Charge Order in NbSe₂

Felix Flicker

U. C. Berkeley

R The English-Speaking Union





- Introduction 1D CDWs
 Weak Coupling NbSe₂
 - Strong Coupling

- gap - pseudogap

- Introduction 1D CDWs
 2D CDWs
 Weak Coupling NbSe₂
 - Strong Coupling

- gap - pseudogap

- 1D CDWs Introduction - 2D CDWs - NbSe₂ Weak Coupling - gap - pseudogap Strong Coupling Conclusions

- 1D CDWs Introduction - 2D CDWs - NbSe₂ Weak Coupling - gap - pseudogap Strong Coupling
- Conclusions



Prototype layered CDW Material

- CDW gap in 1 band only
- wide 'Kohn anomaly'
- particle/hole gap asymmetry
- gap offset
- pseudogap



Prototype layered CDW Material

- CDW gap in 1 band only
- wide 'Kohn anomaly'
- particle/hole gap asymmetry
- gap offset
- pseudogap



Prototype layered CDW Material

- CDW gap in 1 band only
- wide 'Kohn anomaly'
- particle/hole gap asymmetry
- gap offset
- pseudogap



Prototype layered CDW Material

- CDW gap in 1 band only
- wide 'Kohn anomaly'
- particle/hole gap asymmetry
- gap offset
- pseudogap



Prototype layered CDW Material

- CDW gap in 1 band only
- wide 'Kohn anomaly'
- particle/hole gap asymmetry
- gap offset
- pseudogap



Prototype layered CDW Material

- CDW gap in 1 band only
- wide 'Kohn anomaly'
- particle/hole gap asymmetry
- gap offset
- pseudogap



A. Soumyanarayanan et al. PNAS 110(5):1623-1627 (2013)



D. J. Rahn et al. PRB 85, 224532 (2012)



Weak Coupling



F. Weber et al. PRL 107, 107403 (2011)





$$H|A_{\lambda}\rangle = E_{\lambda}S|A_{\lambda}\rangle$$

$$\mathbf{g}_{\mathbf{k},\mathbf{k}'}^{\lambda,\lambda'} \propto \mathbf{v}_{\mathbf{k}}^{\lambda} \left[A_{\mathbf{k}}^{\dagger} S_{\mathbf{k}} A_{\mathbf{k}'} \right]^{\lambda\lambda'} - \left[A_{\mathbf{k}}^{\dagger} S_{\mathbf{k}'} A_{\mathbf{k}'} \right]^{\lambda\lambda'} \mathbf{v}_{\mathbf{k}'}^{\lambda'}$$

C. M. Varma et al. PRB 19, 6130-6141 (1978)







Borisenko et al., PRL 102, 166402 (2002)









- Peierls Mechanism not the natural starting point in D>1
- $g_{kk'}^{\lambda\lambda'}$ needed
- First consistent theoretical explanation of NbSe₂ experimental results

- Peierls Mechanism not the natural starting point in D>1
- $g_{kk'}^{\lambda\lambda'}$ needed
- First consistent theoretical explanation of NbSe₂ experimental results

- Peierls Mechanism not the natural starting point in D>1
- $g_{kk'}^{\lambda\lambda'}$ needed
- First consistent theoretical explanation of NbSe₂ experimental results

- Peierls Mechanism not the natural starting point in D>1
- $g_{kk'}^{\lambda\lambda'}$ needed
- First consistent theoretical explanation of NbSe₂ experimental results

Thanks For Listening

F. Flicker and Jasper van Wezel, "*Charge Order from Orbital Dependent Coupling Evidenced by* NbSe₂", **Nature Communications 6, 7034 (2015)**

Y. Feng, J. van Wezel, J. Wang, F. Flicker, D. M. Silevitch, P. B. Littlewood, and T. F. Rosenbaum, "*Quantum Criticality and the Charge Density Wave Phase Transition in* NbSe₂", **Nature Physics 11, 865 (2015)**

F. Flicker and Jasper van Wezel, "*Charge Ordering Geometries in Uniaxially Strained* NbSe₂", **Physical Review B 92, 201103 (Rapid Communications) (2015)**

F. Flicker and Jasper van Wezel, "Charge Order in Monolayer NbSe2", in preparation