#### Details in GW-BSE

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#### What's GW

To be very concise...  $-i\Sigma \approx GW^{RPA} =$ 

Assuming quasiparticles...  $v(q+G) = \frac{4\pi}{V|q+G|^2}$ .  $\Sigma^{\int_{\text{contour}}} \Sigma^{\text{COH}} + \Sigma^{\text{SEX}}$ ,  $\epsilon_{GG'}(q,\omega) = \delta_{GG'} - v(q+G)\chi_{GG'}(q,\omega)$ ,  $M_{nn'}(k,q,G) = \langle nk+q | \mathrm{e}^{\mathrm{i}(q+G)\cdot r} | n'k \rangle$ 

$$\langle n\mathbf{k} | \Sigma^{\text{SEX}}(\omega) | n'\mathbf{k} \rangle = -\sum_{n''}^{\text{occ}} \sum_{\mathbf{q} \mathbf{G} \mathbf{G}'} M_{n''n}^*(\mathbf{k}, -\mathbf{q}, -\mathbf{G}) M_{n''n'}(\mathbf{k}, -\mathbf{q}, -\mathbf{G}')$$

$$\times \epsilon_{\mathbf{G} \mathbf{G}'}^{-1}(\mathbf{q}, \omega - E_{n'', \mathbf{k} - \mathbf{q}}) v(\mathbf{q} + \mathbf{G}'),$$
(1)

$$\langle n\mathbf{k} | \Sigma^{\text{COH}}(\omega) | n'\mathbf{k} \rangle = \frac{\mathrm{i}}{2\pi} \sum_{n''} \sum_{\mathbf{q}, \mathbf{G}, \mathbf{G}'} M_{n''n}^*(\mathbf{k}, -\mathbf{q}, -\mathbf{G}) M_{n''n'}(\mathbf{k}, -\mathbf{q}, -\mathbf{G}')$$

$$\times \int_0^\infty \mathrm{d}\omega' \frac{[\epsilon_{\mathbf{G}\mathbf{G}'}^{\mathbf{G}}]^{-1}(\mathbf{q}, \omega') - [\epsilon_{\mathbf{G}\mathbf{G}'}^{\mathbf{a}}]^{-1}(\mathbf{q}, \omega')}{\omega - E_{n\mathbf{k}} - \omega' + \mathrm{i} \, 0^+ \, \mathrm{sgn}(E_{n\mathbf{k}})} v(\mathbf{q} + \mathbf{G}').$$
(2)

$$\chi_{\boldsymbol{G}\boldsymbol{G}'}^{r/a}(\boldsymbol{q},\omega) = \sum_{\boldsymbol{k}} \sum_{n} \sum_{n'}^{\text{emp}} M_{nn'}(\boldsymbol{k},\boldsymbol{q},\boldsymbol{G}) M_{nn'}^{*}(\boldsymbol{k},\boldsymbol{q},\boldsymbol{G}') \times \left( \frac{1}{(1+\boldsymbol{F}_{n}) + \boldsymbol{F}_{n} + 10^{+}} + \frac{1}{(1+\boldsymbol{F}_{n}) + 10^{+}} \right).$$
(3)

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#### What's GW

What a splendid equation system!

In practice. . . we always use GPP  $\epsilon(\omega)$  is assumed to be plasmon model-like, so – we feed

$$\chi_{\mathbf{G}\mathbf{G}'}(\mathbf{q},\omega) = \sum_{\mathbf{k}} \sum_{n}^{\text{occ}} \sum_{n'}^{\text{emp}} M_{nn'}(\mathbf{k},\mathbf{q},\mathbf{G}) M_{nn'}^{*}(\mathbf{k},\mathbf{q},\mathbf{G}') \frac{2}{E_{n,\mathbf{k}+\mathbf{q}} - E_{n'\mathbf{k}}}$$
(4)

to analytic expressions of  $\Sigma^{COH, SEX}$ .

**Huge simplification...** Hedin  $\overset{\text{assuming }GW}{\to} GW \overset{\text{QP. approx.}}{\to} (1), (2), (3)$   $\overset{\text{GPP}}{\to}$  we are here

**But still burdensome** *Summation over empty bands – 1000-30000 bands!!!* 

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## The problem of summing over many empty bands

Empty states are important - but why?

One further simplification trick: pseudobands. For  $E_{nk} \gg E_F$ :

$$\begin{split} \{\phi_{n\pmb{k}}\}_{\text{adjacent energy block }n_b} &\to \sum_{n\in \text{block }n_b} \phi_{n\pmb{k}}, \\ \{E_{n\pmb{k}}\}_{\text{adjacent energy block }n_b} &\to \frac{1}{|n_b|} \sum_{n\in \text{block }n_b} E_{n\pmb{k}}. \end{split}$$

4000 bands  $\rightarrow \sim$  400 to  $\sim$  1000 bands.

...but is this snake oil? It makes no sense!!!

# Understanding pseudobands: energy averaging

#### Observation

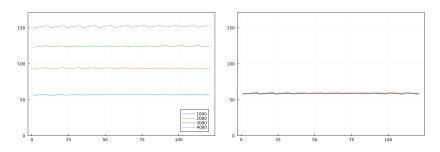
- $\chi \sim MM^* \times \text{some-function}(E_{\text{emp}} E_{\text{occ}});$
- $\Sigma^{\text{COH}} \sim MM^* \times \text{some-function}(\omega E)$ ;
- We are only interested in  $\omega \sim E_{\rm F}$ .
- **★**For  $E_{n\mathbf{k}}\gg E_{\mathsf{F}}$  ( $\omega$ ): energy-dependent factors  $\sim$  const. for all  $\mathbf{k}$ .

Thus for both  $\chi$  and  $\Sigma^{COH}$  involving summation over empty bands:

$$\sum_{n''} M_{n''n}^* M_{n''n'} \times \cdots = \sum_{\text{block } n_b} \cdots \times \sum_{n'' \in \text{block } n_b} M_{n''n}^* M_{n''n'}. \quad (5)$$

In RHS  $E_{\text{emp}}$  is replaced by the average energy.

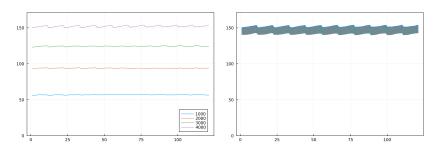
# Understanding pseudobands: energy averaging



**Example** High-energy DFT bands of WTe<sub>2</sub> monolayer (ONCV-SG15, 120 electrons; 80 Ry cutoff,  $20 \times 20 \times 1$  grid). x axis = k index in irreducible 1BZ; fastest varying coordinate =  $k_y$ 

- **■** Bands above 1000 are generally flat (but dispersion  $\uparrow$  as  $n \uparrow$ )
- Low-lying pseudobands blocks are very close in energy (but  $1/(E_{\rm emp}-E_{\rm occ})$  more sensitive to dispersion)

# Understanding pseudobands: energy averaging



**Example** High-energy DFT bands of WTe<sub>2</sub> monolayer (ONCV-SG15, 120 electrons; 80 Ry cutoff,  $20 \times 20 \times 1$  grid). x axis = k index in irreducible 1BZ; fastest varying coordinate =  $k_y$ 

- **■** Bands above 1000 are generally flat (but dispersion  $\uparrow$  as  $n \uparrow$ )
- High-energy blocks are more dispersive (but  $1/(E_{\rm emp}-E_{\rm occ})$  is smaller so no worry)

## Understanding pseudobands: wave function averaging?

For pseudobands to work, we need

$$\sum_{n'' \in \text{block } n_b} M_{n''n}^* M_{n''n'} \sim M_{\text{averaged band},n}^* M_{\text{averaged band},n'}$$

$$= \sum_{n_1'', n_2'' \in \text{block } n_b} M_{n_1''n'} M_{n_2''n}^*. \tag{6}$$

The problem: it of course isn't the case in general. If  $[M_{n''_1n'}M^*_{n''_2n}]_{n''_1n''_2}$  is a random matrix: LHS: RHS  $\approx 0.3$ .

**The question:** Then in which case is (6) correct in some sense?

### The structure of $\phi_{n\mathbf{k}}$

Plane wave basis In BerkeleyGW WFN.h5,

$$\phi_{nk}(\mathbf{r},\sigma) = \frac{1}{\sqrt{V}} \sum_{\mathbf{G}} e^{i(\mathbf{k}+\mathbf{G})\cdot\mathbf{r}} c_{n\mathbf{k},\mathbf{G}\sigma}.$$
 (7)

Thus

$$M_{nn'}(\boldsymbol{k}, \boldsymbol{q}, \boldsymbol{G}) = \langle n\boldsymbol{k} + \boldsymbol{q} | e^{i(\boldsymbol{q} + \boldsymbol{G}) \cdot \boldsymbol{r}} | n' \boldsymbol{k} \rangle = \sum_{\boldsymbol{G}', \sigma} c_{n\boldsymbol{k} + \boldsymbol{q}, \boldsymbol{G} + \boldsymbol{G}' \sigma}^* c_{n\boldsymbol{k}, \boldsymbol{G}' \sigma}.$$
 (8)

**Cutoff** Each k has its own G-grid ( $\sim 30000$  vectors for 80 Ry).

#### Procedure

#### Input

- indices of k, q in k-grid;
- index of  $\boldsymbol{G}$  in  $\boldsymbol{G}$ -grid of  $\boldsymbol{k}$  (expect a  $\boldsymbol{G}$  in GW  $\boldsymbol{G}$ -grid, cutoff = say 30 Ry, not 80 Ry);
- $\bullet$  n, n'.

#### **Procedure**

- $\bigcirc$  find index of k
- ② find the index of  ${m G}+{m G}'$  in  ${m G}$ -grid of  ${m k}+{m q}$ , for each  ${m G}'$  in  ${m G}$ -grid of  ${m k}$
- **3** do summation  $\sum_{\mathbf{G}',\sigma} c_{n\mathbf{k}+\mathbf{q},\mathbf{G}+\mathbf{G}'\sigma}^* c_{n\mathbf{k},\mathbf{G}'\sigma}$ .

**Performance** Main bottleneck: finding G + G'. Using StaticArrays.jl helps a lot!

Under GPP:

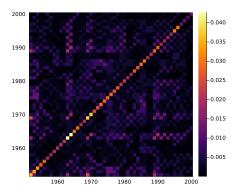
$$\chi_{\boldsymbol{G}\boldsymbol{G}'}(\boldsymbol{q},\omega)^{\text{high band terms}} \approx \sum_{\boldsymbol{k}} \sum_{\text{block } n_{b}}^{\text{emp}} \frac{2}{E_{n,\boldsymbol{k}+\boldsymbol{q}} - E_{\text{average in block } n_{b}}} \times \sum_{n' \in \text{block } n_{b}} \sum_{n}^{\text{occ}} M_{nn'}(\boldsymbol{k},\boldsymbol{q},\boldsymbol{G}) M_{nn'}^{*}(\boldsymbol{k},\boldsymbol{q},\boldsymbol{G}')$$

$$(9)$$

Our goal Finding how diagonal is

$$\sum_{n}^{\text{occ}} M_{nn'}(\boldsymbol{k}, \boldsymbol{q}, \boldsymbol{G}) M_{nn'}^*(\boldsymbol{k}, \boldsymbol{q}, \boldsymbol{G}')$$
 (10)

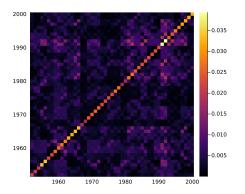
**Case 1:** G = G' In this case  $\sum_{n}^{\text{occ}} M_{nn'}(\mathbf{k}, \mathbf{q}, G) M_{nn'}^*(\mathbf{k}, \mathbf{q}, G')$  is large and fairly diagonal



Parameters  $\mathbf{G} = (0, 1, -14), \mathbf{k} = \mathbf{k}_2 = (0, 0.00, 0), \mathbf{q} = (0, 0.00, 0)$ 

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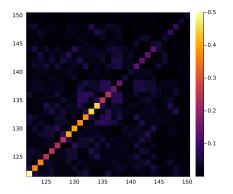
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Parameters  $\mathbf{G} = (0, 1, -14), \mathbf{k} = \mathbf{k}_2 = (0, 0.05, 0), \mathbf{q} = (0, 0.10, 0)$ 

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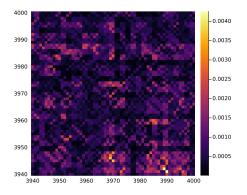
**Case 1:** G = G' In this case  $\sum_{n}^{\text{occ}} M_{nn'}(\mathbf{k}, \mathbf{q}, G) M_{nn'}^*(\mathbf{k}, \mathbf{q}, G')$  is large and fairly diagonal



Strikingly,  $\sum_{n}^{\text{occ}} M_{nn'}(\mathbf{k}, \mathbf{q}, \mathbf{G}) M_{nn'}^*(\mathbf{k}, \mathbf{q}, \mathbf{G}')$  is still very diagonal for bands near Fermi surface!

## Numerical charaterization of $\sum_{n}^{\text{occ}} M_{nn'}(\mathbf{k}, \mathbf{q}, \mathbf{G}) M_{nn'}^*(\mathbf{k}, \mathbf{q}, \mathbf{G}')$

Case 2:  $G \neq G' \sum_{n}^{\text{occ}} M_{nn'}(\mathbf{k}, \mathbf{q}, G) M_{nn'}^*(\mathbf{k}, \mathbf{q}, G')$  is very non-diagonal, but since the terms's random phases cancel each other so it's small



# Half-way generalization about pseudobands in $\chi$

- pseudobands works when it's necessary to do so
- What prevents pseudobands from working around Fermi surface is the energy dispersion

... but do we have any theoretical explanation for this?