Constrained RPA

Jinyuan Wu

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Motivation

Strongly correlated electrons are ...

- Hard to treat *ab initio*
- Should be described by a lattice model

$$H = -t \sum_{\langle \boldsymbol{i},\boldsymbol{j} \rangle,\sigma} c^{\dagger}_{\boldsymbol{i}\sigma} c_{\boldsymbol{j}\sigma} + \text{h.c.} + U \sum_{\boldsymbol{i}} n_{\boldsymbol{i}\uparrow} n_{\boldsymbol{i}\downarrow}. \tag{1}$$

But should all electrons be included?

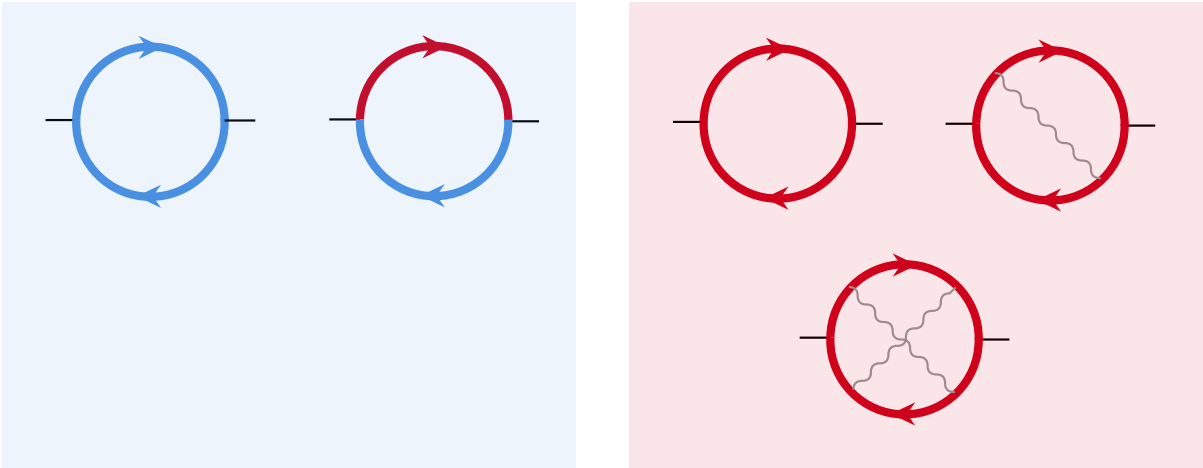
- d electron: correlated ones
- r electron: "trivial" ones

Diagrammatics

Screening by d and r = first screened by r and then d

$$W = (1 - vP)^{-1}v = (1 - W_r P_d)^{-1} W_r, \quad W_r = (1 - vP_r)^{-1}v, \qquad (2)$$

- W is effective interaction
- where P_d is the polarization within d subspace,
- and $P_r = P - P_d$

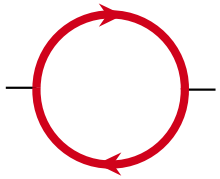


P_r
RPA works

P_d
RPA doesn't work

Diagrammatics

- P_d hard to obtain $\Rightarrow$ we left it to an effective interaction without doing resummation for it
- P_r can be done by `epsilon` $\Rightarrow$ W_r can be obtained by `epsilon`
- But



  is not in P_r, or otherwise we have double counting
- That's what we call **constrained RPA (cRPA)**
- A program to calculate a single ring diagram is needed

Band part

In GW:

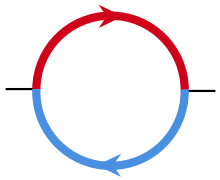
$$\begin{aligned} \varepsilon^{GW} &= \varepsilon^{0} + \Sigma^{\text{Hartree}} + \Sigma^{GW} \\ &= \underbrace{\varepsilon^{0} + \Sigma^{\text{Hartree}} + V^{\text{xc}}}_{\texttt{WFN}} + \underbrace{\Sigma^{GW}}_{\texttt{sigma}} - \underbrace{V^{\text{xc}}}_{\texttt{vxc.dat}} \end{aligned} \tag{3}$$

In the effective model ...

- We want ε^{0}
- So the same $-V^{\text{xc}}$ procedure is needed
- A problem to calculate a single ring diagram is needed.

Risks in accuracy

Uncontrolled approximation: does RPA work for



The hopping between d electrons is small $\Rightarrow$ it's safe to do so[1]

[1] F Aryasetiawan et al. "Frequency-dependent local interactions and low-energy effective models from electronic structure calculations". In: *Physical Review B* 70.19 (2004). Fig. 3 and Fig. 4 in the article are swapped in the published version: compare them with Fig 2, p. 195104.

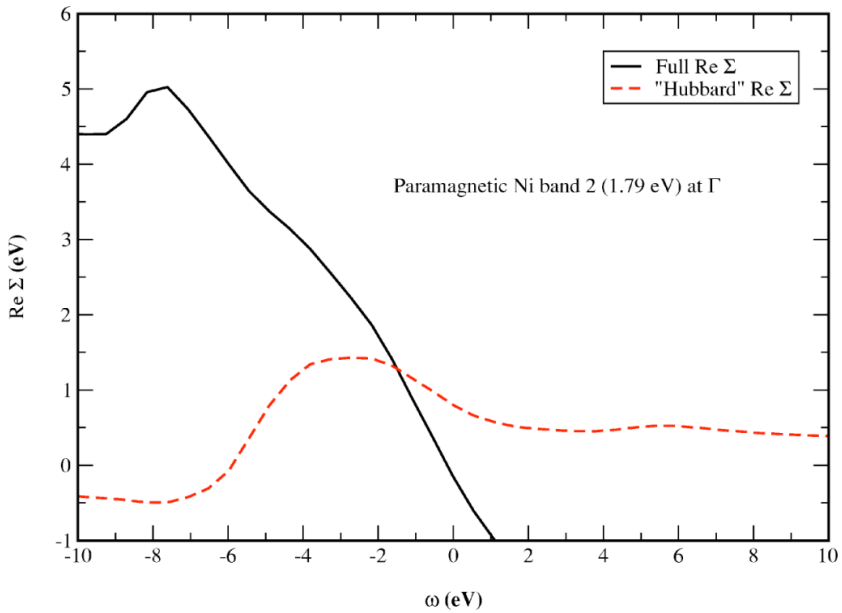
Dynamic interaction in Hubbard model

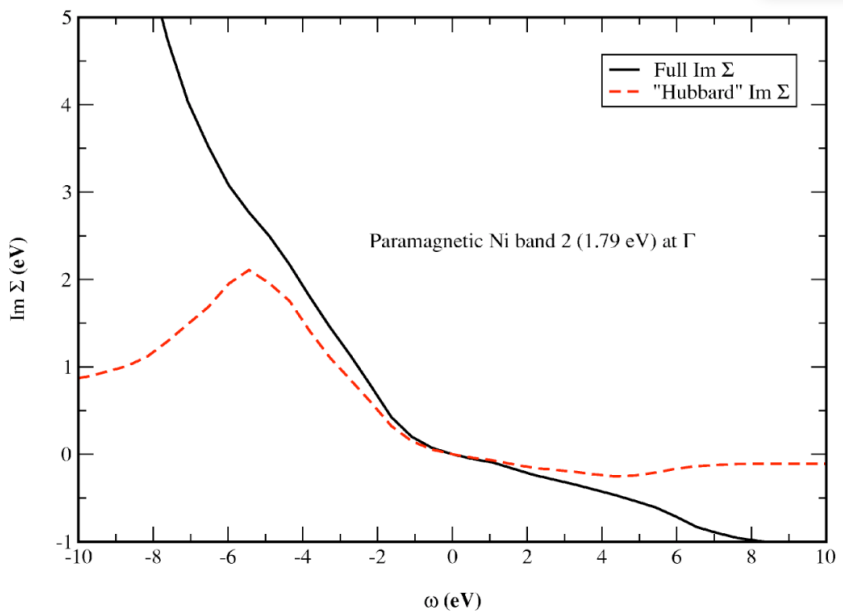
Recall that $W = W(\boldsymbol{r}, \boldsymbol{r}, \omega) \Rightarrow U_{\boldsymbol{ijkl}}(\omega) \Rightarrow$ retarded interaction!

- Hamiltonian form: interactions are always immediate (bosonic auxiliary field required to create retardation)
- . . . or path integral formalism is to be used
- Either it, it's slow!!!

Dynamic interaction in Hubbard model

Is it possible to just enforce $\omega = 0$ . . . But it's not accurate![2]





Ni self-energy, from static Hubbard $U = W_r(\omega = 0)$, and from full $W(\omega)$

[2] Aryasetiawan et al., "Frequency-dependent local interactions and low-energy effective models from electronic structure calculations".

Dynamic interaction in Hubbard model

Why $\mathrm{Im}\,\Sigma$ is better captured by static U than $\mathrm{Re}\,\Sigma$?
Because $\mathrm{Im}\,\Sigma \propto$ DOS TODO: why??
And then high-frequency electrons contribute to low-frequency $\mathrm{Re}\,\Sigma$

Possible way to make the Hubbard model static

Key point: correct the single-electron Hamiltonian using the retarded interaction

- Def: G_d = Green function for d electrons; $\tilde{G}_d$ = Green function for d electrons corrected by $W - W_d$; W = RPA screened Coulomb interaction; W_d = static Hubbard U screened by G_d; $\tilde{W}_d$ = static Hubbard U screened by $\tilde{G}_d$;
- Ignoring hopping between d and r subspaces caused by W (so the $\mathrm{i}\, W(G - G_d)$ term is irrelevant for d electrons), we have

$$\tilde{G}_d^{-1} - \mathrm{i}\, \tilde{G}_d \tilde{W}_d = G_d^{-1} - \mathrm{i}\, G_d W_d - \mathrm{i}\, G_d (W - W_d). \tag{4}$$

- From

$$W = (1 - UP)^{-1} U, \quad P = -\mathrm{i}\, GG \Rightarrow G^{-1} - \mathrm{i}\, GW = (1 - UP)^{-1} G^{-1} \tag{5}$$

we get the final equation: (here $\tilde{P}_d = -\mathrm{i}\, \tilde{G}_d \tilde{G}_d$, $P_d = -\mathrm{i}\, G_d G_d$)

$$(1 - U\tilde{P}_d)^{-1} \tilde{G}_d^{-1} = (1 - UP_d^{-1})^{-1} G_d^{-1} - \mathrm{i}\, G_d (W - W_d). \tag{6}$$

Existing implementations

- **VASP** `ALGO=CRPA` selects constrained RPA calculations, available as of VASP.6.4
- **ABINIT** `ucrpa` before version 9, as an option in the RPA module
- **BerkeleyGW** ✗Unfortunately . . .