Omega|^2}{2\omega} \end{split}$$

Two-photon resonance

$$H_{\text{eff}} = \frac{E_s + E_p}{2} + \left[\frac{E_s - E_p + 2\omega}{2} - \frac{1}{\omega} \left(\frac{4}{3} |\Omega|^2 + \frac{K^2}{2} \right) \right] \sigma_z + \frac{1}{\omega} \Re \left(\Omega(\Lambda_s - \Lambda_p) \right) \sigma_x - \frac{1}{\omega} \Im \left(\Omega(\Lambda_s - \Lambda_p) \right) \sigma_y$$

$$H_{\text{eff}} = \varepsilon_0 + \vec{B}(\vec{k}) \cdot \vec{\sigma}$$

= $\varepsilon_0 + \vec{B}_z(\vec{k}) \cdot \vec{\sigma_z} + \vec{B}_x(\vec{k}) \cdot \vec{\sigma_x} + \vec{B}_y(\vec{k}) \cdot \vec{\sigma_y}$

with

$$\varepsilon_0 = \frac{E_s + E_p}{2}
\vec{B}_z(\vec{k}) = \frac{E_s - E_p + 2\omega}{2} - \frac{1}{\omega} \left(\frac{4}{3} |\Omega|^2 + \frac{K^2}{2} \right)
\vec{B}_x(\vec{k}) = \frac{1}{\omega} \Re \left(\Omega(\Lambda_s - \Lambda_p) \right)
\vec{B}_y(\vec{k}) = -\frac{1}{\omega} \Im \left(\Omega(\Lambda_s - \Lambda_p) \right)$$

Outline

$$H_{\rm eff} = \vec{B}(\vec{k}) \cdot \vec{\sigma} = \varepsilon(\mathbf{k}) \hat{\mathbf{n}}(\mathbf{k}) \cdot \tilde{\sigma}$$

Chern numbers

$$H_{\mathrm{eff}} = \vec{B}(\vec{k}) \cdot \vec{\sigma} = \varepsilon(\mathbf{k}) \hat{\mathbf{n}}(\mathbf{k}) \cdot \tilde{\sigma}$$

$$C_{\pm} = \frac{\pm 1}{4\pi} \int_{\mathsf{FB7}} \mathbf{\hat{n}} \cdot (\partial_{\mathbf{k_x}} \mathbf{\hat{n}} \times \partial_{\mathbf{k_y}} \mathbf{\hat{n}}) \mathbf{d^2 k}$$

Outline

Chern numbers

One-photon resonance

None topological non-trivial states found.

Chern numbers

One-photon resonance

None topological non-trivial states found.

Tne-photon resonance

For small potential, V < 1.5, non-zero Chern number cases found. Shaking amplitudes f and frequencies ω allowed for topological non-trivial states differs for different potential depths. While the range of f is approximately several k_0 , the range of ω allowed from tens to hundruds units comparing to the energy unit of the system.

Thank you!