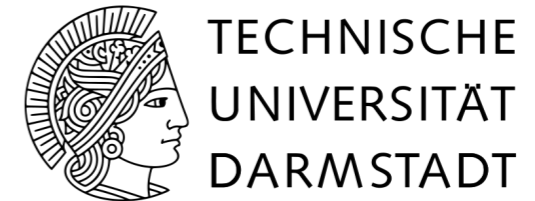


# Gorkov-Green's function approach to open-shell nuclei

Vittorio Somà (EMMI/TU Darmstadt)



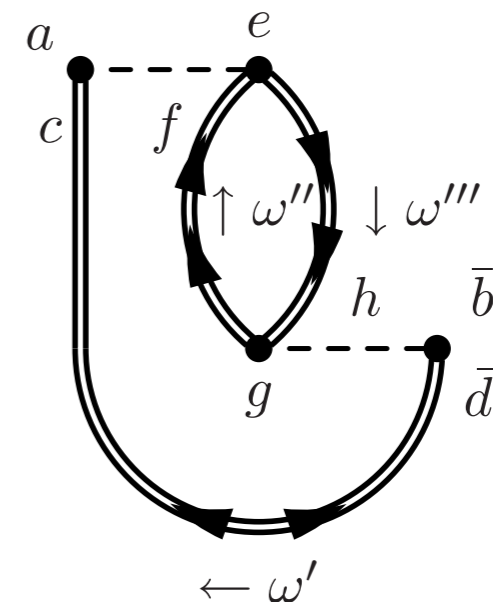
*Collaborators:*

Carlo Barbieri (Surrey, UK)

Thomas Duguet (CEA Saclay, France)

*Based on:*

VS, Duguet, Barbieri PRC 84 064317 (2011)

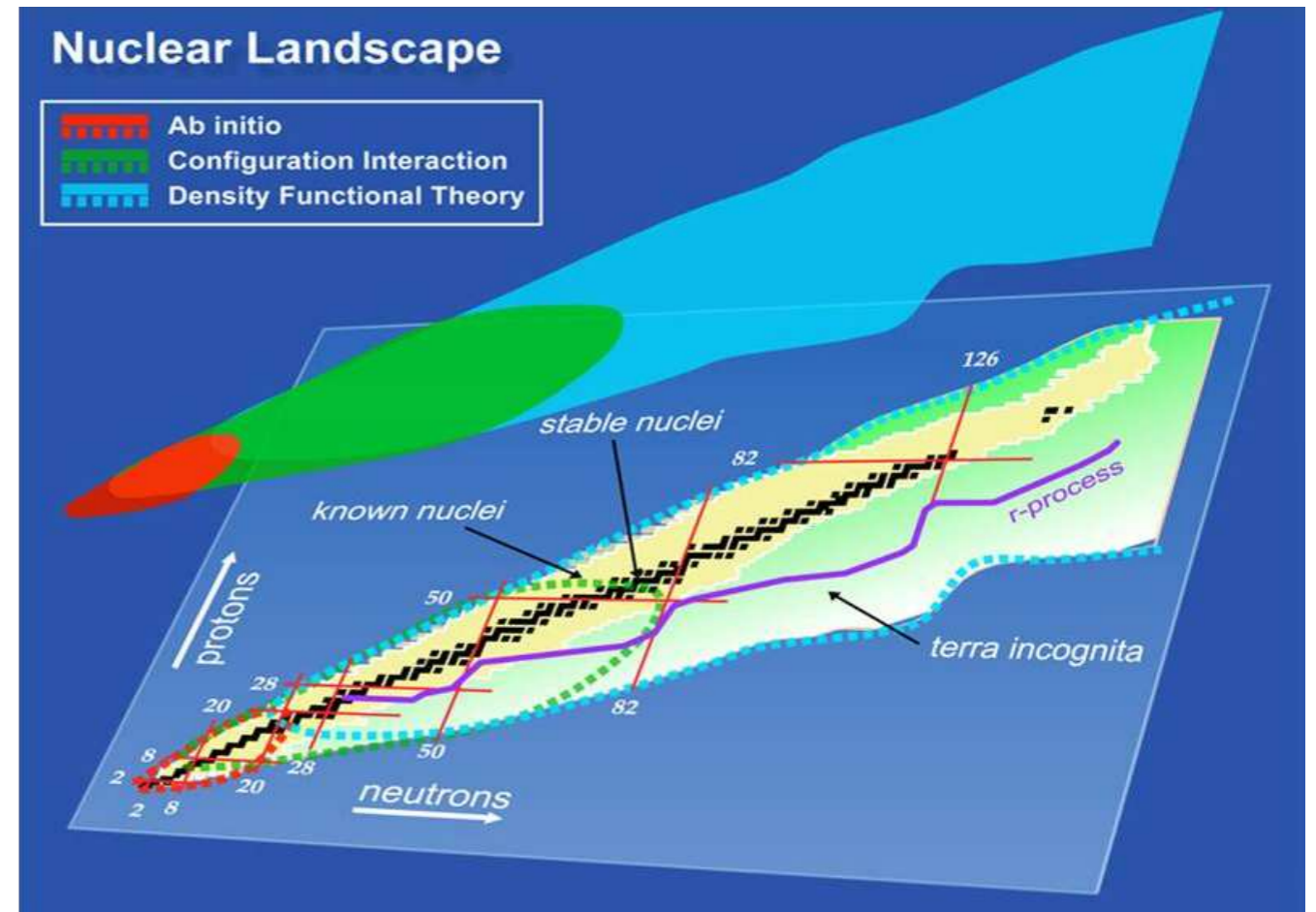


“Facets of Strong-Interaction Physics”

Hirscheegg 2012

# Towards a unified description of nuclei

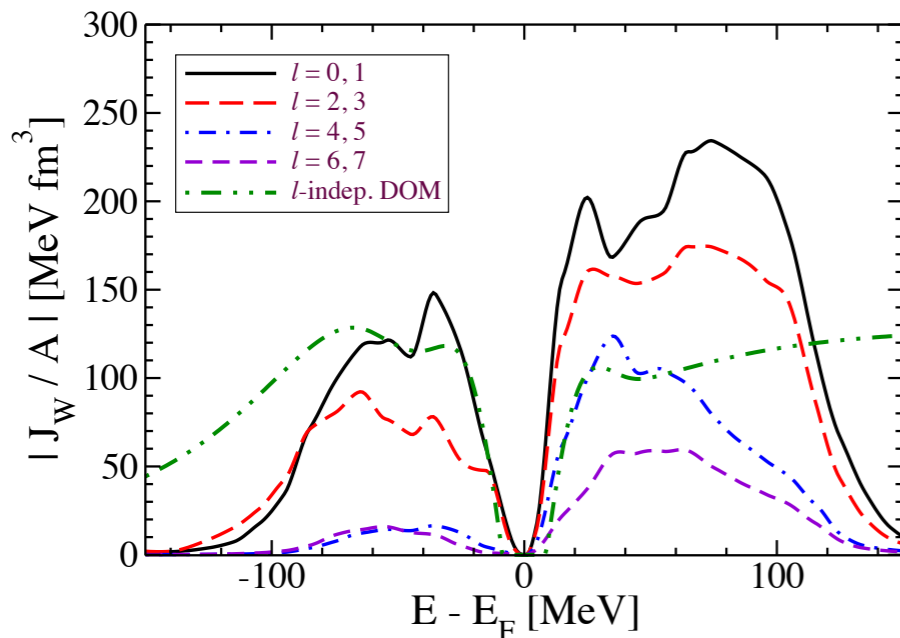
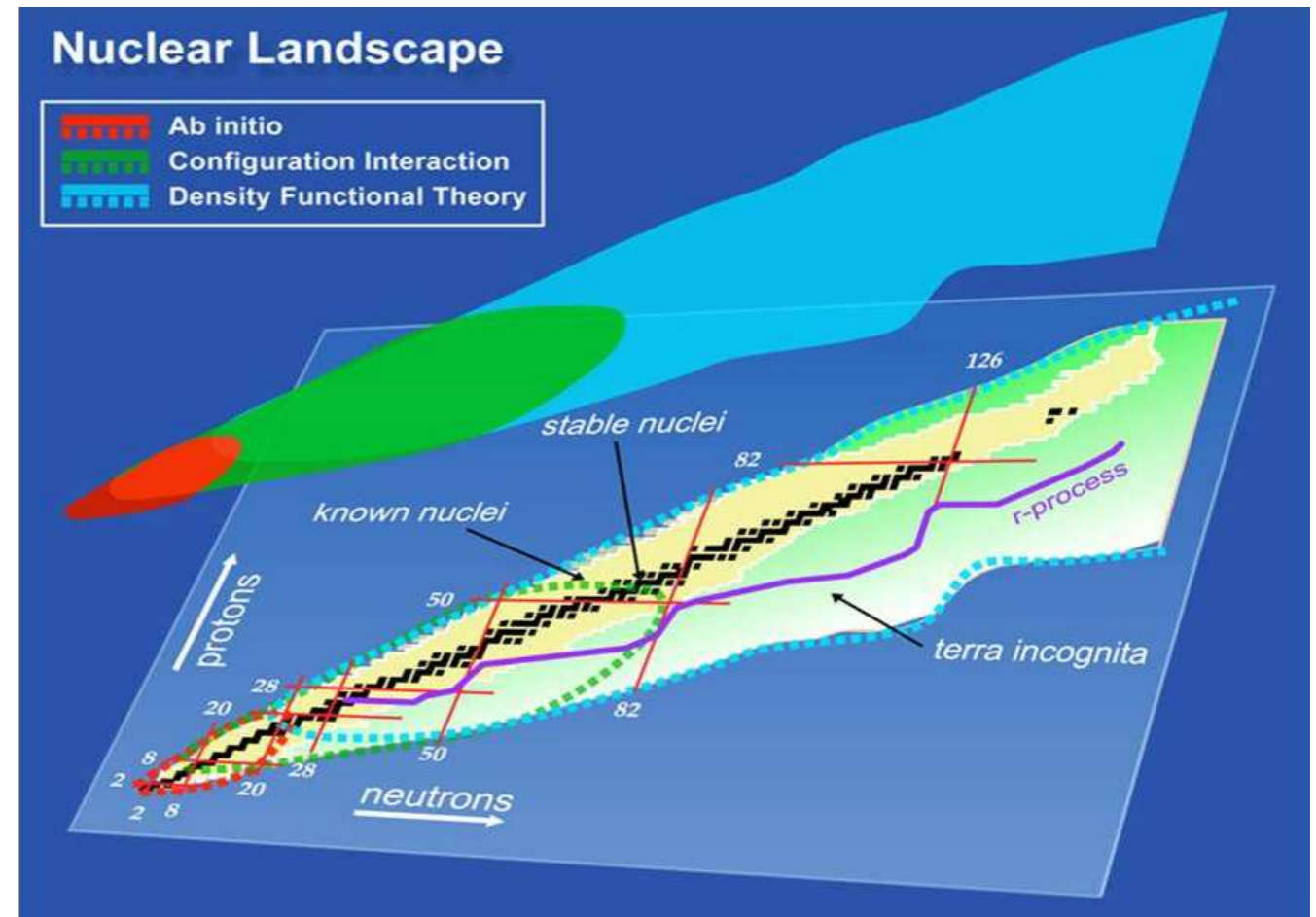
- ❖ How to extend to open-shell?
- ❖ How to link with EDF?
- ❖ How to calculate reactions?



# Towards a unified description of nuclei

Ab-initio Green's functions

- ✿ How to extend to open-shell?
  - ➡ this talk
- ✿ How to link with EDF?
  - ➡ work in progress
- ✿ How to calculate reactions?
  - ➡ TD-GF [Rios *et al.* 2011]
  - ➡ link to DOM



[Waldecker, Barbieri, Dickhoff 2011]

# State-of-the-art ab-initio nuclear structure theory

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✱ Methods for an ab-initio description of medium-mass nuclei

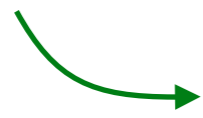
(1) Self-consistent Dyson-Green's function [Barbieri, Dickhoff, ...]

(2) Coupled-cluster [Dean, Hagen, Hjorth-Jensen, Papenbrock, ...]

(3) In-medium similarity renormalization group [Tsukiyama, Bogner, Schwenk, ...]



Similar level of accuracy



But limited to to doubly-closed-shell  $\pm 1$  and  $\pm 2$  nuclei

# State-of-the-art ab-initio nuclear structure theory

---

## ✱ Methods for an ab-initio description of medium-mass nuclei

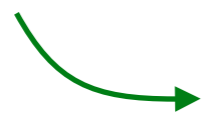
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Similar level of accuracy



But limited to to doubly-closed-shell  $\pm 1$  and  $\pm 2$  nuclei

---

## ✱ Truly open-shell nuclei

(a) Multi-reference methods: IMSRG + CI, MR-CC

(b) Single-reference methods: explicit account of pairing mandatory



Dyson-Green's functions



Gorkov-Green's functions

# Connection to *non-empirical* EDF

- \* Standard EDF parameterizations (e.g. Skyrme, Gogny, relativistic)
  - Successful in major shell where adjusted
  - Lack predictive power in new regions of interest

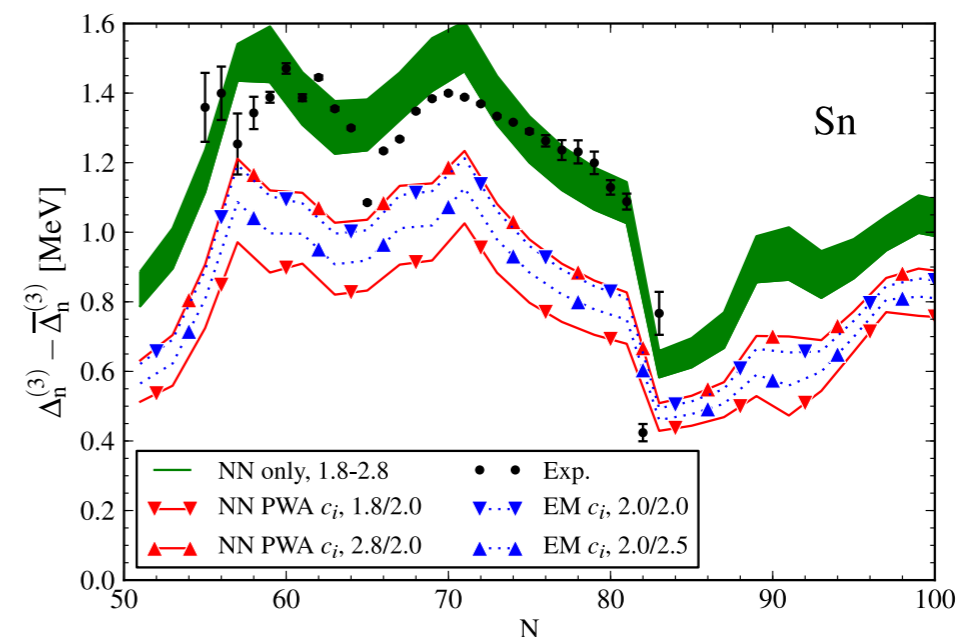
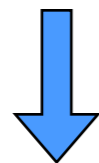


Efforts to extend and connect with more fundamental approaches

- \* Non-empirical EDF from low-momentum interactions

[Lesinski *et al.* 2011]

- Pairing channel [Lesinski *et al.*]
- Particle-hole channel [Gebremariam *et al.*]



Benchmarks needed from many-body methods that share the same features

# Green's functions: essential features

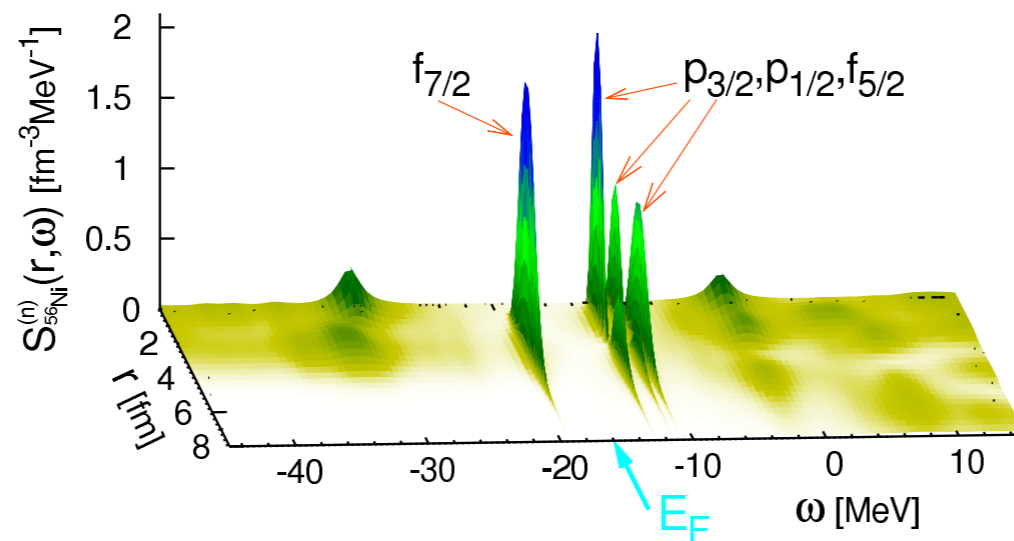
$$\Sigma = \text{[Diagram 1]} + \text{[Diagram 2]} + \text{[Diagram 3]} + \text{[Diagram 4]} + \dots$$

- ⇒ Ab-initio, self-consistent approach
- ⇒ Direct connection to observables
- ⇒ Improvability (**diagrammatic expansion**)
- ⇒ Control over many-body requirements (**conserving approximations**)

# Dyson Green's functions

## \* Spectral function

$$S_a^-(\omega) \equiv \sum_k \left| \langle \psi_k^{N-1} | a_a | \psi_0^N \rangle \right|^2 \delta(\omega - (E_0^N - E_k^{N-1})) = \frac{1}{\pi} \text{Im} G_{aa}(\omega)$$



[Barbieri 2009]

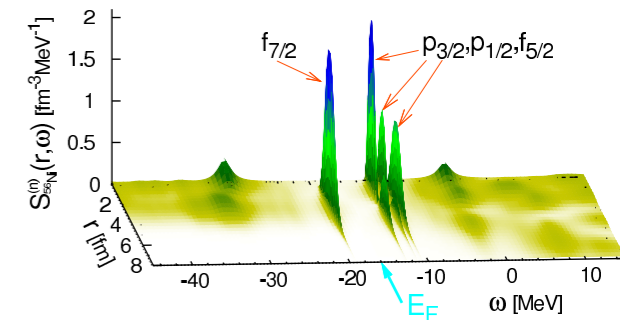




# Dyson Green's functions

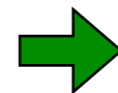
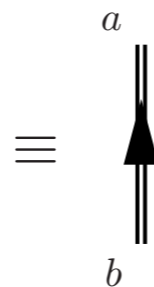
## \* Spectral function

$$S_a^-(\omega) \equiv \sum_k \left| \langle \psi_k^{N-1} | a_a | \psi_0^N \rangle \right|^2 \delta(\omega - (E_0^N - E_k^{N-1})) = \frac{1}{\pi} \text{Im} G_{aa}(\omega)$$



## \* Green's function

$$i G_{ab}(t, t') \equiv \langle \Psi_0^N | T \{ a_a(t) a_b^\dagger(t') \} | \Psi_0^N \rangle$$

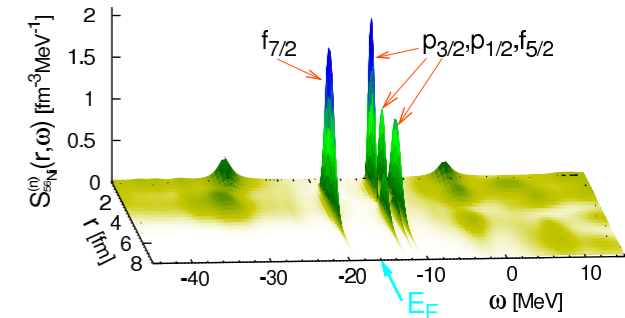


- ⇒ N-particle ground state
- ⇒ One nucleon addition and removal ( $N \pm 1$  systems)

# Dyson Green's functions

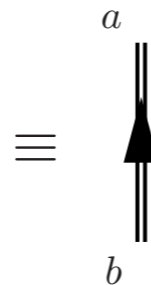
## \* Spectral function

$$S_a^-(\omega) \equiv \sum_k \left| \langle \psi_k^{N-1} | a_a | \psi_0^N \rangle \right|^2 \delta(\omega - (E_0^N - E_k^{N-1})) = \frac{1}{\pi} \text{Im} G_{aa}(\omega)$$



## \* Green's function

$$i G_{ab}(t, t') \equiv \langle \Psi_0^N | T \{ a_a(t) a_b^\dagger(t') \} | \Psi_0^N \rangle$$



- N-particle ground state
- One nucleon addition and removal ( $N \pm 1$  systems)

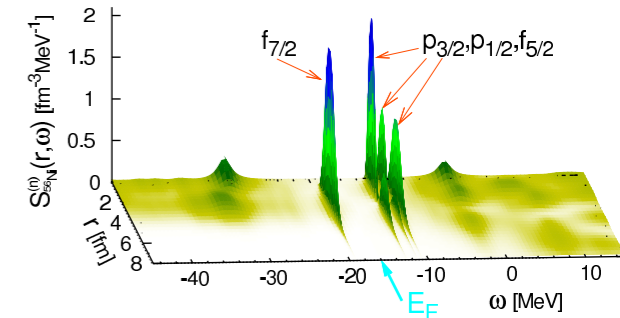
→ Contains all structure information probed by nucleon transfer

$$G_{ab}(\omega) \equiv \sum_k \frac{\langle \psi_0^N | a_a | \psi_k^{N+1} \rangle \langle \psi_k^{N+1} | a_b^\dagger | \psi_0^N \rangle}{\omega - (E_k^{N+1} - E_0^N) + i\eta} + \sum_k \frac{\langle \psi_0^N | a_b^\dagger | \psi_k^{N-1} \rangle \langle \psi_k^{N-1} | a_a | \psi_0^N \rangle}{\omega - (E_0^N - E_k^{N-1}) - i\eta}$$

# Dyson Green's functions

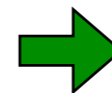
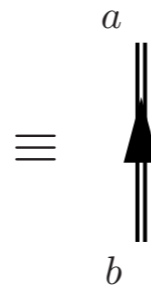
## \* Spectral function

$$S_a^-(\omega) \equiv \sum_k \left| \langle \psi_k^{N-1} | a_a | \psi_0^N \rangle \right|^2 \delta(\omega - (E_0^N - E_k^{N-1})) = \frac{1}{\pi} \text{Im} G_{aa}(\omega)$$



## \* Green's function

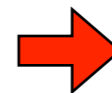
$$i G_{ab}(t, t') \equiv \langle \Psi_0^N | T \{ a_a(t) a_b^\dagger(t') \} | \Psi_0^N \rangle$$



- N-particle ground state
- One nucleon addition and removal ( $N \pm 1$  systems)

## \* Dyson equation

$$G_{ab}(\omega) = G_{ab}^{(0)}(\omega) + \sum_{cd} G_{ac}^{(0)}(\omega) \Sigma_{cd}^*(\omega) G_{db}(\omega)$$



Solution breaks down when pairing instabilities appear

# Going open-shell: Gorkov ansatz

---

- \* Formulate the expansion scheme around a Bogoliubov vacuum

→ Zeroth order already incorporates pairing

- \* Auxiliary many-body state  $|\Psi_0\rangle \equiv \sum_N^{\text{even}} c_N |\psi_0^N\rangle$

→ Mixes various particle numbers

→ Introduce a “grand-canonical” potential  $\Omega = H - \mu N$

→  $|\Psi_0\rangle$  minimizes  $\Omega_0 = \langle \Psi_0 | \Omega | \Psi_0 \rangle$

under the constraint  $N = \langle \Psi_0 | N | \Psi_0 \rangle$

→  $\Omega_0 = \sum_{N'} |c_{N'}|^2 \Omega_0^{N'} \approx E_0^N - \mu N$

# Gorkov Green's functions and equations

✱ Set of 4 Green's functions

$$i G_{ab}^{11}(t, t') \equiv \langle \Psi_0 | T \{ a_a(t) a_b^\dagger(t') \} | \Psi_0 \rangle \equiv \begin{array}{c} a \\ \uparrow \uparrow \\ b \end{array}$$

$$i G_{ab}^{21}(t, t') \equiv \langle \Psi_0 | T \{ \bar{a}_a^\dagger(t) a_b^\dagger(t') \} | \Psi_0 \rangle \equiv \begin{array}{c} \bar{a} \\ \downarrow \downarrow \\ b \end{array}$$

$$i G_{ab}^{12}(t, t') \equiv \langle \Psi_0 | T \{ a_a(t) \bar{a}_b(t') \} | \Psi_0 \rangle \equiv \begin{array}{c} a \\ \uparrow \downarrow \\ \bar{b} \end{array}$$

$$i G_{ab}^{22}(t, t') \equiv \langle \Psi_0 | T \{ \bar{a}_a^\dagger(t) \bar{a}_b(t') \} | \Psi_0 \rangle \equiv \begin{array}{c} \bar{a} \\ \downarrow \uparrow \\ \bar{b} \end{array}$$

[Gorkov 1958]



$$\mathbf{G}_{ab}(\omega) = \mathbf{G}_{ab}^{(0)}(\omega) + \sum_{cd} \mathbf{G}_{ac}^{(0)}(\omega) \boldsymbol{\Sigma}_{cd}^*(\omega) \mathbf{G}_{db}(\omega)$$

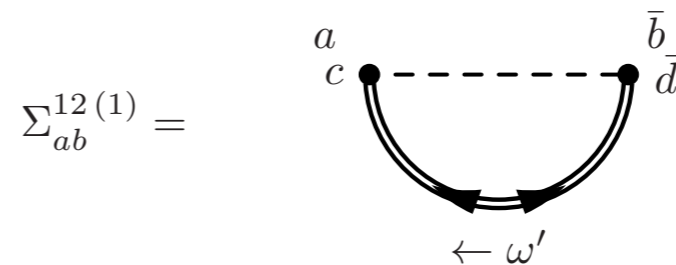
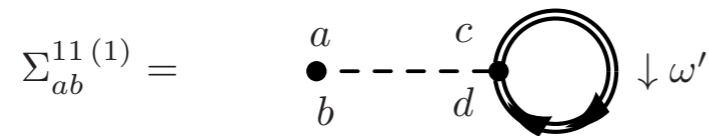
Gorkov equations

$$\boldsymbol{\Sigma}_{ab}^*(\omega) \equiv \begin{pmatrix} \Sigma_{ab}^{*11}(\omega) & \Sigma_{ab}^{*12}(\omega) \\ \Sigma_{ab}^{*21}(\omega) & \Sigma_{ab}^{*22}(\omega) \end{pmatrix}$$

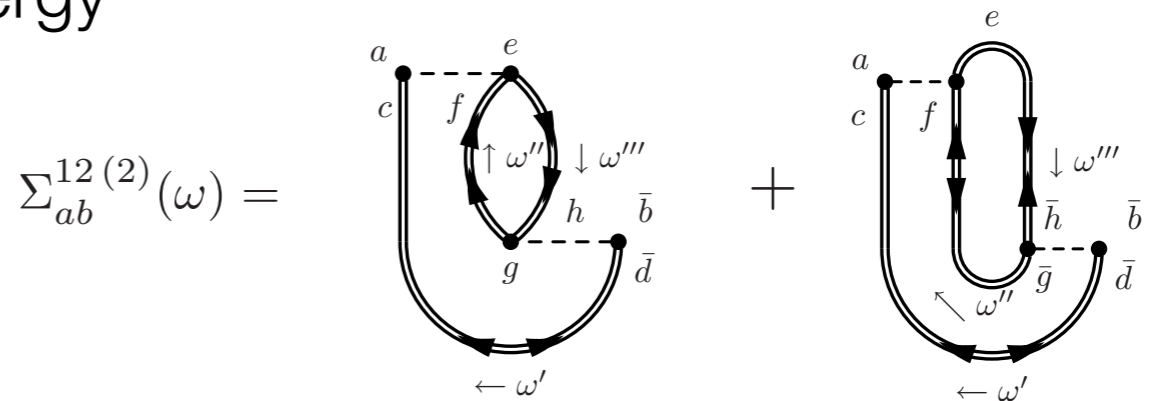
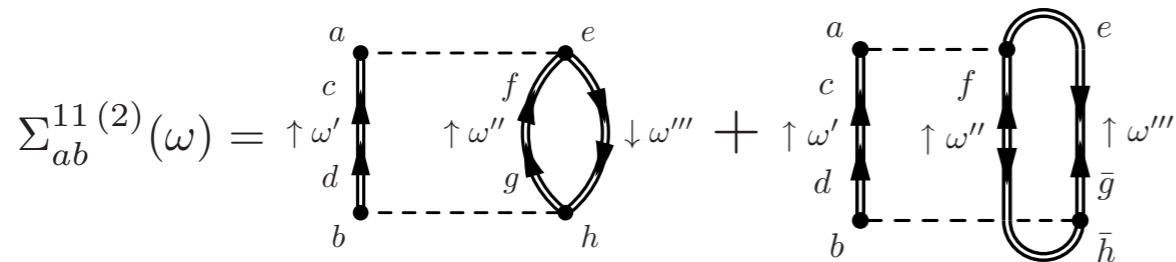
$$\boldsymbol{\Sigma}_{ab}^*(\omega) \equiv \boldsymbol{\Sigma}_{ab}(\omega) - \mathbf{U}_{ab}$$

# 1<sup>st</sup> & 2<sup>nd</sup> order diagrams and eigenvalue problem

\* 1<sup>st</sup> order  $\Rightarrow$  energy-independent self-energy



\* 2<sup>nd</sup> order  $\Rightarrow$  energy-dependent self-energy



\* Gorkov equations  $\longrightarrow$  eigenvalue problem

$$\sum_b \begin{pmatrix} t_{ab} - \mu_{ab} + \Sigma_{ab}^{11}(\omega) & \Sigma_{ab}^{12}(\omega) \\ \Sigma_{ab}^{21}(\omega) & -t_{ab} + \mu_{ab} + \Sigma_{ab}^{22}(\omega) \end{pmatrix} \Big|_{\omega_k} \begin{pmatrix} \mathcal{U}_b^k \\ \mathcal{V}_b^k \end{pmatrix} = \omega_k \begin{pmatrix} \mathcal{U}_a^k \\ \mathcal{V}_a^k \end{pmatrix}$$

$$\begin{aligned} \mathcal{U}_a^{k*} &\equiv \langle \Psi_k | \bar{a}_a^\dagger | \Psi_0 \rangle \\ \mathcal{V}_a^{k*} &\equiv \langle \Psi_k | a_a | \Psi_0 \rangle \end{aligned}$$

# Gorkov equations

---

$$\sum_b \begin{pmatrix} t_{ab} - \mu_{ab} + \Sigma_{ab}^{11}(\omega) & \Sigma_{ab}^{12}(\omega) \\ \Sigma_{ab}^{21}(\omega) & -t_{ab} + \mu_{ab} + \Sigma_{ab}^{22}(\omega) \end{pmatrix} \Big|_{\omega_k} \begin{pmatrix} \mathcal{U}_b^k \\ \mathcal{V}_b^k \end{pmatrix} = \omega_k \begin{pmatrix} \mathcal{U}_a^k \\ \mathcal{V}_a^k \end{pmatrix}$$



$$\begin{pmatrix} T - \mu + \Lambda & \tilde{h} & \mathcal{C} & -\mathcal{D}^\dagger \\ \tilde{h}^\dagger & -T + \mu - \Lambda & -\mathcal{D}^\dagger & \mathcal{C} \\ \mathcal{C}^\dagger & -\mathcal{D} & E & 0 \\ -\mathcal{D} & \mathcal{C}^\dagger & 0 & -E \end{pmatrix} \begin{pmatrix} \mathcal{U}^k \\ \mathcal{V}^k \\ \mathcal{W}_k \\ \mathcal{Z}_k \end{pmatrix} = \omega_k \begin{pmatrix} \mathcal{U}^k \\ \mathcal{V}^k \\ \mathcal{W}_k \\ \mathcal{Z}_k \end{pmatrix}$$

Energy *independent* eigenvalue problem

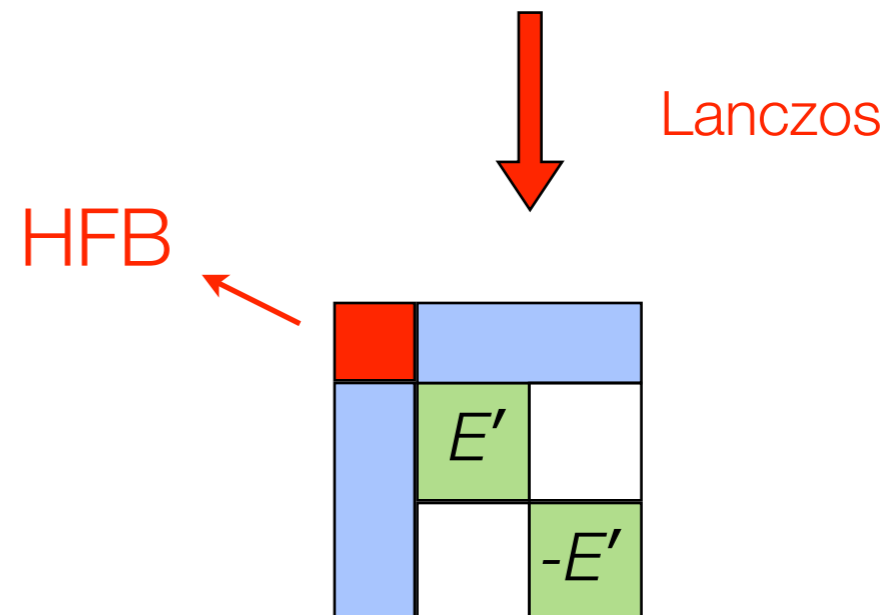
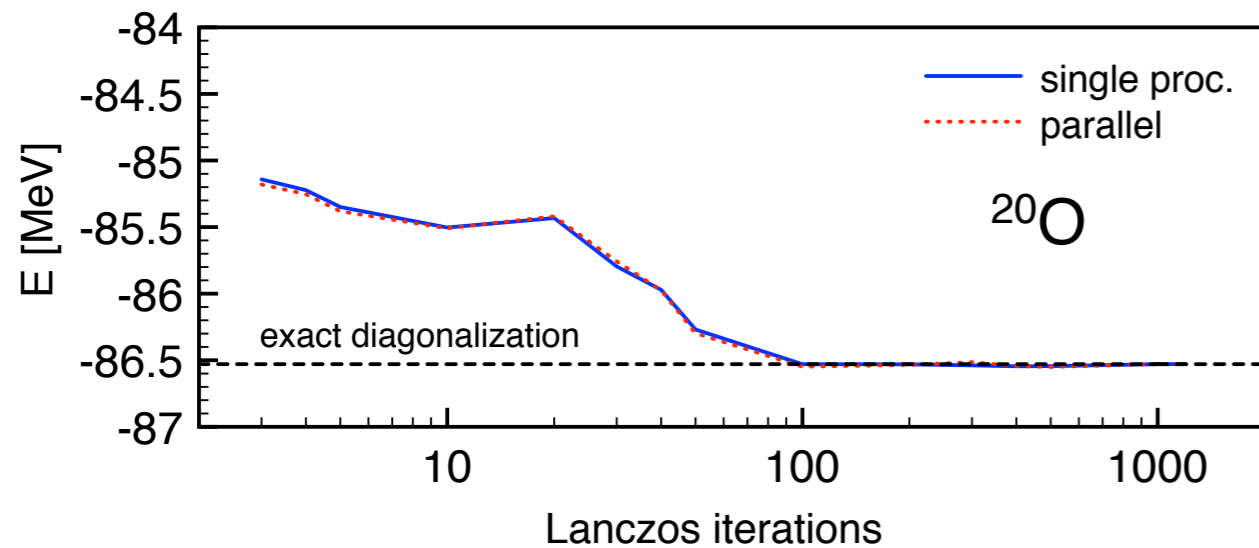
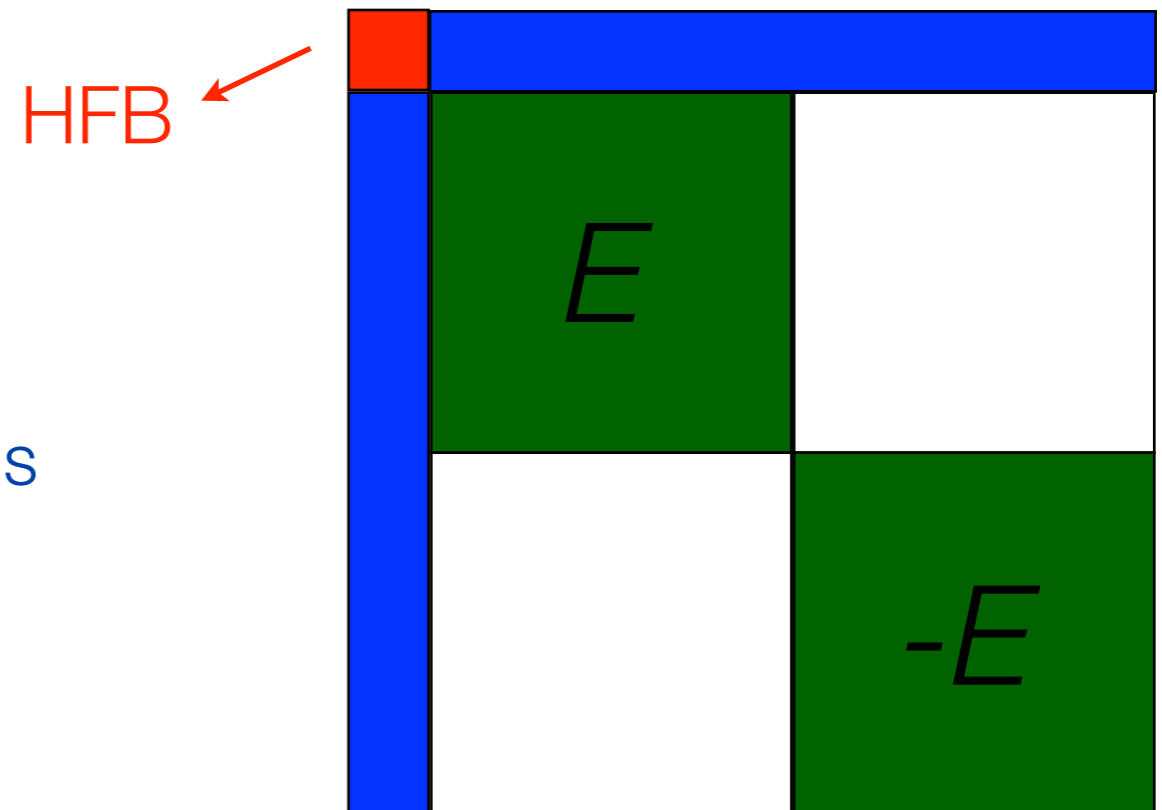
with the normalization condition

$$\sum_a \left[ |\mathcal{U}_a^k|^2 + |\mathcal{V}_a^k|^2 \right] + \sum_{k_1 k_2 k_3} \left[ |\mathcal{W}_k^{k_1 k_2 k_3}|^2 + |\mathcal{Z}_k^{k_1 k_2 k_3}|^2 \right] = 1$$

# Lanczos projection

$$\begin{pmatrix} T - \mu + \Lambda & \tilde{h} & C & -D^\dagger \\ \tilde{h}^\dagger & -T + \mu - \Lambda & -D^\dagger & C \\ C^\dagger & -D & E & 0 \\ -D & C^\dagger & 0 & -E \end{pmatrix} \begin{pmatrix} U^k \\ V^k \\ W_k \\ Z_k \end{pmatrix} = \omega_k \begin{pmatrix} U^k \\ V^k \\ W_k \\ Z_k \end{pmatrix}$$

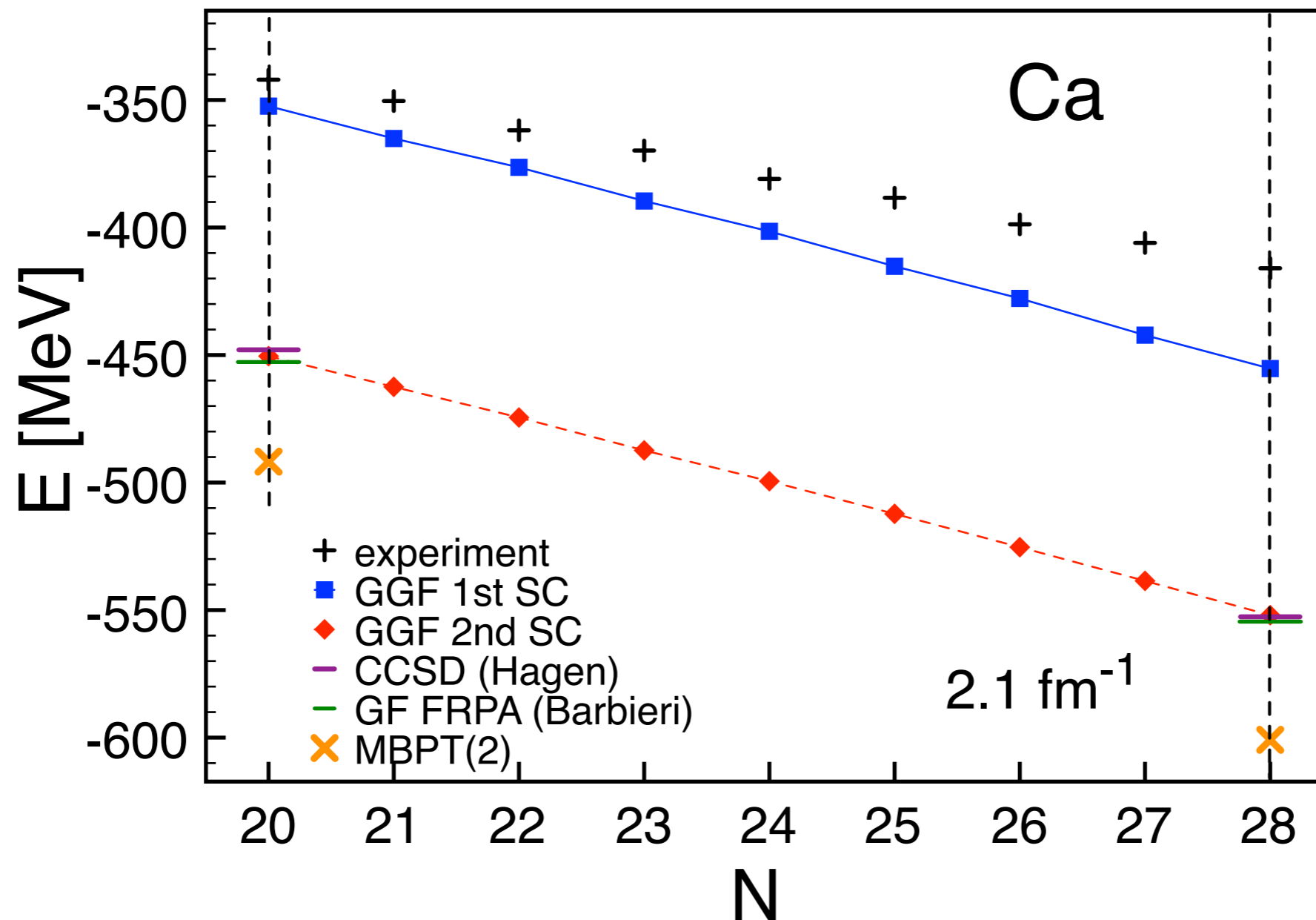
- Conserves moments of spectral functions
- Equivalent to exact diagonalization for  $N_L \rightarrow \dim(E)$





# Binding energies

✱ Systematic along isotopic/isotonic chains become available



# Spectrum and spectroscopic factors

## \* Separation energy spectrum

$$G_{ab}^{11}(\omega) = \sum_k \left\{ \frac{\mathcal{U}_a^k \mathcal{U}_b^{k*}}{\omega - \omega_k + i\eta} + \frac{\bar{\mathcal{V}}_a^{k*} \bar{\mathcal{V}}_b^k}{\omega + \omega_k - i\eta} \right\}$$

Lehmann representation

where

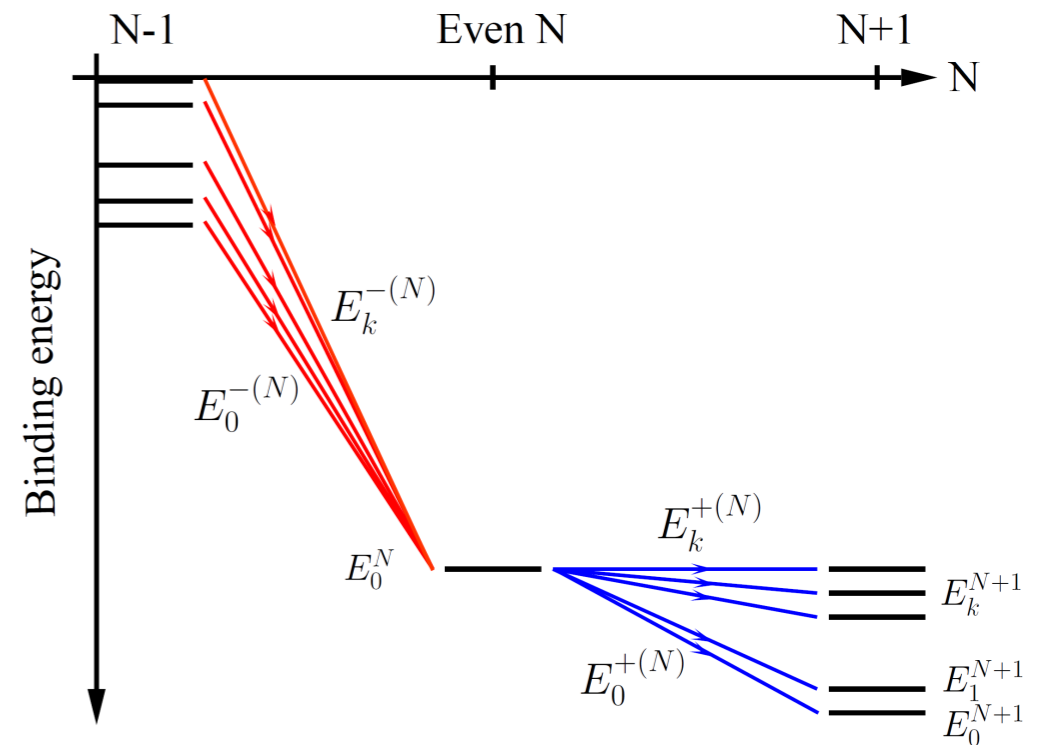
$$\begin{cases} \mathcal{U}_a^{k*} \equiv \langle \Psi_k | a_a^\dagger | \Psi_0 \rangle \\ \mathcal{V}_a^{k*} \equiv \langle \Psi_k | \bar{a}_a | \Psi_0 \rangle \end{cases}$$

and

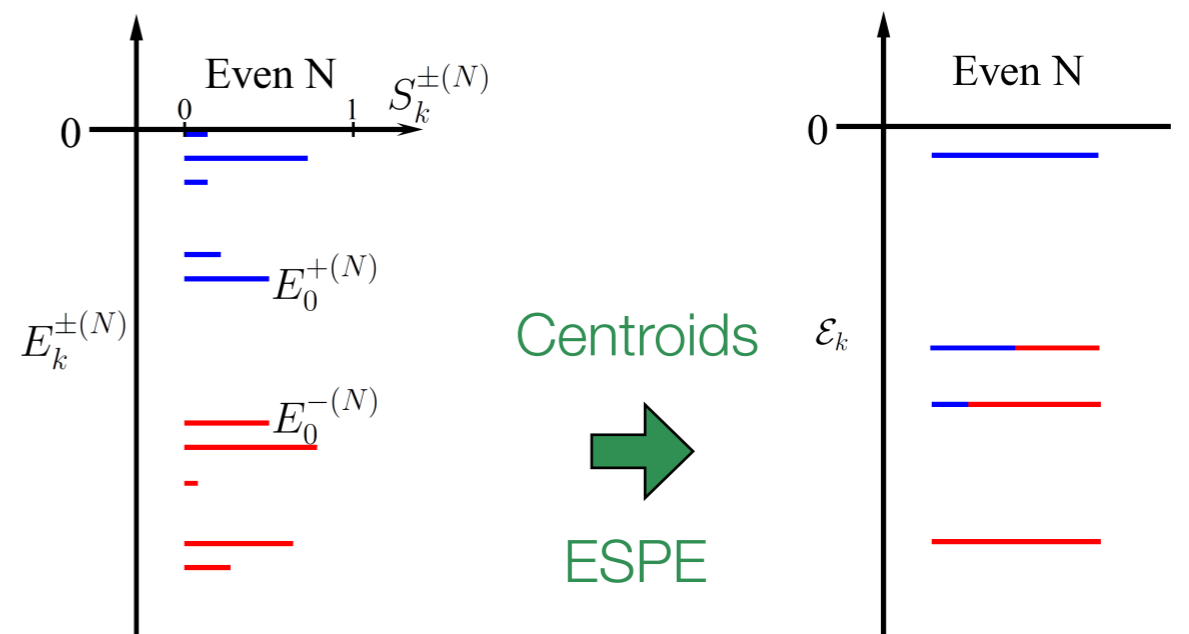
$$\begin{cases} E_k^{+(N)} \equiv E_k^{N+1} - E_0^N \\ E_k^{-(N)} \equiv E_0^N - E_k^{N-1} \end{cases}$$

## \* Spectroscopic factors

$$\begin{aligned} \mathcal{S}_k^+ &\equiv \sum_a |\langle \psi_k | a_a^\dagger | \psi_0 \rangle|^2 = \sum_a |\mathcal{U}_a^k|^2 \\ \mathcal{S}_k^- &\equiv \sum_a |\langle \psi_k | a_a | \psi_0 \rangle|^2 = \sum_a |\mathcal{V}_a^k|^2 \end{aligned}$$

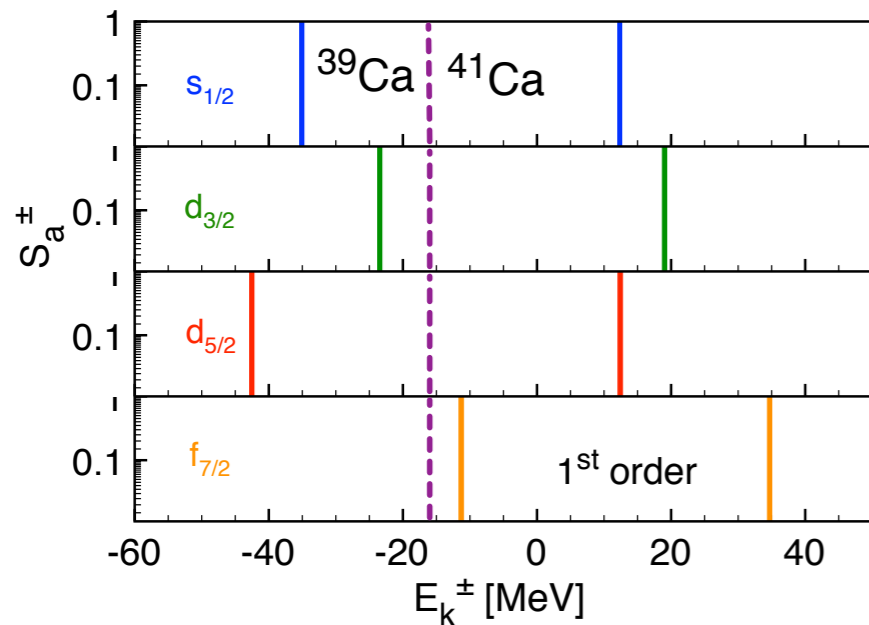


Separation energies + transfer strengths



# Spectral function

Dyson 1<sup>st</sup> order (HF)

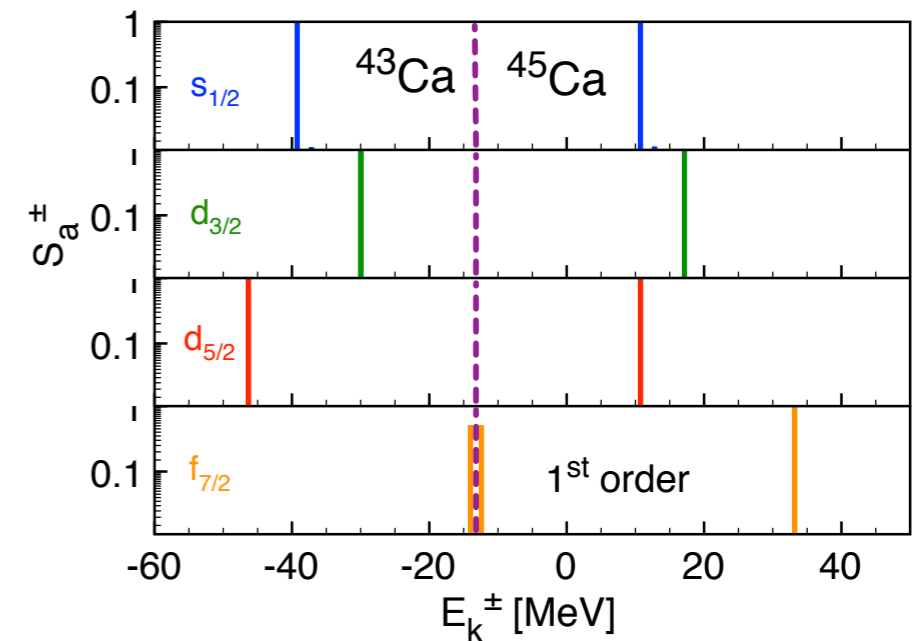


Fragmentation

Static pairing

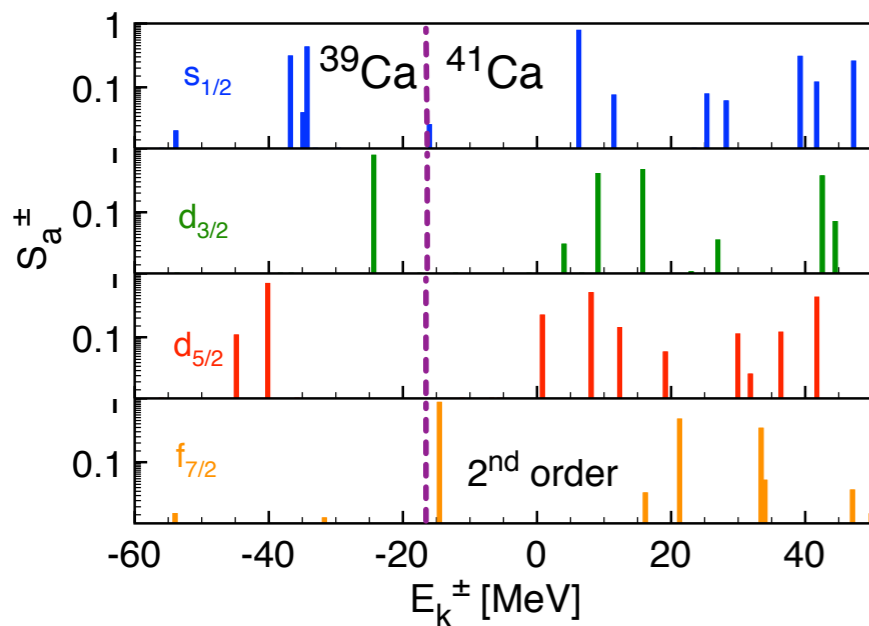


Gorkov 1<sup>st</sup> order (HFB)

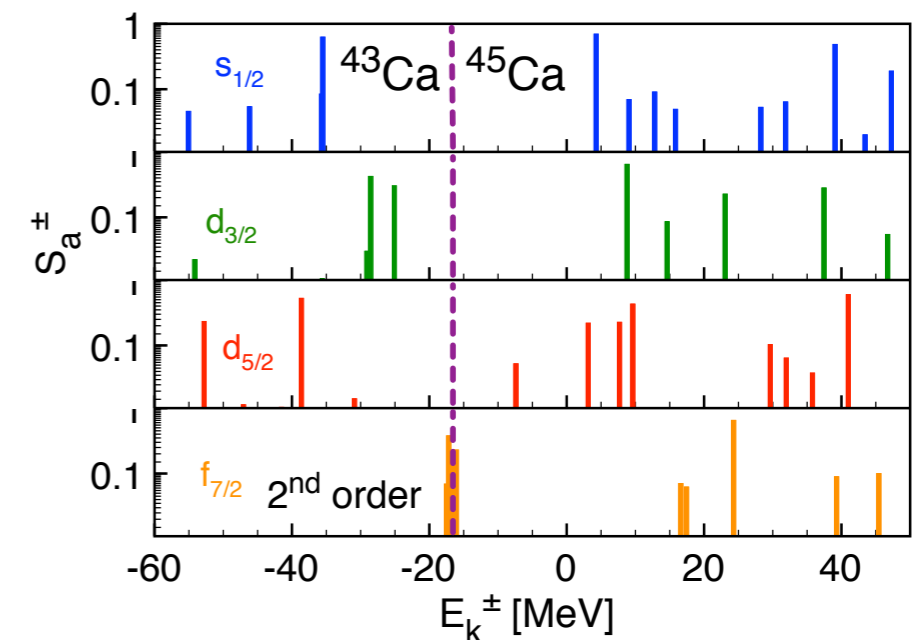


Dynamical fluctuations

Dyson 2<sup>nd</sup> order



Gorkov 2<sup>nd</sup> order



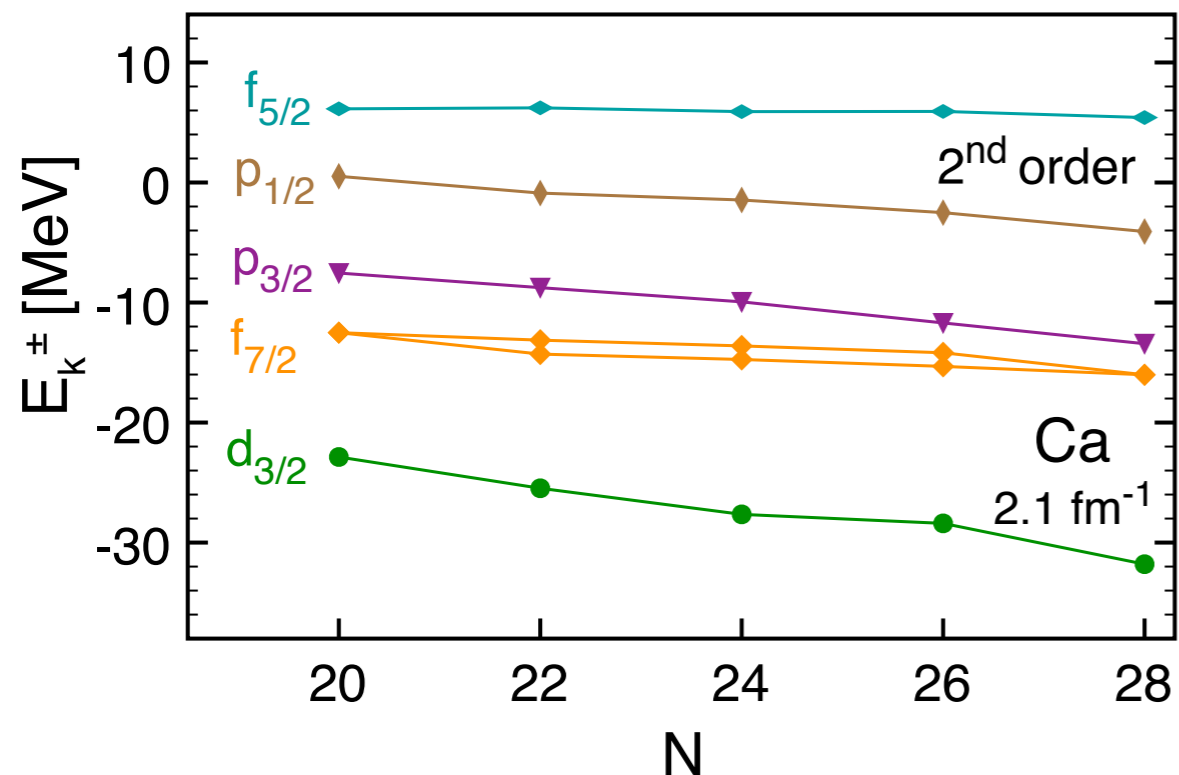
# Shell structure evolution

✱ ESPE collect fragmentation of “single-particle” strengths from both  $N \pm 1$

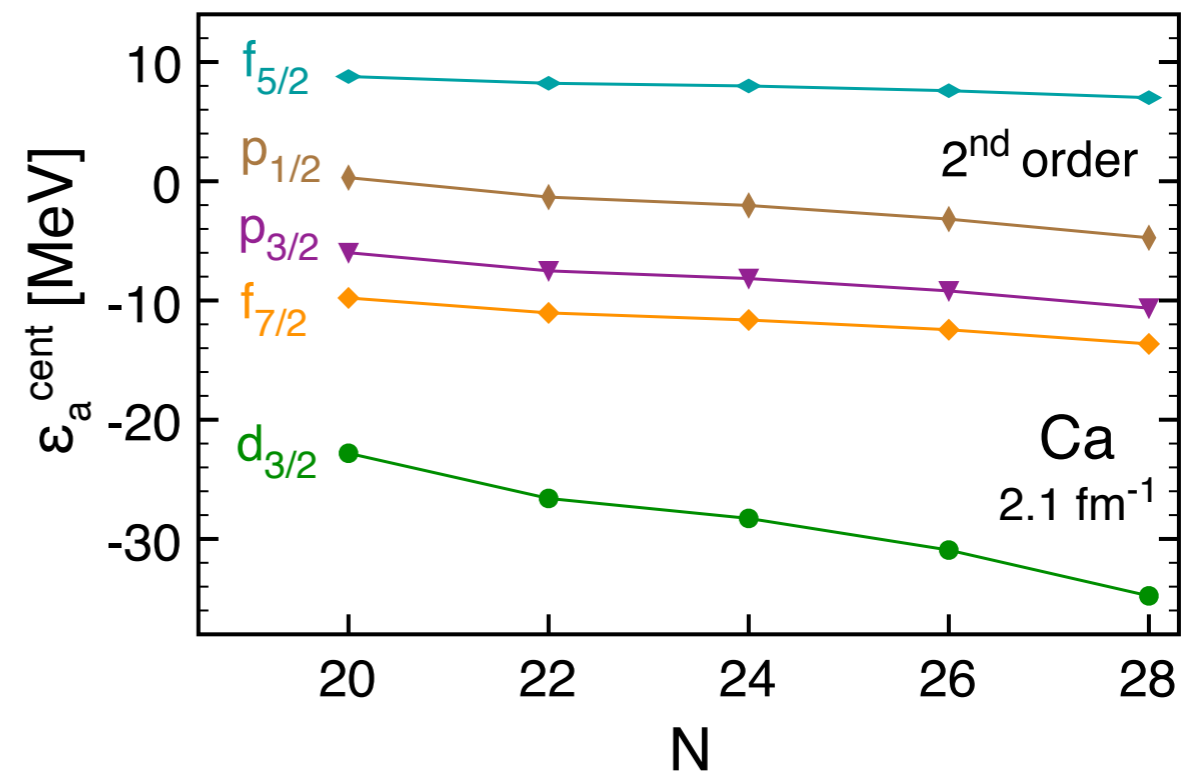
$$\epsilon_a^{cent} \equiv h_{ab}^{cent} \delta_{ab} = t_{aa} + \sum_{cd} \bar{V}_{acad}^{NN} \rho_{dc}^{[1]} + \sum_{cdef} \bar{V}_{acdaef}^{NNN} \rho_{efcd}^{[2]} \equiv \sum_k \mathcal{S}_k^{+a} E_k^+ + \sum_k \mathcal{S}_k^{-a} E_k^-$$

[Baranger 1970, Duguet *et al.* 2011]

Quasiparticle peaks



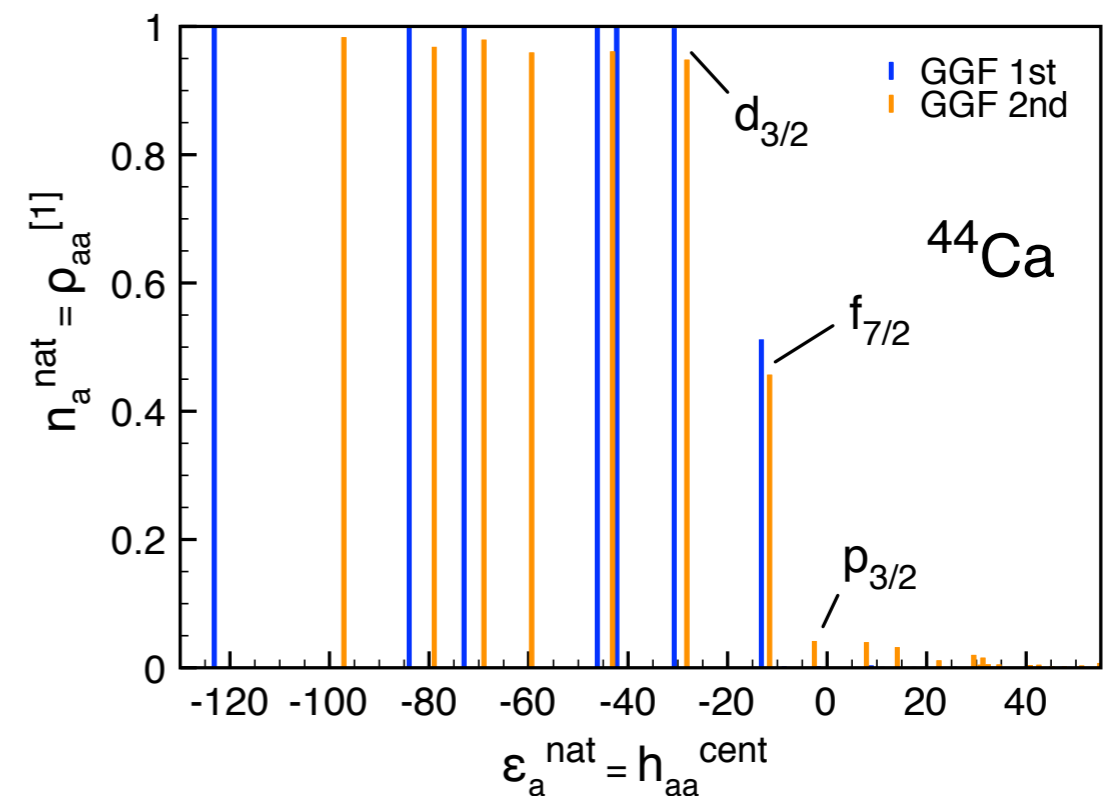
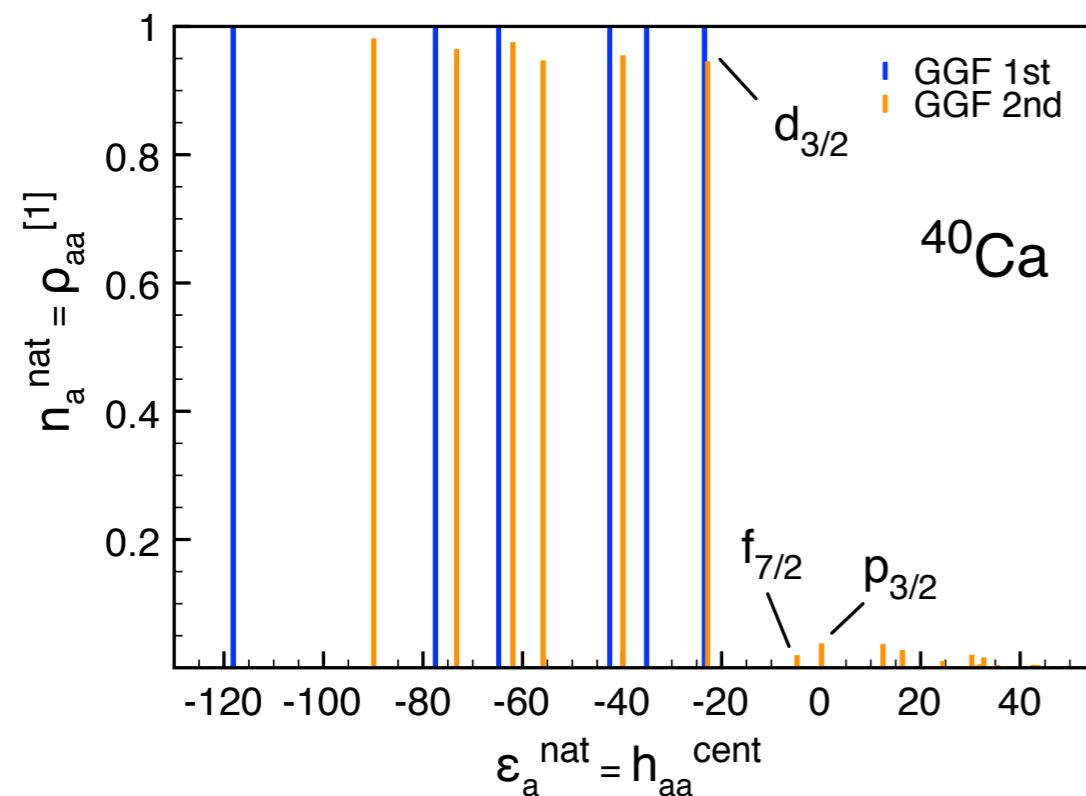
Centroids



# Natural single-particle occupation

✱ Natural orbit  $a$ :  $\rho_{ab}^{[1]} = n_a^{\text{nat}} \delta_{ab}$

✱ Associated energy:  $\varepsilon_a^{\text{nat}} = h_{aa}^{\text{cent}}$

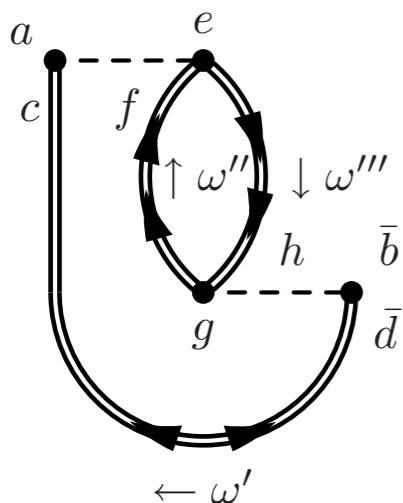
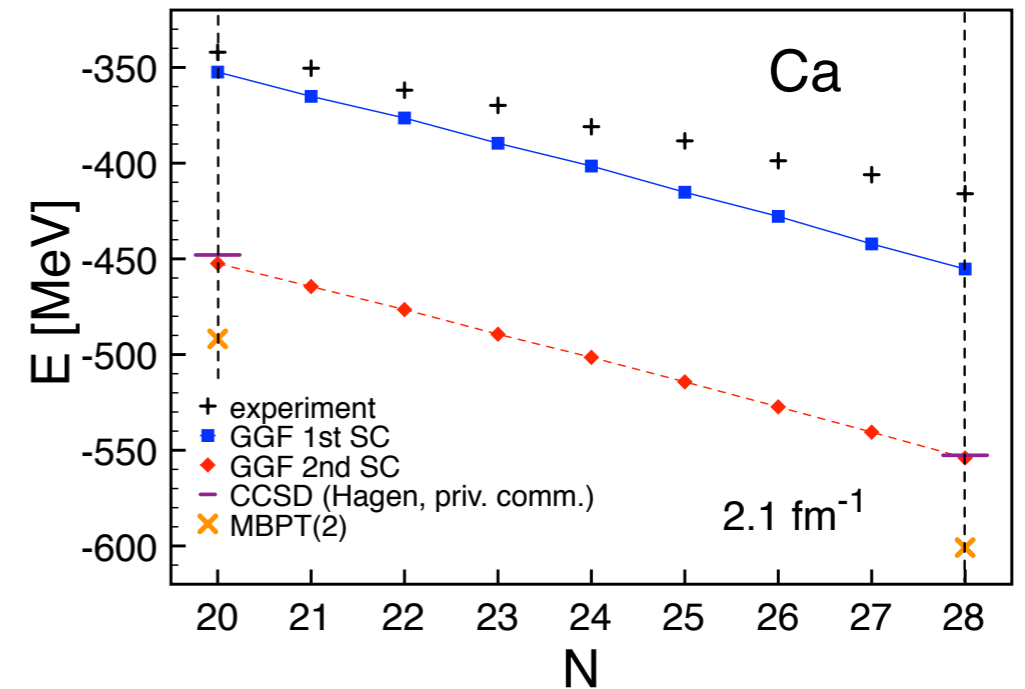


✱ Dynamical correlations similar for doubly-magic and semi-magic

✱ Static pairing essential to open-shells

# Conclusions & Outlook

- ✱ Gorkov-Green's functions:  
first ab-initio **open-shell** calculations
- ✱ Provide optical potentials  
for **reaction models**
- ✱ Provide constraints for next-generation  
**Energy Density Functionals**



- ✱ Implementation of three-body forces
- ✱ Formulation of **particle-number restored** Gorkov theory
- ✱ Improvement of the self-energy expansion