

Moiré Superlattices at Fractional Band Fillings: Particle-Hole Duality and Quantum Geometry

Ahmed Abouelkomsan

AA, Zhao Liu & Emil Bergholtz, Phys. Rev. Lett. **124**, 106803 (2020)

Zhao Liu, **AA** & Emil Bergholtz, Phys. Rev. Lett. **126**, 026801 (2021)

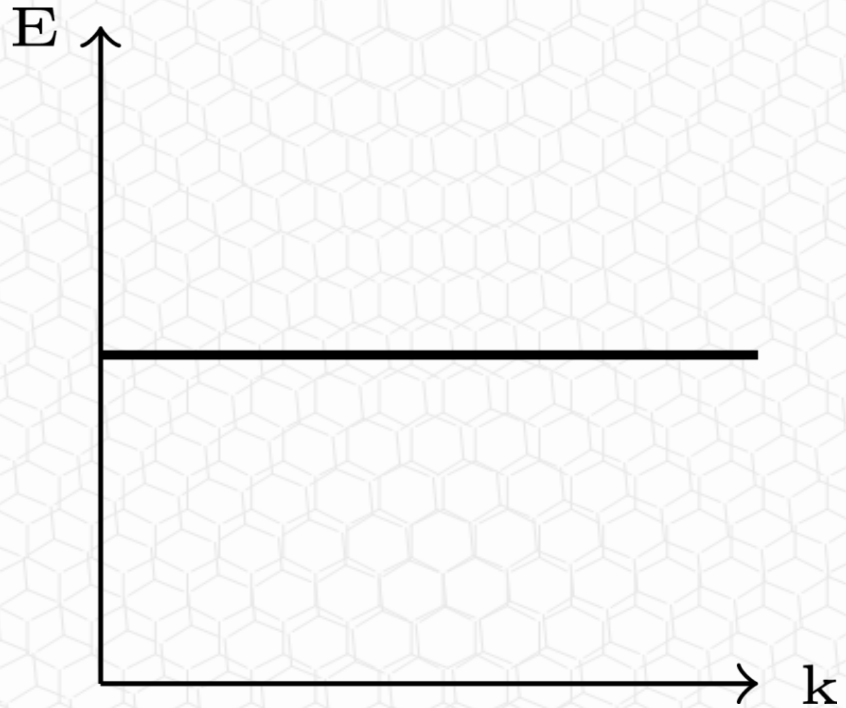
AA, Kang Yang & Emil Bergholtz, arXiv:2202.10467

*Knut and Alice
Wallenberg
Foundation*



The Problem

A Flat Band in a Moiré System



The Problem

A Flat Band



Fractional filling



Fractional band filling \rightarrow non integer filling per Moiré unit cell

The Problem

A Flat Band



Add interactions at
fractional band fillings

$$V(\mathbf{q}) \sim \frac{1}{\sqrt{\mathbf{q}^2 + \kappa^2}}$$



What is the underlying phase ?

Fractional Chern insulators, Charge density waves, Fermi
Liquids, etc. ???

Fractional Chern Insulators, why bother?

- Broadly fractional Chern insulators are lattice analogues of the fractional quantum Hall effect.

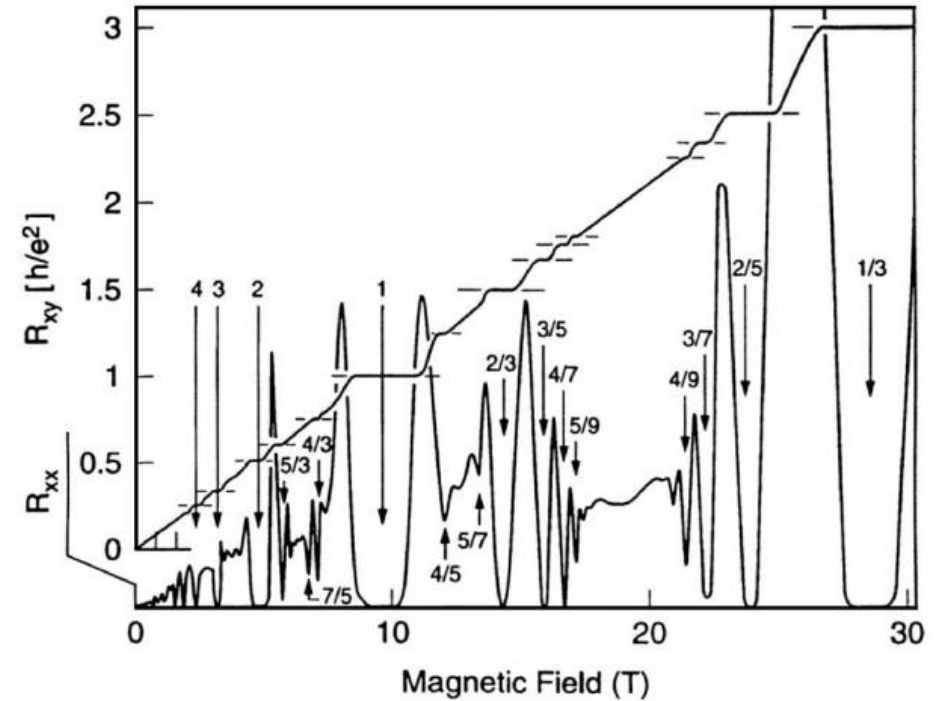


Fig from : Stormer, Physica B **177**, 1-4 (1992)

- Klitzing et al. Phys. Rev. Lett. **45**, 494 (1980)
- Tsui et al. Phys. Rev. Lett. **48**, 1559 (1982)

Fractional Chern Insulators, why bother?

- Overcomes challenges with the conventional FQHE experimental setup!

*Very strong magnetic
fields*



$|\mathbf{B}| \sim 10 - 30$ Tesla



*Extremely low
temperatures*

$T \lesssim 1$ Kelvin

Fractional Chern Insulators, why bother?

- No magnetic field required!

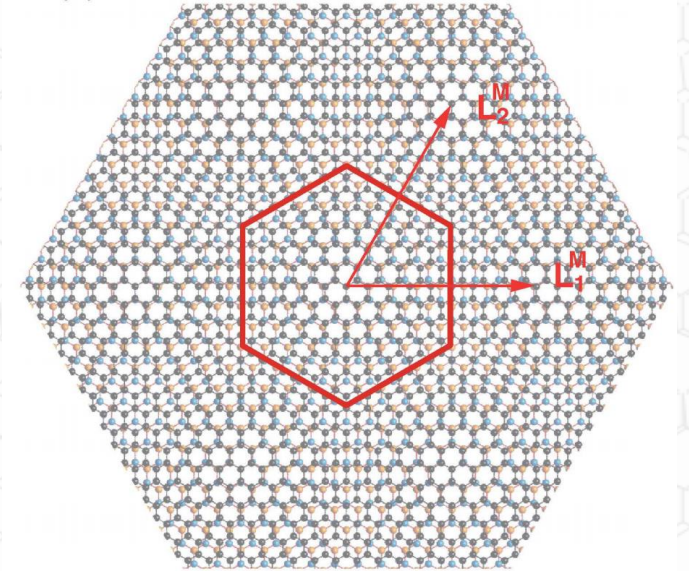
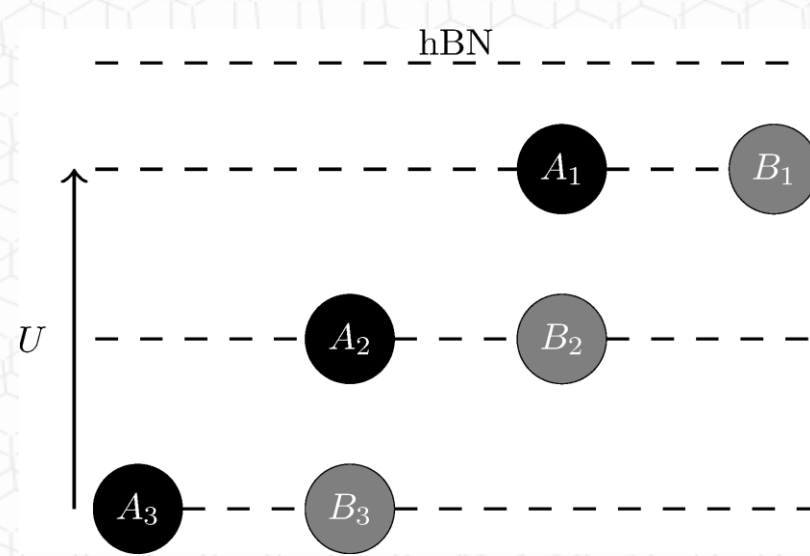
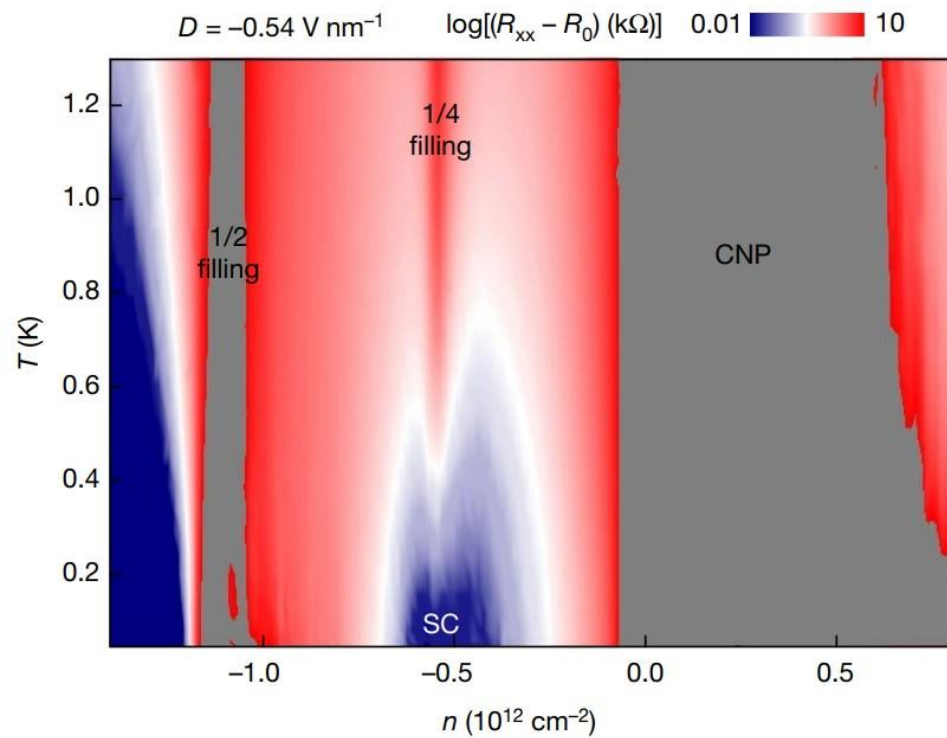


- Interactions on the lattice scale are greater than the magnetic length scale -> Higher energy gap!
- A step towards high temperature topological phases.



- More than FQHE!
- Higher Chern number FCIs are possible -> no mapping to decoupled Landau levels! ,
PRL, **109**, 186805 (2012)

Trilayer Graphene aligned with Boron Nitride



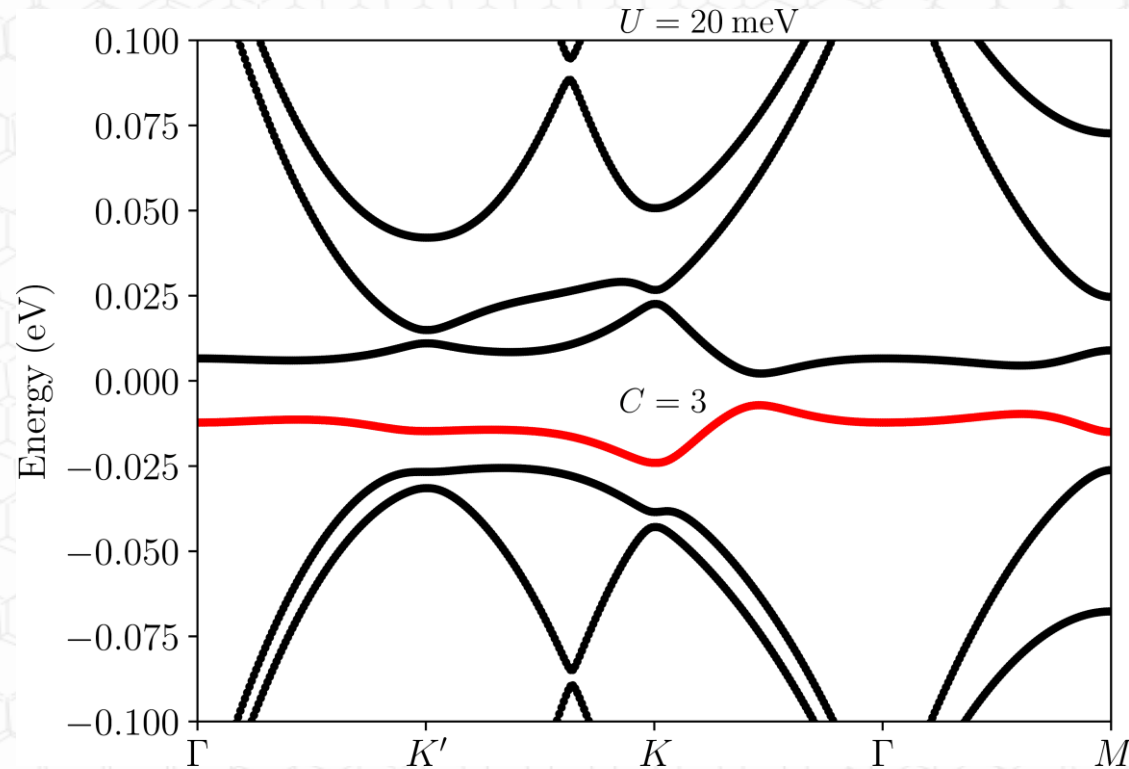
Phys. Rev. B **90**, 155406 (2014)

- Correlated Insulators and superconductors!

Chen et al. Nature **572**, (2019)

Trilayer Graphene aligned with Boron Nitride

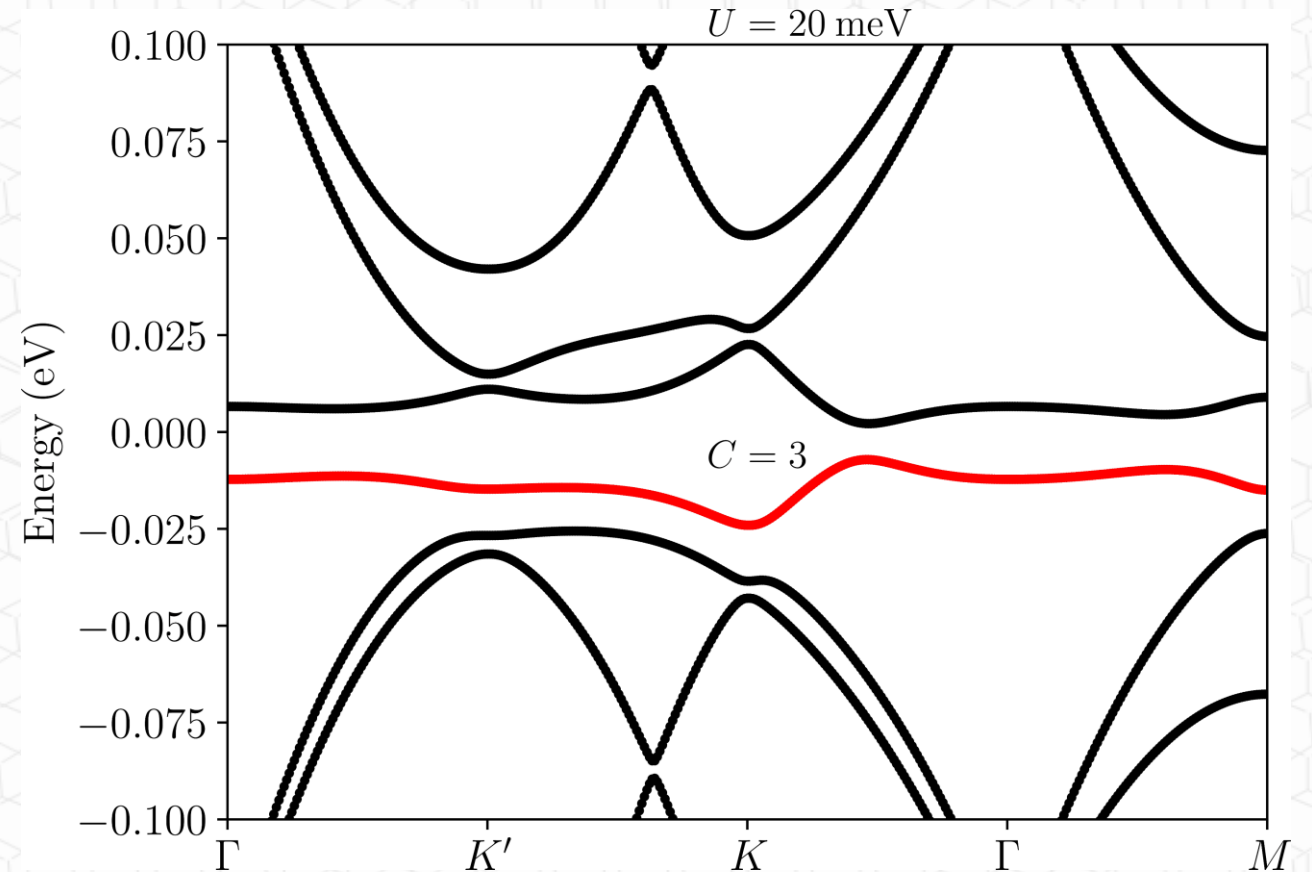
Phys. Rev. Lett. 122, 016401 (2019)



What happens when the red band is fractionally filled?

Trilayer Graphene aligned with Boron Nitride

- Are FCI states possible?
- No numerical evidence!
- Why?



Another perspective

$$H_{\text{proj}} = \sum_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4} V_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4} c_{\mathbf{k}_1}^\dagger c_{\mathbf{k}_2}^\dagger c_{\mathbf{k}_3} c_{\mathbf{k}_4}$$

$\{c_{\mathbf{k}_i}\}$ are band operators!

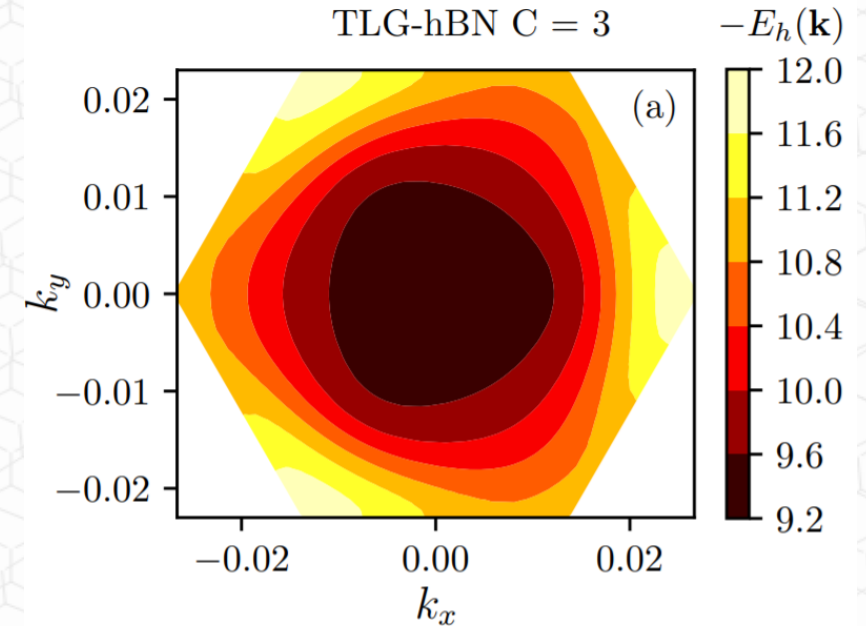
Particle-Hole Transformation, $c_{\mathbf{k}} \rightarrow d_{\mathbf{k}}^\dagger$

$$H_{\text{proj}} \rightarrow \sum_{\mathbf{k}} E_h(\mathbf{k}) d_{\mathbf{k}}^\dagger d_{\mathbf{k}} + \sum_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4} V_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4}^* d_{\mathbf{k}_1}^\dagger d_{\mathbf{k}_2}^\dagger d_{\mathbf{k}_3} d_{\mathbf{k}_4}$$

$$E_h(\mathbf{k}) = \sum_{\mathbf{k}'} (V_{\mathbf{k}' \mathbf{k} \mathbf{k}' \mathbf{k}} + V_{\mathbf{k} \mathbf{k}' \mathbf{k} \mathbf{k}'} - V_{\mathbf{k} \mathbf{k}' \mathbf{k}' \mathbf{k}} - V_{\mathbf{k}' \mathbf{k} \mathbf{k} \mathbf{k}'})$$

Dispersive!
for projected interactions

Unlike Landau levels!



Another perspective

$$H_{\text{proj}} = \sum_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4} V_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4} c_{\mathbf{k}_1}^\dagger c_{\mathbf{k}_2}^\dagger c_{\mathbf{k}_3} c_{\mathbf{k}_4}$$

Particle-Hole Transformation,

$$H_{\text{proj}} \rightarrow \sum_{\mathbf{k}} E_h(\mathbf{k}) d_{\mathbf{k}}^\dagger d_{\mathbf{k}} + \sum_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4} V_{\mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \mathbf{k}_4}^* d_{\mathbf{k}_1}^\dagger d_{\mathbf{k}_2}^\dagger d_{\mathbf{k}_3} d_{\mathbf{k}_4}$$

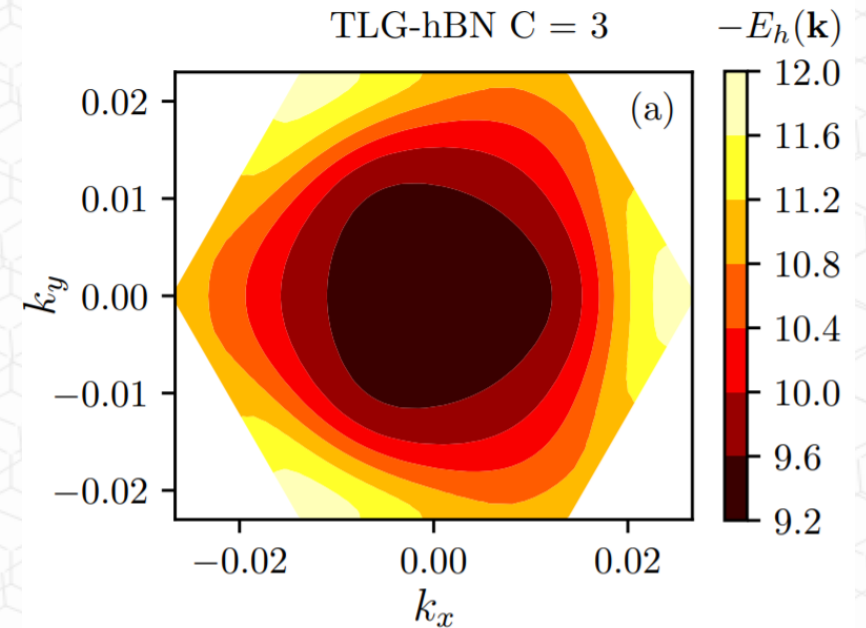
$$E_h(\mathbf{k}) = \sum_{\mathbf{k}'} (V_{\mathbf{k}' \mathbf{k} \mathbf{k}' \mathbf{k}} + V_{\mathbf{k} \mathbf{k}' \mathbf{k} \mathbf{k}'} - V_{\mathbf{k} \mathbf{k}' \mathbf{k}' \mathbf{k}} - V_{\mathbf{k}' \mathbf{k} \mathbf{k} \mathbf{k}'})$$

What role does $E_h(\mathbf{k})$ play?

Is there a quantity that $E_h(\mathbf{k})$ correlates with?

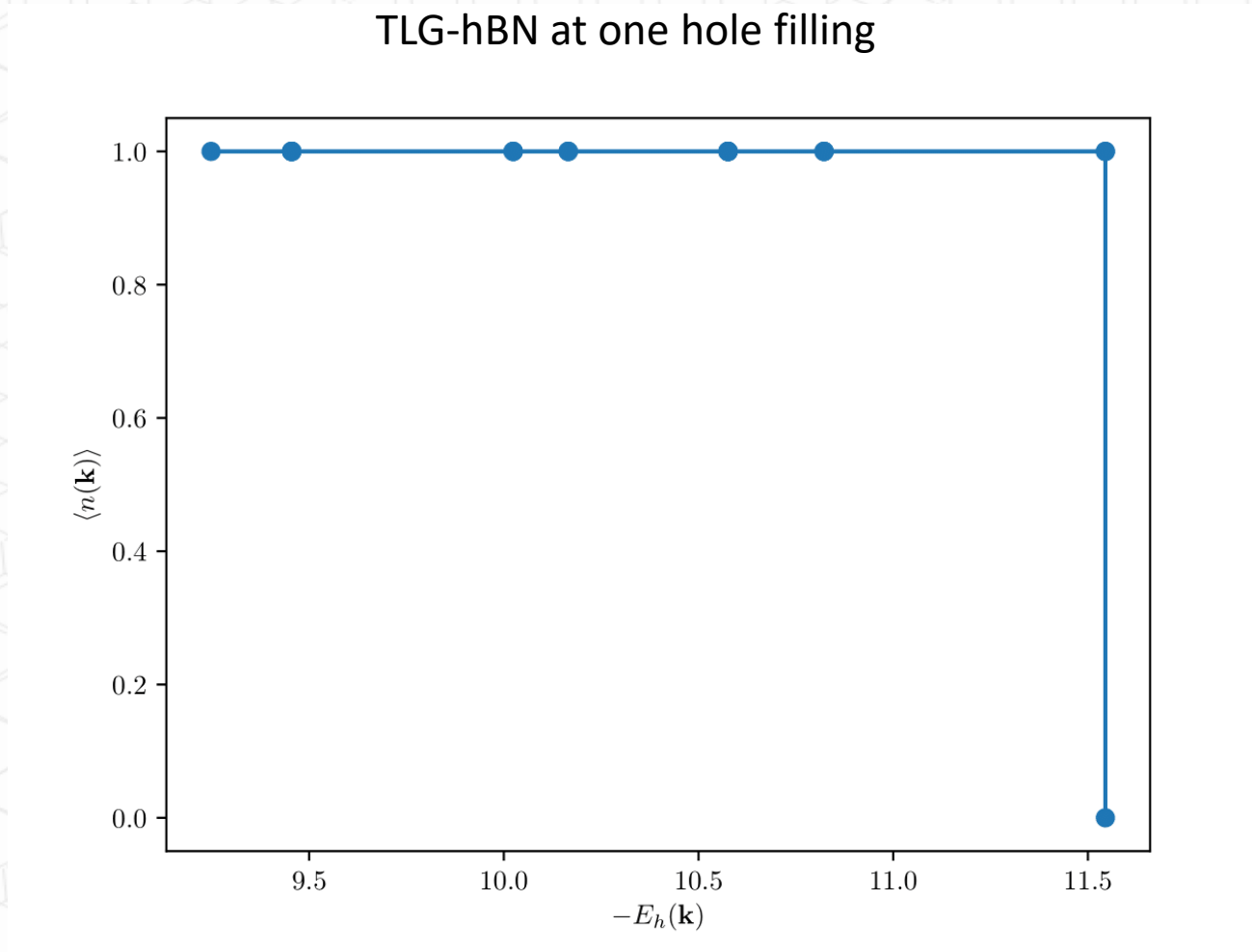
Look at the electron occupation in the GS of H_{proj}

vs $E_h(\mathbf{k})$



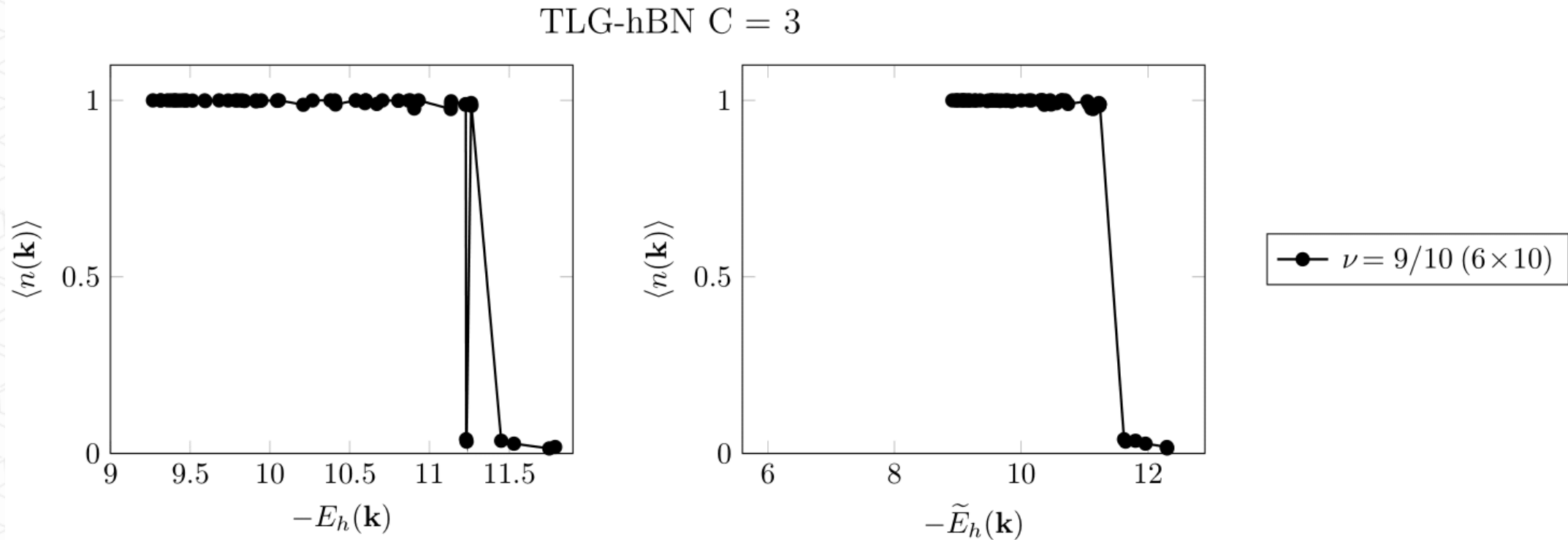
$$\frac{W_h(\text{Hole dispersion bandwidth})}{W(\text{Flatband bandwidth})} \sim 5$$

Trivial example:



$\langle n(\mathbf{k}) \rangle = \langle c_{\mathbf{k}}^\dagger c_{\mathbf{k}} \rangle$ is electron occupation in the many-body ground state obtained from exact diagonalization

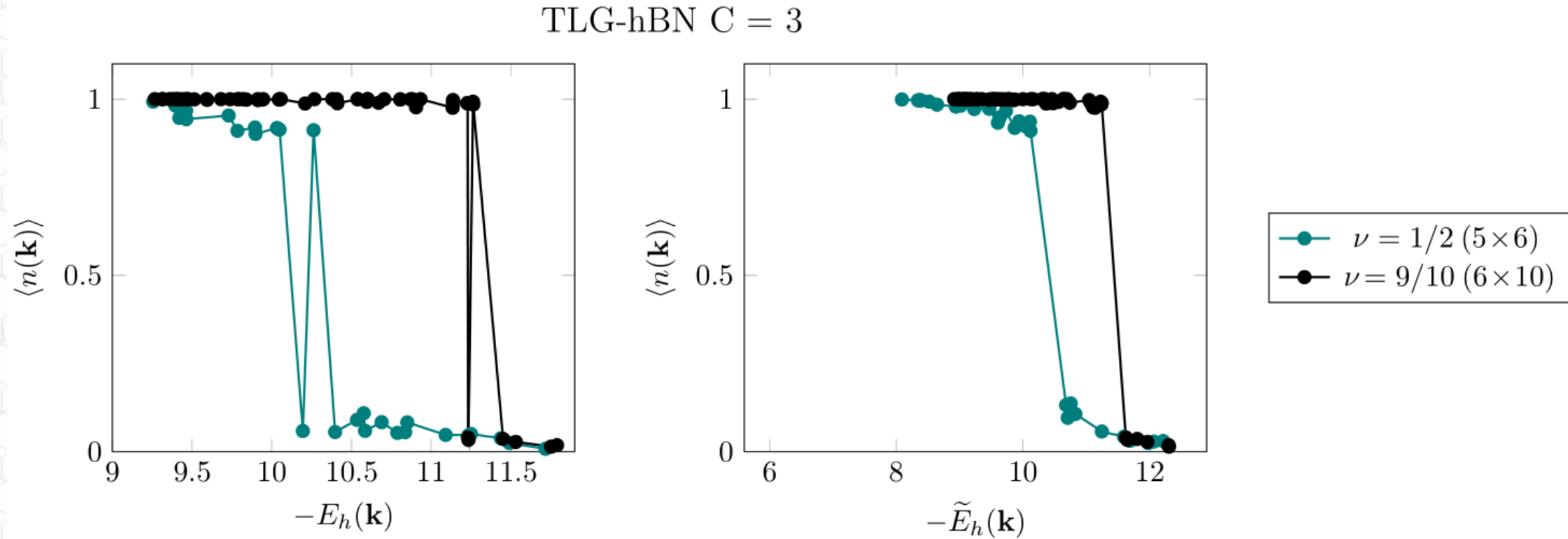
Emergent Fermi Liquids in TLG-hBN



$\langle n(\mathbf{k}) \rangle = \langle c_{\mathbf{k}}^\dagger c_{\mathbf{k}} \rangle$ is electron occupation in the many-body ground state obtained from exact diagonalization

$\tilde{E}_h(\mathbf{k})$ is the renormalized single-hole energy after Hartree-Fock mean field of the interaction

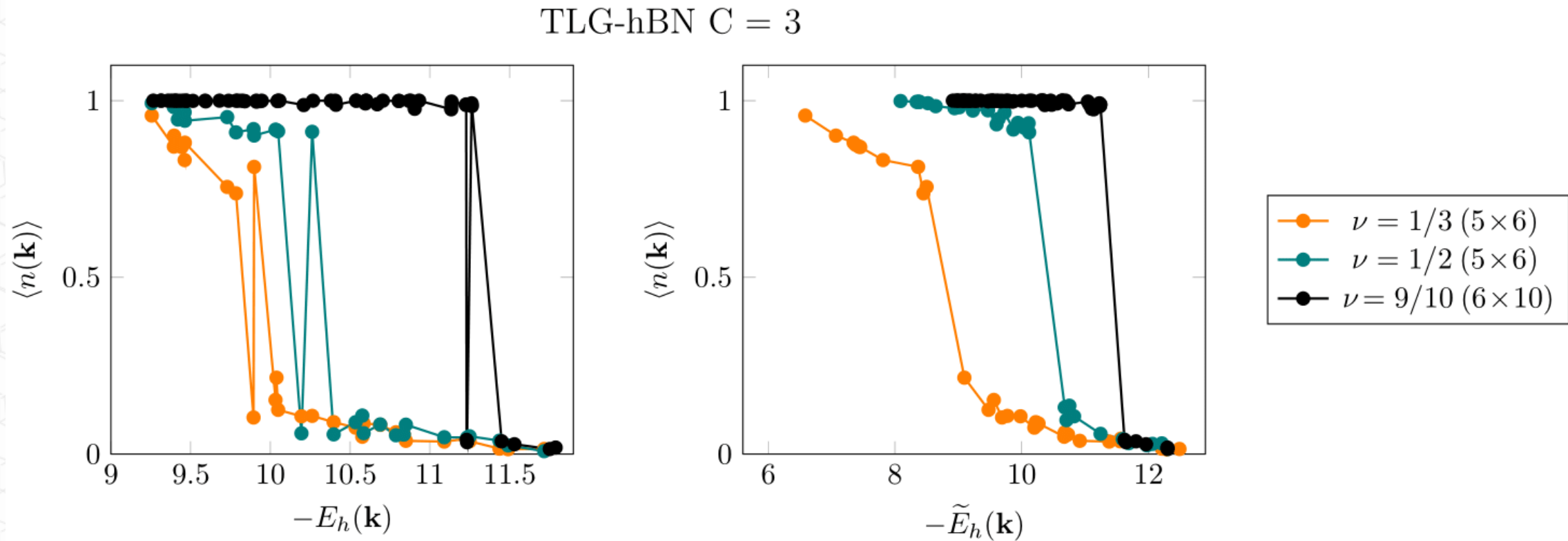
Emergent Fermi Liquids in TLG-hBN



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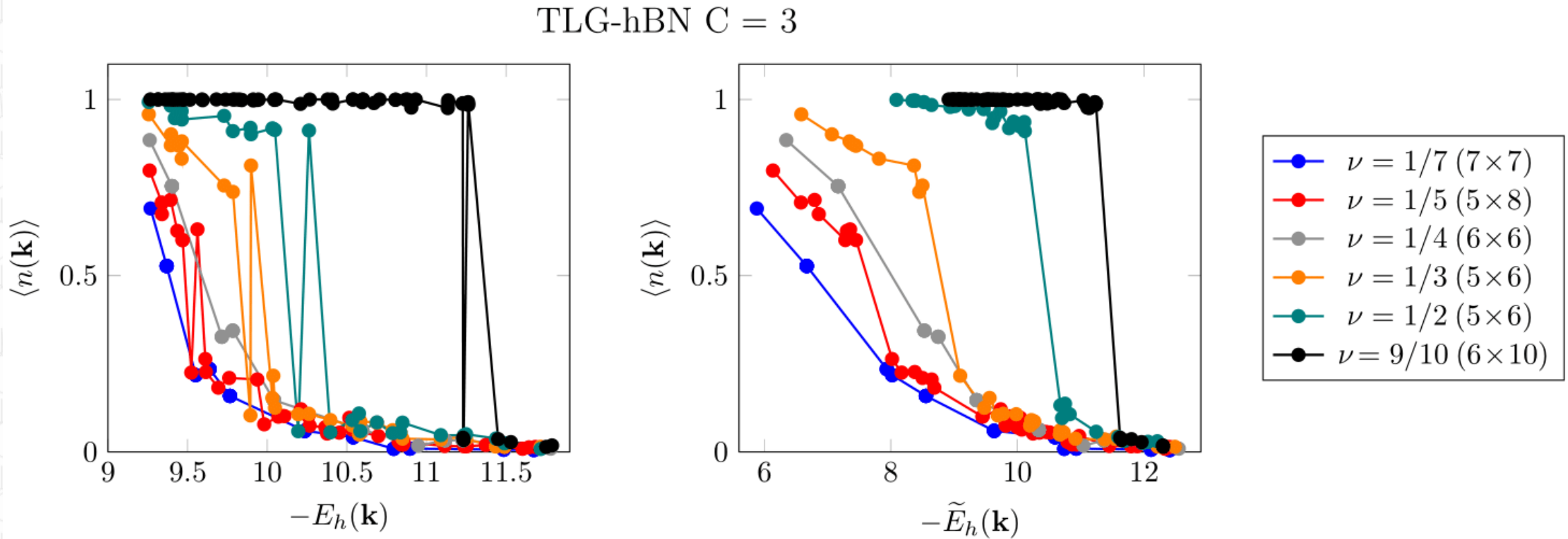
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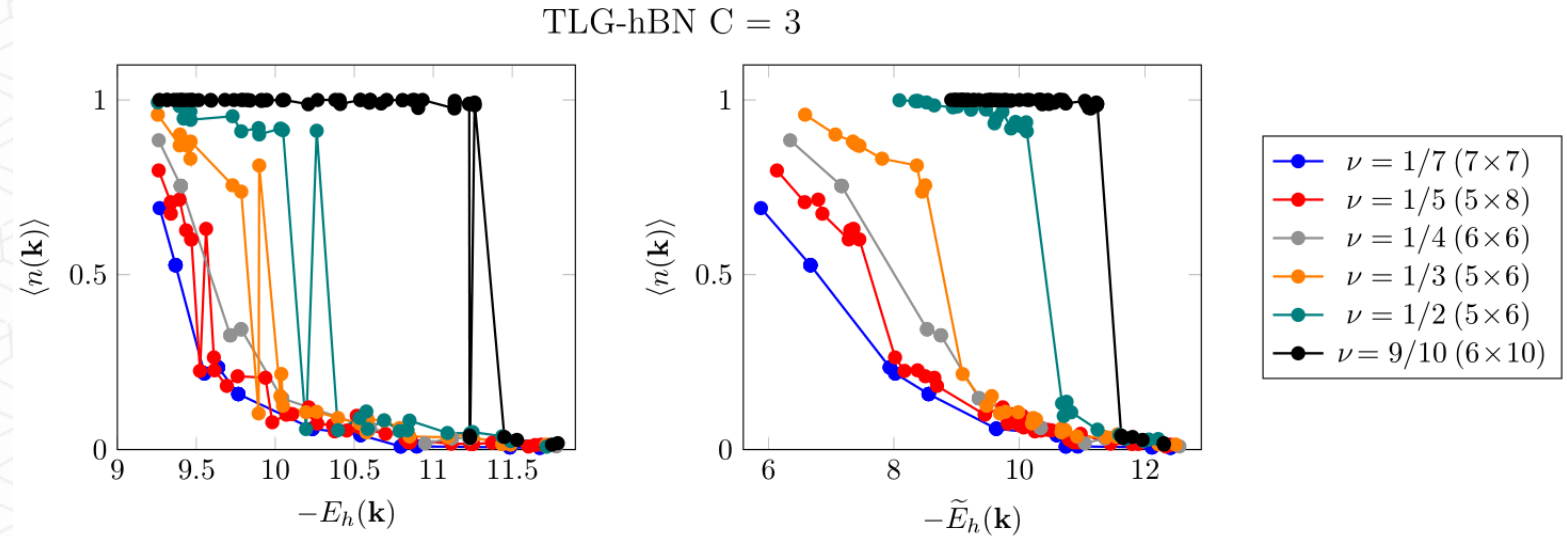
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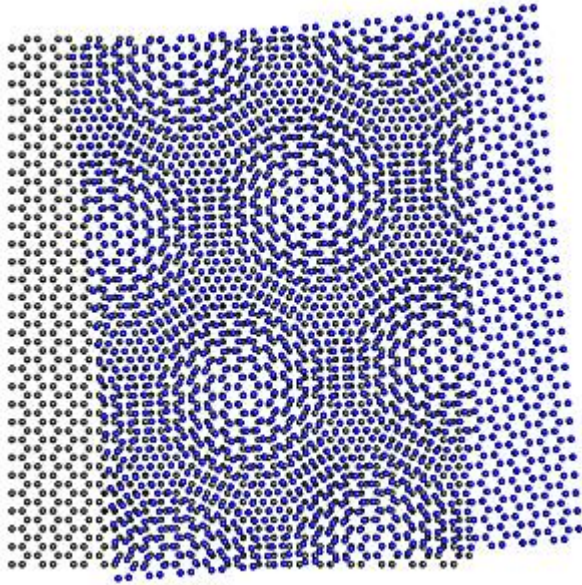
What happened here?



- The problem is weakly interacting in terms of holes!
- Emergent Fermi Liquids from an initial strongly interacting problem.
- The hole dispersion dictates the underlying physics.
- Guiding principle : Electrons prefer to occupy states with the largest hole-energy.

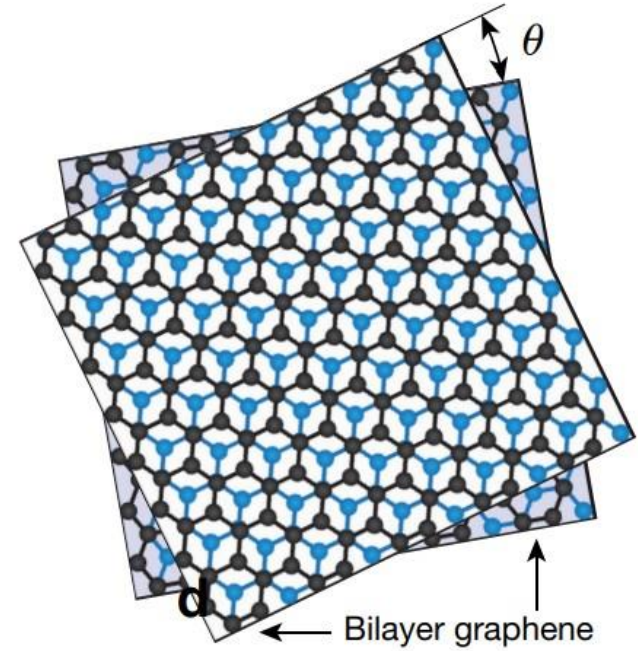
FCIs in other Moiré systems

Twisted Bilayer Graphene



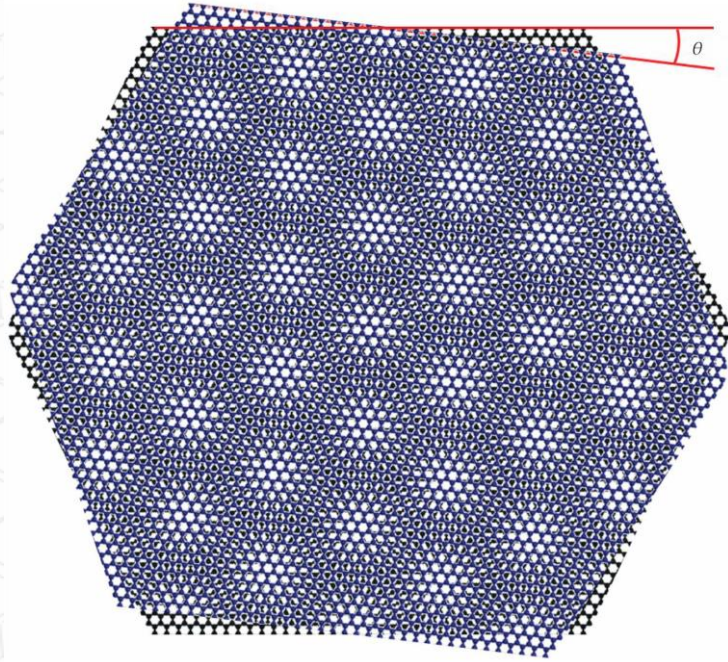
Credit : NIST

Twisted Double Bilayer Graphene



Credit : Liu et al, *Nature* **583**, (2020)

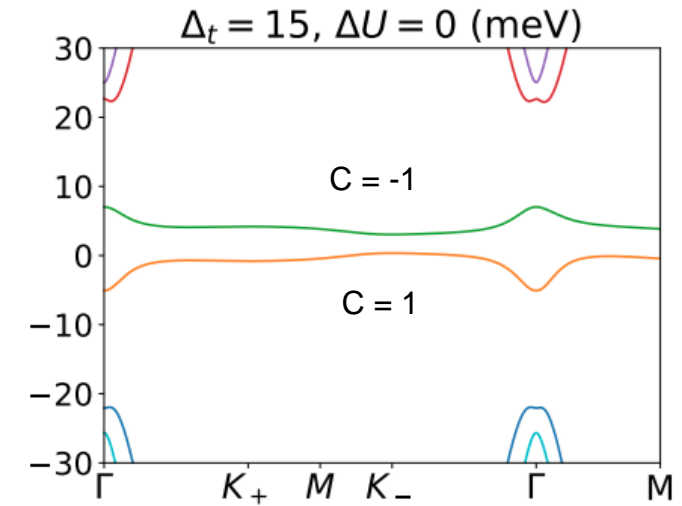
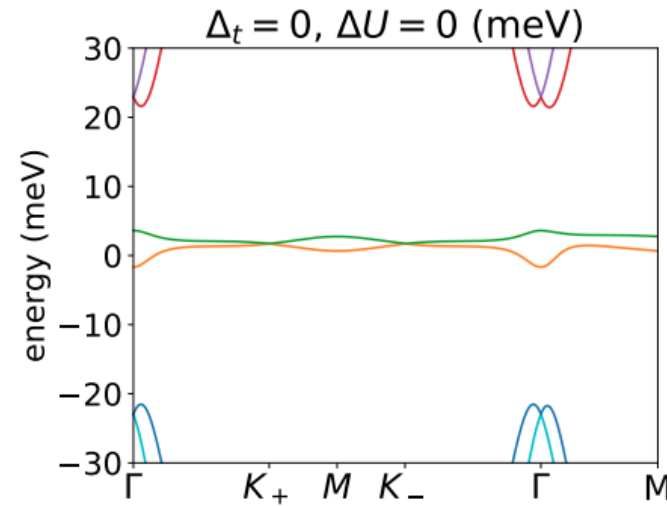
Twisted Bilayer Graphene aligned with Boron Nitride



No alignment with BN

$$\theta \approx 1.05^\circ$$

Aligned with BN



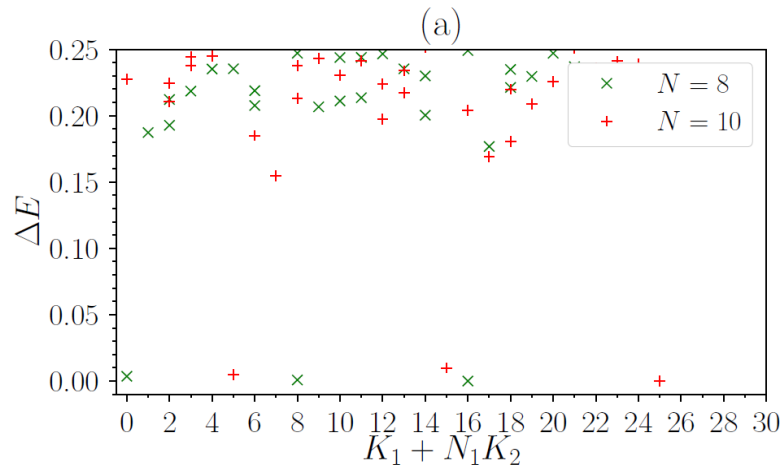
Courtesy of ICFO/ Xiaobo Lu.

arXiv:1908.00986

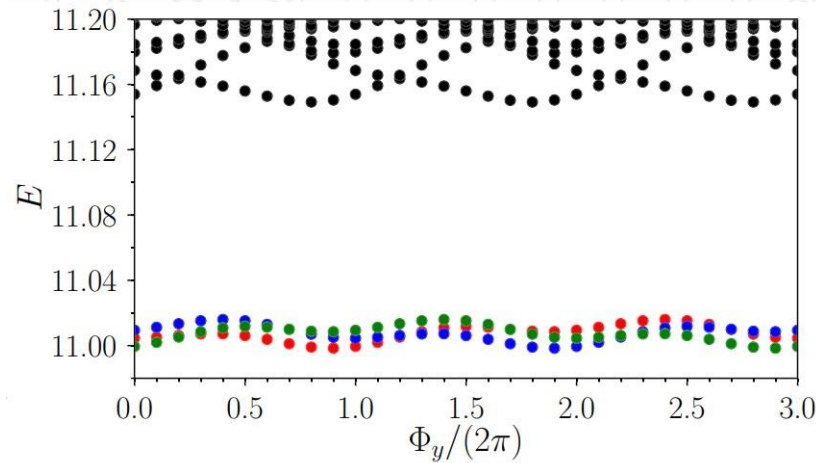
Yes, FCIs, finally!

In twisted bilayer graphene aligned with boron nitride — but only at **slightly weaker inter-layer tunnelling than in current experiments...**

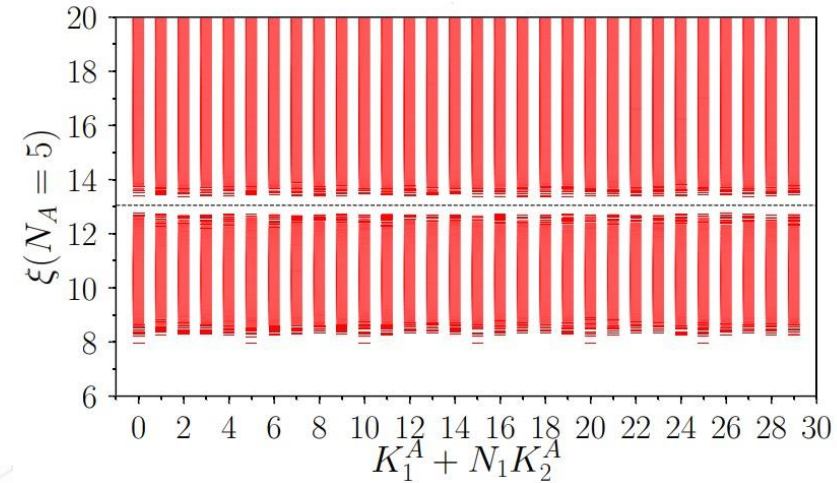
Ground state degeneracy on a torus



Spectral flow of ground states



Particle entanglement spectrum

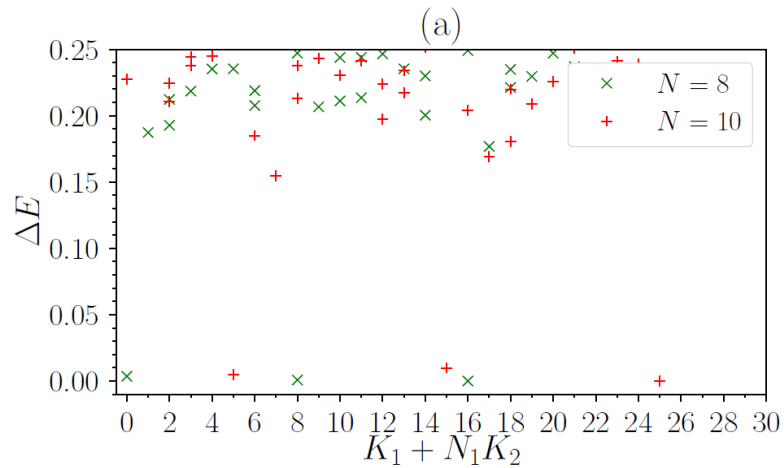


- Laughlin like state at filling $\nu = 1/3$

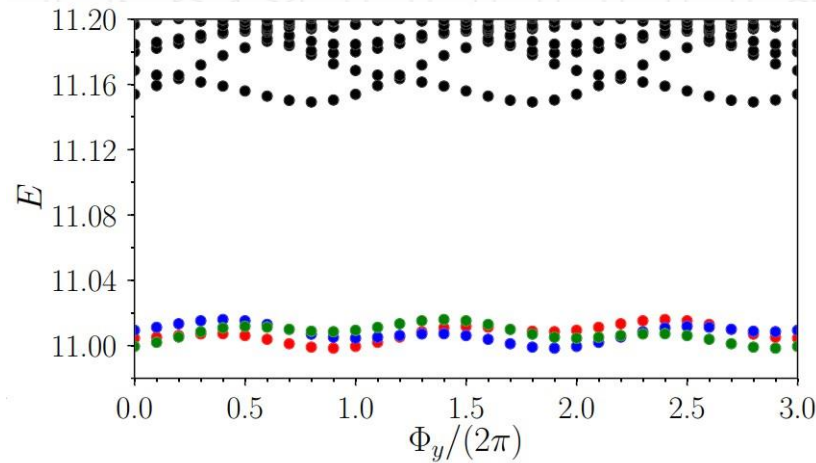
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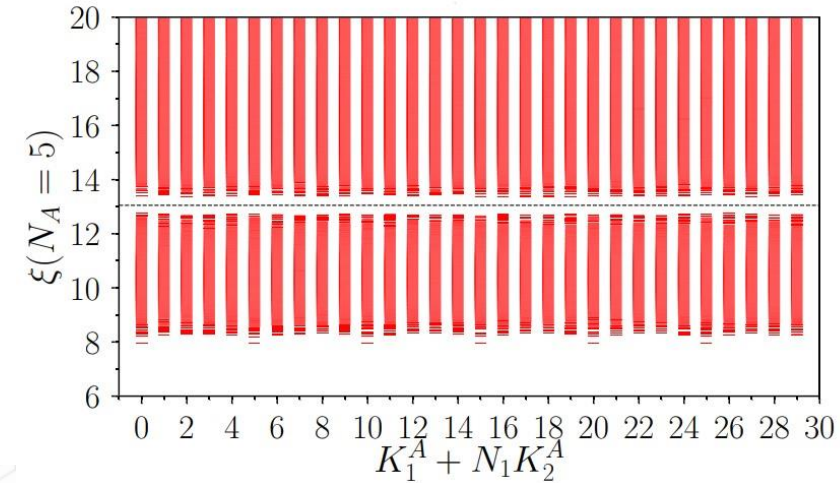
Ground state degeneracy on a torus



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Particle entanglement spectrum

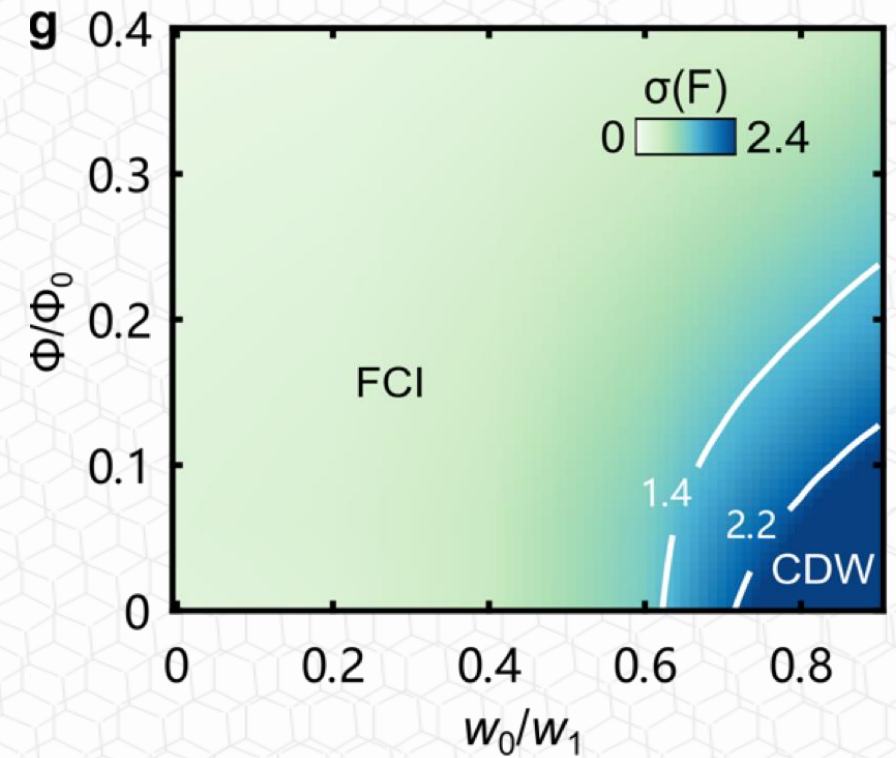
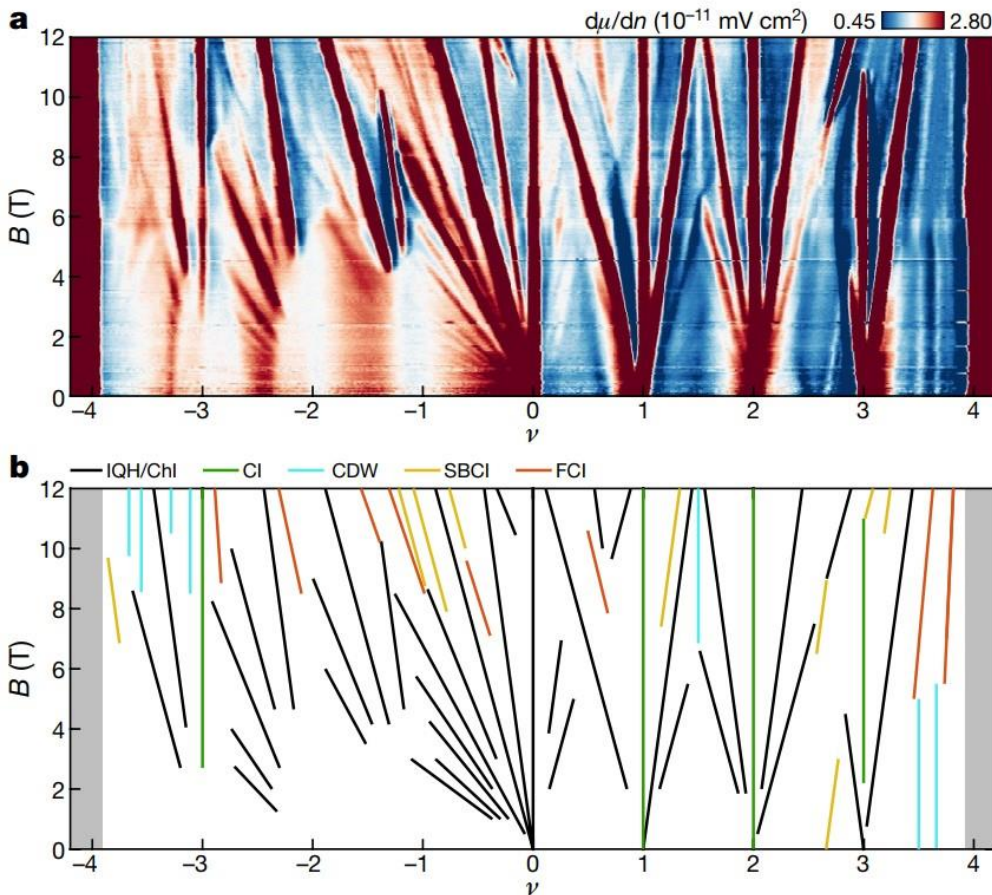


- Corroborated by other works
 - Spin polarization confirmed by Repellin and Senthil, Phys. Rev. Research 2, 023238 (2020)
 - Analytical understanding in the “chiral limit” by Ledwith et. al., Phys. Rev. Research 2, 023237 (2020)

Fractional Chern insulators in magic-angle twisted bilayer graphene

[Yonglong Xie](#) , [Andrew T. Pierce](#), [Jeong Min Park](#), [Daniel E. Parker](#), [Eslam Khalaf](#), [Patrick Ledwith](#), [Yuan Cao](#), [Seung Hwan Lee](#), [Shaowen Chen](#), [Patrick R. Forrester](#), [Kenji Watanabe](#), [Takashi Taniguchi](#), [Ashvin Vishwanath](#), [Pablo Jarillo-Herrero](#)  & [Amir Yacoby](#) 

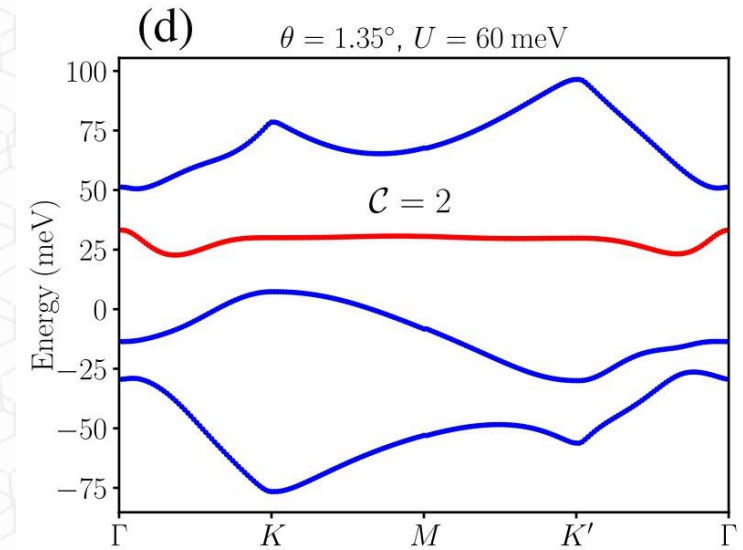
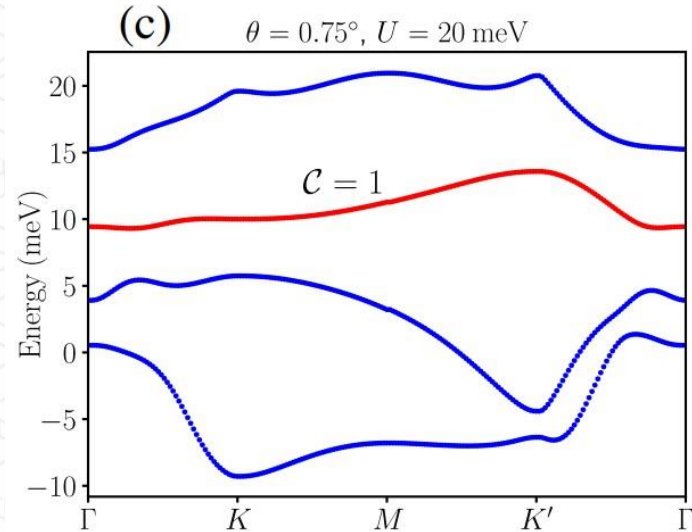
[Nature](#) **600**, 439–443 (2021) | [Cite this article](#)



- Weak field (5 Tesla), similar effect as changing inter-layer tunnelling

$$\nu = t \frac{\phi}{\phi_0} + s$$

A series of FCIs in tunable TDBG



- Spin-polarized Laughlin like state at $\nu = 1/3$
- Laughlin state particle-hole conjugate at $\nu = 2/3$
- Spin-singlet Halperin 332 state at $\nu = 2/5$
- Possibly Halperin 332 particle-hole conjugate at $\nu = 3/5$
- Spin-polarized FCI at $\nu = 1/3$ in $C = 2$ band!
- It could be thought of as a weakly interacting state of composite fermions with negative flux attachment!

Quantum Geometry : Another look at the competition

$$\eta_{\mu\nu}(\mathbf{k}) = g_{\mu\nu}(\mathbf{k}) + \frac{i}{2} \epsilon_{\mu\nu} F(\mathbf{k})$$

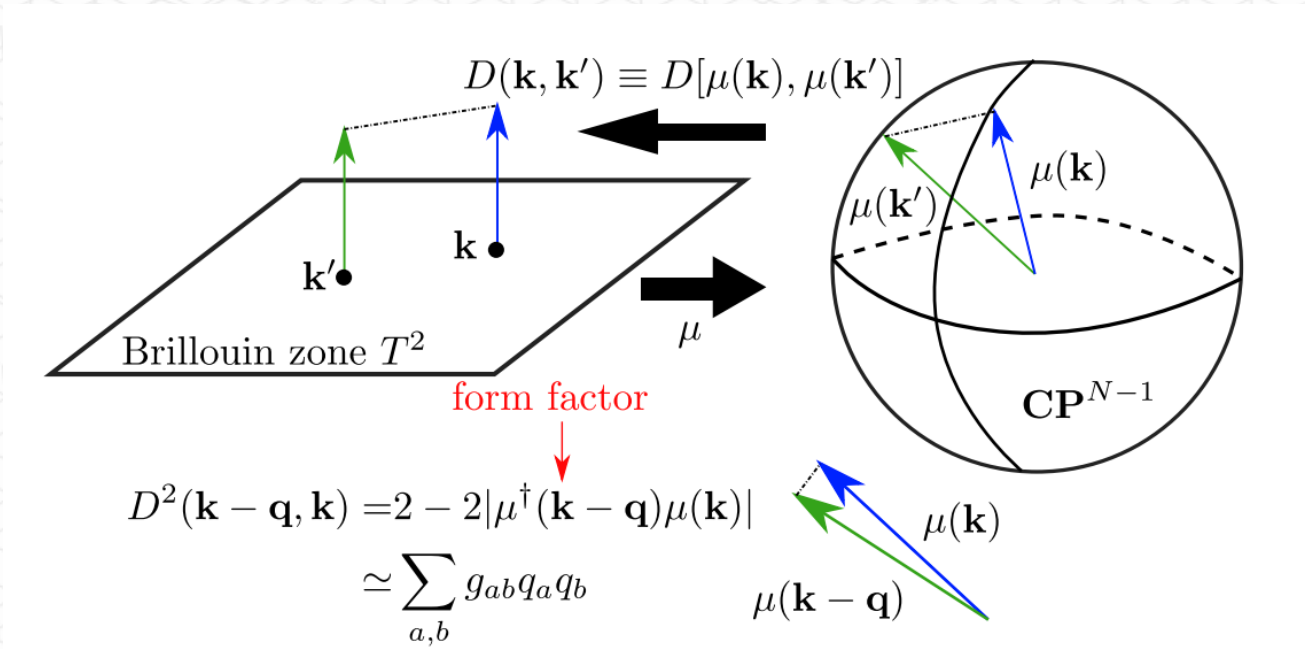
Quantum Geometric
Tensor

“Fubini-Study Metric”

Berry Curvature

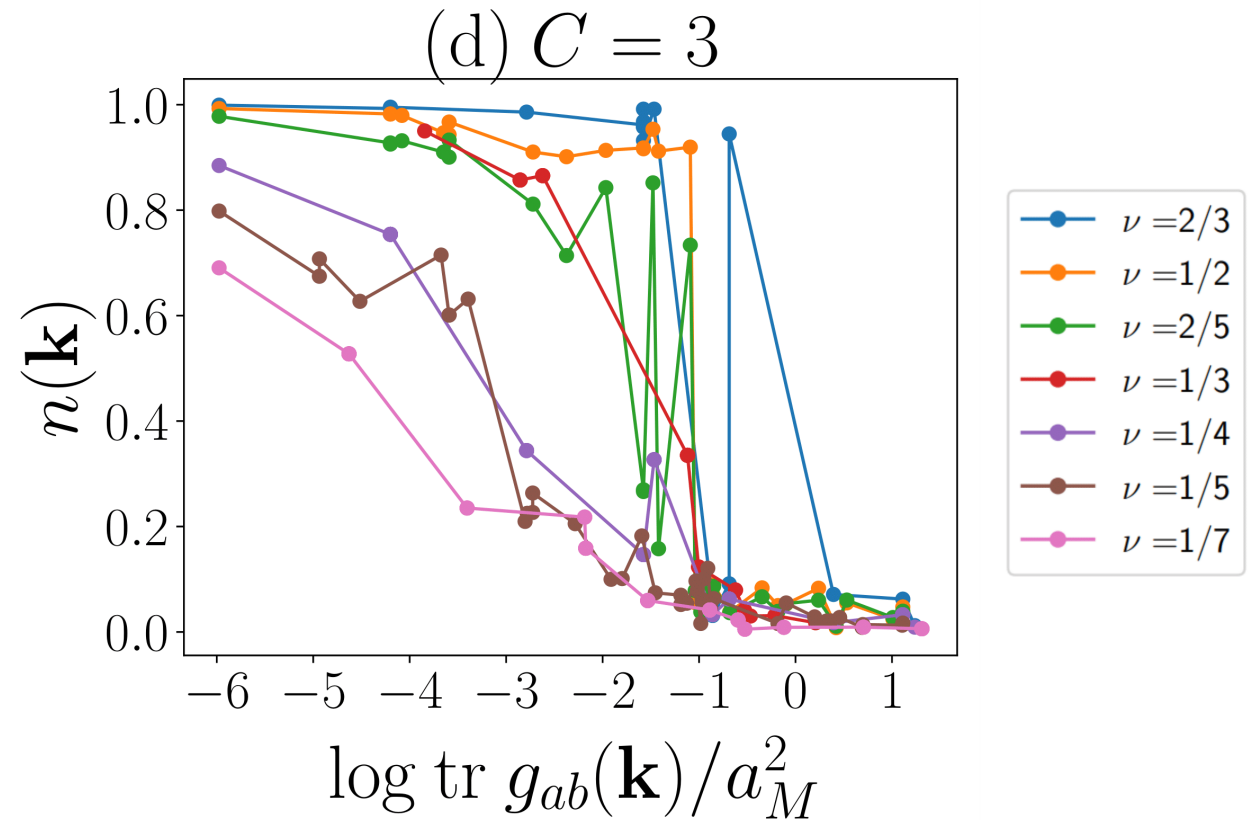
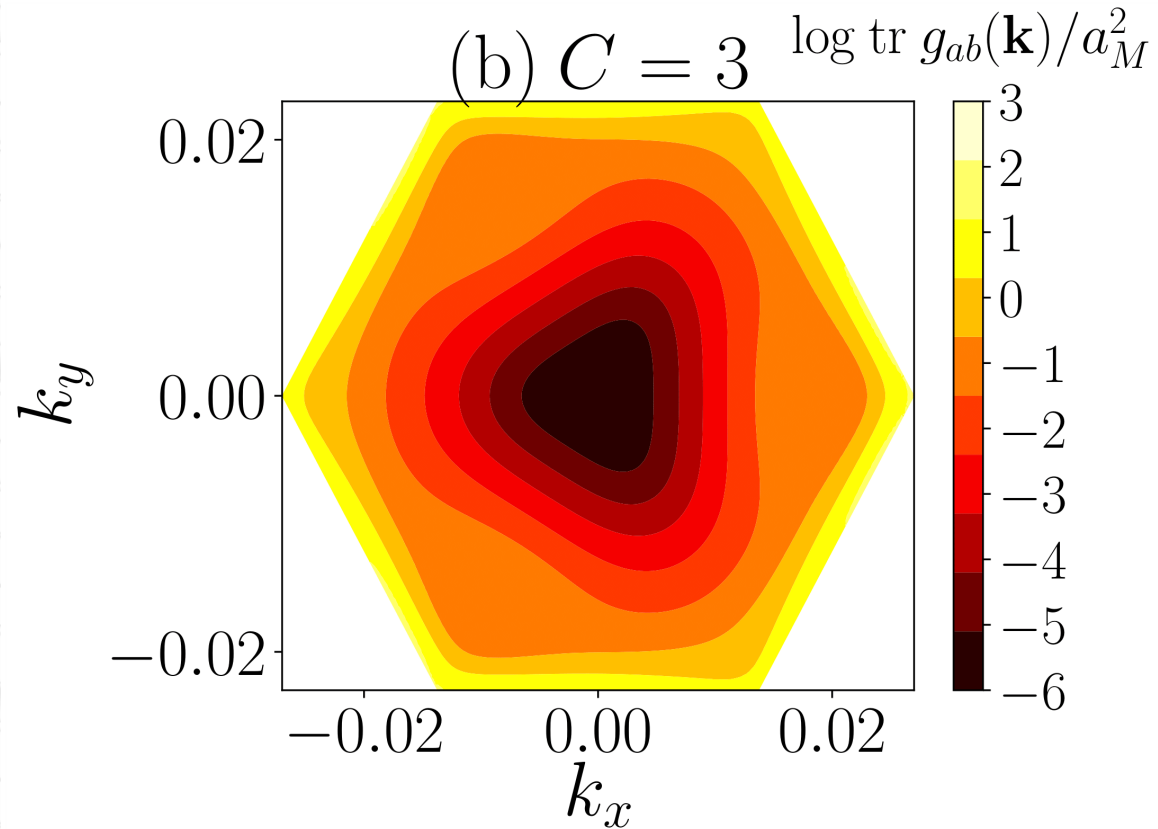
Fubini-Study metric

- Natural interpretation in terms of the distance between Bloch states



- It appears in the hole dispersion $E_h(\mathbf{k}) = \sum_{\mathbf{q}} V(\mathbf{q}) |\mu^\dagger(\mathbf{k} - \mathbf{q})\mu(\mathbf{k})|^2 \approx \sum_{\mathbf{q}} V(\mathbf{q}) (1 - q_a q_b g_{ab}(\mathbf{k}))$
- It appears also in the Fock term, $E_k(\mathbf{k}) \approx \sum_{\mathbf{q}} V(\mathbf{q}) |\mu^\dagger(\mathbf{k} - \mathbf{q})\mu(\mathbf{k})|^2 \langle c_{\mathbf{k}+\mathbf{q}}^\dagger c_{\mathbf{k}+\mathbf{q}} \rangle \approx \sum_{\mathbf{q}} V(\mathbf{q}) (1 - q_a q_b g_{ab}(\mathbf{k})) \langle c_{\mathbf{k}+\mathbf{q}}^\dagger c_{\mathbf{k}+\mathbf{q}} \rangle$
- How does the metric $g_{ab}(\mathbf{k})$ affect the electron occupation $\langle n(\mathbf{k}) \rangle = \langle c_{\mathbf{k}}^\dagger c_{\mathbf{k}} \rangle$?

Let's look again



- Guiding principle :

Electrons tend to occupy states with the highest hole-energy. \longleftrightarrow Electrons tend to occupy states with the lowest Fubini-Study metric trace!

Take home message

- Moiré systems are promising platforms for fractional Chern insulators.
- The particle-hole asymmetry of interactions in a single band has dramatic consequences.
- The Fubini-Study metric is a relevant quantity to the low energy physics of Moiré materials.