

Non-Empirical Pairing Energy Functional for nuclei

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Outline

1 Introduction

- Nuclear theory: goals and methods
- Energy Density Functional methods

2 Non-empirical energy functional

- The pairing part of the EDF as a first step
- Low momentum interactions
- Separable operator representation of $V_{\text{low } k} + V_{\text{Coul}}$

3 Pairing gaps in finite nuclei

- Implementation for finite nuclei calculations
- Results including nuclear, Coulomb and CSB terms

4 Results for soft versus hard NN interactions

- Dependence of pairing gaps on the RG scale Λ
- Fully microscopic calculations in infinite nuclear matter

5 Summary and outlook

Theory of nuclei

Ultimate goals

- Comprehensive and unified description of all nuclei
- **From basic interactions between nucleons + link to QCD**
- Understand states of nuclear matter in astrophysical environments

Difficulties

- Self-bound, two-components quantum many-fermions system
- Complex interaction from low-energy regime of QCD
 - Tensor and spin-orbit components
 - Unnaturally large scattering lengths
 - NNN unavoidable
 - Repulsive core and strong tensor at short distances?
- Unified description from deuteron to SHE nuclei to NS
- Need to extrapolate to unknown regions

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Ultimate goal

Ground state

Mass, deformation



Spectroscopy

Spectroscopy



Collective modes

RPA, QRPA, GCM



Reaction properties

Fusion, transfer...



Heavy elements

Fission, fusion, SHE



Exotic behaviors

Drip-lines, halos

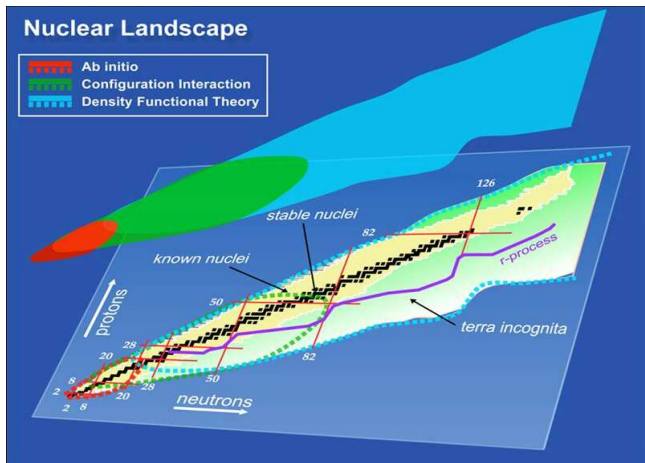


Astrophysics

NS, SN, r-process



Which theoretical method(s)?



- No “one size fits all” theory for nuclei
- All theoretical approaches need to be linked

Energy Density Functional (EDF) approaches

Basic elements

- Approaches **not** based on a correlated wave-function
- Energy is postulated to be a functional of one-body density (matrices)
- **Symmetry breaking** is at the heart of the method
- Two formulations (i) **Single-Reference** (ii) **Multi-Reference**

Pros

- Use of full single-particle space
- Quantal + collective picture
- Universality of EDF ($A \gtrsim 16$)
- Ground-state description
- Static (smooth) correlations

Difficulties

- No universal parametrization
- Empirical \neq predictive power
- Spectroscopy / odd nuclei
- Dynamical (fluctuating) correlations
- Limited accuracy ($\sigma_{2135}^{mass} \approx 700$ keV)

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Pairing correlations in nuclei

$T = 1$ pairing in nuclei

- nn/pp superfluidity impacts all low-energy properties of nuclei
- Non-perturbative channel to be treated explicitly

Most impacted observables

- Lowest two-qp states in even-even nuclei \approx measures the "gap"
- Odd-even mass staggering (OEMS) \approx measures the "gap"
- Collective excitations
 - Moment of inertia of rotational bands
 - Low-lying vibrational states
 - Shape isomers from intruders
- Pair transfer
- Competition of pro- and anti-halo effects on the one-body density
- Neutron star physics
 - Glitches in the inner crust
 - Neutrino emission process
 - Heat diffusion

Single-reference EDF methods (1)

Elements of formalism

- $\mathcal{E}[\rho, \kappa^*, \kappa]$ = functional of one-body density matrices

$$\rho_{ji} = \langle \Phi | c_i^\dagger c_j | \Phi \rangle \quad ; \quad \kappa_{ji} = \langle \Phi | c_i c_j | \Phi \rangle$$

- $|\Phi\rangle$ = **auxiliary**/symmetry-breaking/product **state of reference**
- Minimizing $\mathcal{E}[\rho, \kappa^*, \kappa]$ leads to Hartree-Fock-Bogoliubov-like equations

$$\begin{pmatrix} h - \lambda & \Delta \\ -\Delta^* & -h^* + \lambda \end{pmatrix} \begin{pmatrix} U_i \\ V_i \end{pmatrix} = E_i \begin{pmatrix} U_i \\ V_i \end{pmatrix}$$

- Effective potentials and vertices are defined through

$$h_{ij} \equiv \frac{\delta \mathcal{E}}{\delta \rho_{ji}} \equiv \sum_{kl} \overline{v}_{ikjl}^{ph} \rho_{lk} \quad ; \quad \Delta_{ij} \equiv \frac{\delta \mathcal{E}}{\delta \kappa_{ij}^*} \equiv \frac{1}{2} \sum_{kl} \overline{v}_{ijkl}^{pp} \kappa_{kl}$$

- $\overline{v}^{ph} / \overline{v}^{pp}$ = Consistent many-body expansion in terms of NN/NNN
- Quasiparticle w.f. (U_i, V_i) , energy E_i , densities...

Single-reference EDF methods (2)

Empirical parameterizations of $\mathcal{E}[\rho, \kappa, \kappa^*]$; e.g. Skyrme or Gogny

$$\mathcal{E}[\rho, \kappa, \kappa^*] \equiv \text{[Diagram 1]} \cdot \dots \cdot \text{[Diagram 2]} + \text{[Diagram 3]} + \text{[Diagram 4]}$$

Skyrme, Gogny...
DDDI, Gogny...

- Tremendous successes for known nuclei
- Existing parameterizations are over/under constrained
- "Asymptotic freedom" as one enters "the next major shell"

Pairing part: example of EDF from DDDI

$$\mathcal{E}^{\kappa\kappa} = \sum_q \int d\vec{r} A^{\tilde{\rho}\tilde{\rho}}(\vec{r}) |\tilde{\rho}_q(\vec{r})|^2 \equiv - \sum_q \int d\vec{r} \Delta_q(\vec{r}) \tilde{\rho}_q^*(\vec{r})$$

$$\Delta_{ij}^q \equiv \frac{\delta \mathcal{E}^{\kappa\kappa}}{\delta \kappa_{ij}^{q*}} = \int d\vec{r} \Psi_{ij}^{q*}(\vec{r}) \Delta_q(\vec{r})$$

- $\tilde{\rho}_q(\vec{r})$ = local pair density / $A^{\tilde{\rho}\tilde{\rho}}(\vec{r})$ = density-dependent coupling
- (Quasi-) local pairing EDF must be regularized/renormalized

Single-reference EDF methods (3)

Performance of existing pairing EDFs

- Moment of inertia of super-deformed bands = success story of the 90'
- OEMS \approx ok but missing systematic/detailed characterization
- QP excitations = missing systematic characterization
- Divergence of predictions in the "next major shell"

Crucial undergoing works

- Enrich the analytical structure of empirical functionals
- Improve fitting protocols = data, algorithm and post-analysis

One can also propose a complementary approach. . .

- Known data hardly constrain non-trivial characteristics of pairing EDF
- Interesting not to rely entirely on fitting data

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Constructing non-empirical EDFs for nuclei

Long term objective

Build non-empirical EDF in place of existing models

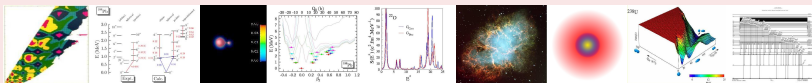
Empirical



Predictive?



Finite nuclei



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Empirical



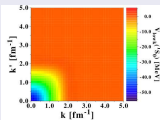
Predictive?

Non-empirical

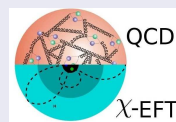


Predictive...

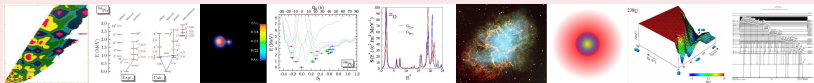
Low- k NN+NNN



QCD / χ -EFT



Finite nuclei



Long term project and collaboration

Design non-empirical Energy Density Functionals

- Bridge with "ab-initio" many-body techniques
- Calculate properties of heavy/complex nuclei from NN+NNN
- Controlled calculations with theoretical error bars



First step: pairing part of the EDF

Motivations for a non-empirical approach

- Empirical schemes lack predictive power
- Microscopic origin of superfluidity in finite nuclei?
 - Contribution from the direct term of V_{NN} (1S_0 , 1D_2 , 3PF_2)?
 - Coupling to density/spin/isospin fluctuations: 40%?
 - Absolute value/isotopic trend is of great interest

Start with v^{pp} built at 1st order in V_{NN} (nuclear + Coulomb)

- Single channel (1S_0) dominates at sub-nuclear densities
- Virtual state at $E \simeq 0$ makes V_{NN} almost separable in 1S_0

Nuclear interactions and the Renormalization Group

Approach

- $V(\vec{k}, \vec{k}', \Lambda = \infty) = V^{\text{hard}}(\vec{k}, \vec{k}')$
- Run down Λ
- Keep $\delta^{S L J}(k)$ and E_{Deuteron}

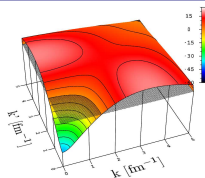
General Properties

- Vacuum interaction
- Universal $V_{NN}(\Lambda \approx 2) \equiv V_{\text{low } k}$
- $V_{\text{low } k}$ is perturbative

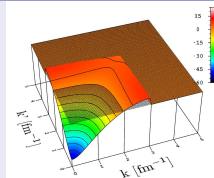
Crucial points

- $H = V_{NN}(\Lambda) + V_{NNN}(\Lambda) + \dots$
- $\partial_{\Lambda} A \neq 0 \Rightarrow$ missing pieces
- Ex: omitted $NNN(\Lambda)$

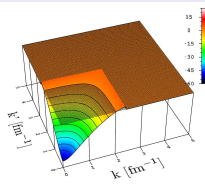
Convergence of the RG flow



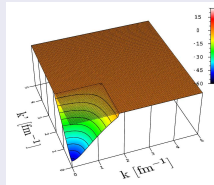
$\Lambda = 5.0 \text{ fm}^{-1}$



$\Lambda = 3.0 \text{ fm}^{-1}$



$\Lambda = 2.4 \text{ fm}^{-1}$



$\Lambda = 1.8 \text{ fm}^{-1}$

Finite nuclei calculations

Low-momentum interactions for finite nuclei calculations

- High-precision bare interactions with regularized hard-core
- Good starting point for structure calculations through EDF method?

$V_{\text{low } k}$ is given as tables of numbers

Produce analytical operatorial representation

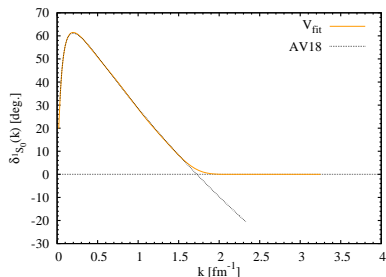
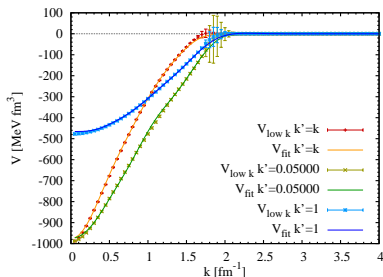
- Why?
 - Interest to understand encoded operatorial structure
 - Perform integrals analytically in codes
- Which representation?
 - Gaussian/Gogny-like (V. Rotival)
 - Sum of separable terms (T. Lesinski)

Separable representation of $V_{\text{low } k}(\Lambda) + V_{\text{Coul}}$

- High precision separable representation of rank n

$$V_n^1 S_0(k, k', \Lambda) = \sum_{\alpha, \beta=1}^n g_\alpha(k) \lambda_{\alpha\beta} g_\beta(k')$$

- Fit of $g_\alpha(k)$ and $\lambda_{\alpha\beta}$ to $V_{\text{low } k}^1 S_0(k, k', \Lambda)$ and $\delta^1 S_0(k)$
- For $\Lambda = 1.8/4.0/\infty$ fm $^{-1}$ (rank 2/4/15) and smooth cutoff



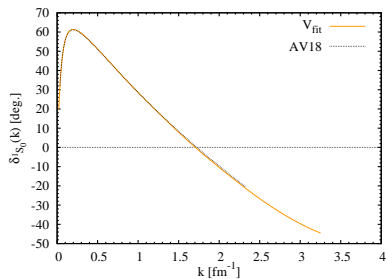
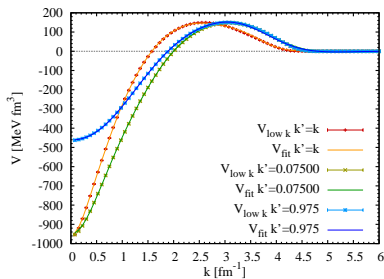
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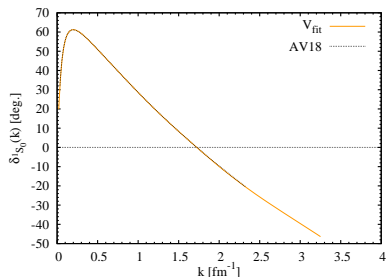
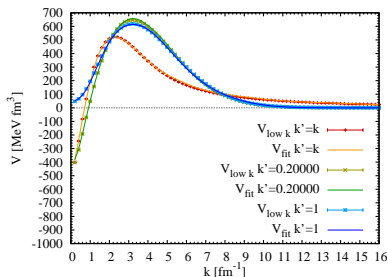
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- \exists separable representation of $V_{\text{Coul}, \ell=0}^a(k, k')$

Coulomb interaction

Need to incorporate Coulomb effects on proton gaps

- Only one such published calculation so far: Madrid group (Gogny)
- Simplified treatment of e.m. interaction (Coulomb)

Truncate the Coulomb interaction at $r = a > 2R_{\text{nucleus}}$

- A separable expansion exists (keep 1S_0 part here)

$$V_{\text{Coul}, \ell=0}^a(k, k') = 4\pi e^2 a^2 \sum_{n=0}^{\infty} (2n+1) j_n^2\left(\frac{ak}{2}\right) j_n^2\left(\frac{ak'}{2}\right),$$

$$\lambda_{\alpha\beta} = e^2 a^2 (2\alpha+1) \delta_{\alpha\beta}$$

$$g_{\alpha}(k) = \sqrt{4\pi} j_{\alpha}^2\left(\frac{ak}{2}\right)$$

$$G_{\alpha}(r) = \frac{1}{\sqrt{\pi} a^2 r} P_{\alpha}\left(1 - 2\left(\frac{r}{a}\right)^2\right) \text{ for } r \leq a$$

- ~ 15 terms needed (peanuts !)

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EDF calculations in spherical nuclei (1)

- Separable force in coordinate-space [$\mathbf{r} = \mathbf{r}_1 - \mathbf{r}_2$, $\mathbf{R} = (\mathbf{r}_1 + \mathbf{r}_2)/2$]

$$\langle \mathbf{r}'_1 \mathbf{r}'_2 | V_n^{S_0} | \mathbf{r}_1 \mathbf{r}_2 \rangle = \sum_{\alpha, \beta}^n G_\alpha(r') \lambda_{\alpha\beta} G_\beta(r) \delta(\mathbf{R}' - \mathbf{R}),$$

- Coordinate-space form factor $G_\alpha(r) =$ fourier transform of $g_\alpha(k)$
- Pairing functional

$$\mathcal{E}^{\kappa\kappa} = \sum_q \frac{1}{2} \int d^3\mathbf{R} \sum_{\alpha, \beta=1}^n \check{\rho}_\alpha^{q*}(\mathbf{R}) \lambda_{\alpha\beta} \check{\rho}_\beta^q(\mathbf{R})$$

- One defines **effective** pair densities $\check{\rho}_\alpha^q(\mathbf{R})$ through

$$\check{\rho}_\alpha^q(\mathbf{R}) = \int d^3\mathbf{r} G_\alpha(r) \sum_\sigma (-)^{\frac{1}{2}-\sigma} \kappa^q(\mathbf{R} + \mathbf{r}/2, \sigma; \mathbf{R} - \mathbf{r}/2, -\sigma)$$

- ➡ Incorporate the finite range/non-locality of the interaction
- ➡ Induce non-local pairing field and density
- ➡ BUT the functional depends only on *effective* pair densities *locally* !

EDF calculations in spherical nuclei (2)

- Define reduced two-body wave-functions (spin-singlet part)

$$\check{\Psi}_{ij}^{q\alpha}(\mathbf{R}) \equiv \int d^3\mathbf{r} G_\alpha(r) \Psi_{ij}^q(\mathbf{R} + \mathbf{r}/2, \mathbf{R} - \mathbf{r}/2)$$

$$\Psi_{ij}^q(\mathbf{r}, \mathbf{r}') \equiv \sum_{\sigma} (-)^{s-\sigma} \phi_i(\mathbf{r}, \sigma, q) \phi_j(\mathbf{r}', -\sigma, q).$$

- The ϕ_i are **basis functions** : the $\check{\Psi}_{ij}^{q\alpha}(\mathbf{R})$ are computed **once**
- Build densities and pairing field matrix elements

$$\check{\Delta}_\alpha^q(\mathbf{R}) \equiv \frac{1}{2} \sum_{\beta}^n \lambda_{\alpha\beta} \check{\rho}_\beta^q(\mathbf{R}) \equiv \frac{1}{2} \sum_{\beta}^n \lambda_{\alpha\beta} \sum_{ij} \check{\Psi}_{ij}^{q\beta}(\mathbf{R}) \kappa_{ij}^q$$

$$\Delta_{ij}^q = \sum_{\alpha}^n \int d^3\mathbf{R} \check{\Psi}_{ij}^{q,\alpha}(\mathbf{R}) \check{\Delta}_\alpha^q(\mathbf{R})$$

- ➔ **Pseudo-local pairing problem!**

EDF calculations in spherical nuclei (3)

- New spherical code BSLHFB (T. Lesinski, unpublished)
 - Handle finite-range/non-local forces for systematic calculations
 - Calculations almost as cheap as for a local EDF
 - Basis of spherical bessel functions $j_\ell(kr)$
 - Well suited for drip-line physics

- Calculations
 - Results for 470 nuclei predicted spherical (Gogny-D1S)
 - Pairing complemented with (SLy5) Skyrme EDF ; $m_0^* = 0.7m$
 - $k_{\max} \sim 4.0 \text{ fm}^{-1}$, $R_{box} = 20 \text{ fm}$, $j_{max} = 45/2$

- Comparison of theoretical and experimental pairing gaps

- ✓ Reminder: nothing in the pairing channel is adjusted in nuclei

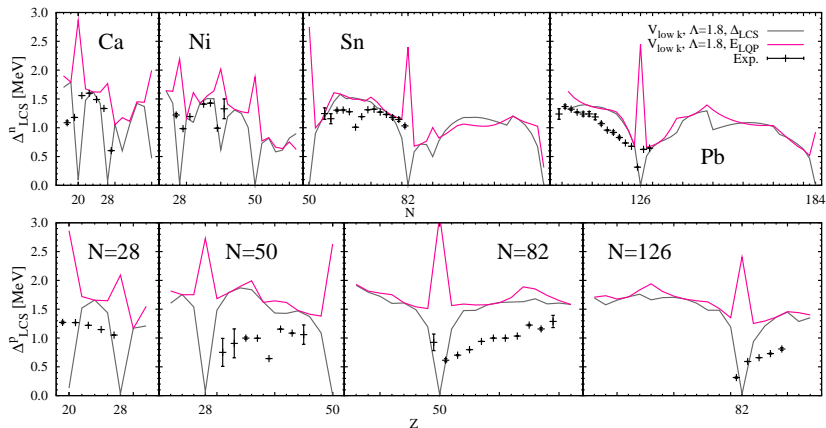
[T. Duguet and T. Lesinski, Eur. Phys. J. Special Topics **156** (2008) 207]

[T. Lesinski, T. Duguet, K. Bennaceur, J. Meyer, arXiv:0809.2895]

Gaps from $v^{pp} = V_{NN}(\Lambda = 1.8)$

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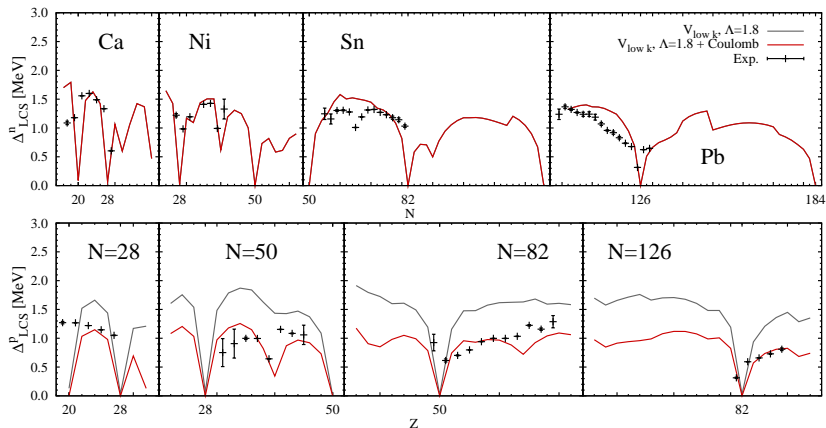


- Neutron gaps Δ^n are consistently close to experimental data
- Proton gaps Δ^p overestimates experimental data systematically

Gaps from $v^{pp} = V_{NN}(\Lambda = 1.8) + V_{Coul,\ell=0}^a$

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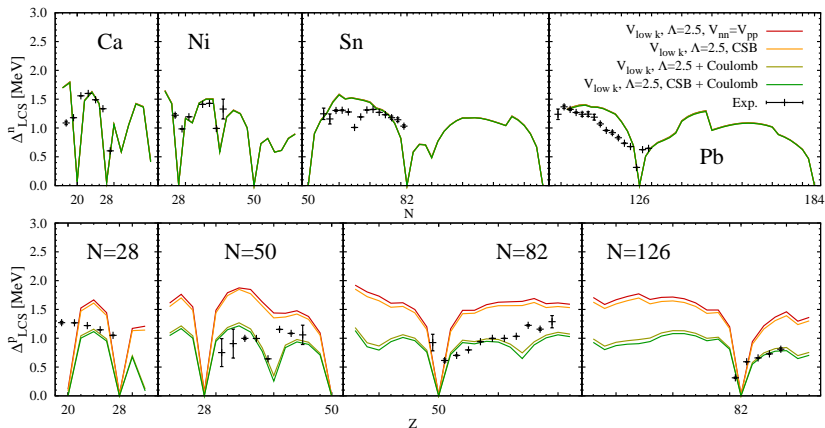


- Coulomb decreases Δ^P by $\sim 40\%$ to bring them close to experiment
- Few masses in the next major shell will be extremely valuable

Gaps from $v^{pp} = V_{NN}(\Lambda = 1.8) + V_{Coul,\ell=0}^a + CSB$

[T. Duguet and T. Lesinski, Eur. Phys. J. Special Topics **156** (2008) 207]

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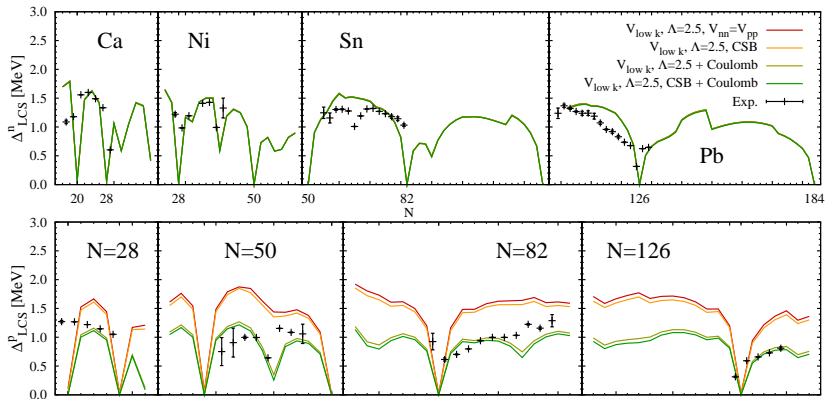


- NN is different in $T_z = \pm 1 \iff CSB$
- Effect of CSB on Δ^P negligible compared to Coulomb

Gaps from $v^{pp} = V_{NN}(\Lambda = 1.8) + V_{Coul,\ell=0}^a + CSB$

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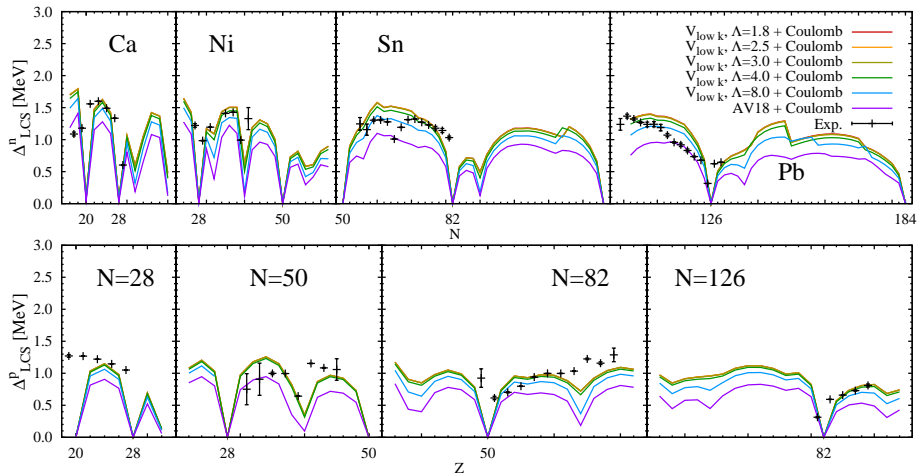


Neutron gap in ^{120}Sn from SLy4 + AV18 - Milan group

- $\Delta_{AV18}^n / \Delta_{low\ k}^n \approx 2/3!$
- Is there an inconsistency? What are the reasons for such a difference?

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Dependence of Δ^q on the RG cut-off Λ 

- Calculations with SLy4 + $V_{NN}(\Lambda)$ for varying Λ

- No variation for $\Lambda \in [1.8, 4] \text{ fm}^{-1}$ but arises for $\Lambda > 4 \text{ fm}^{-1}$ with $\Delta^q \searrow$

Consistency of ph and pp self-energies

The Λ dependence of physical observable characterizes

- Missing pieces in the Hamiltonian one keeps at each Λ
- Correlations missing in the many-body calculation
- Effects of bad approximations at the level one is working at

Fully microscopic calculations in infinite nuclear matter - K. Hebeler et al.

- Different many-body expansions for large and small Λ
 - Results at large and small Λ likely to be different at a given order
 - v^{ph} and v^{pp} must be produced at same Λ
- Impact of using empirical v^{ph} such that $m_r^*(k_F^r) \cong m_r^*(k \approx k_F^r, k_F)$

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 - Low momentum interactions
 - Separable operator representation of $V_{\text{low } k} + V_{\text{Coul}}$
- 3 Pairing gaps in finite nuclei
 - Implementation for finite nuclei calculations
 - Results including nuclear, Coulomb and CSB terms
- 4 Results for soft versus hard NN interactions
 - Dependence of pairing gaps on the RG scale Λ
 - Fully microscopic calculations in infinite nuclear matter
- 5 Summary and outlook

Summary

Pairing gaps in finite nuclei from vacuum NN + Coul.

- Based on low-momentum interactions from RG methods
- First systematic calculations in finite nuclei

First set of results

- Lowest order accounts for the magnitude of experimental gaps
- Coulomb essential for proton gaps ($\sim 40\%$)
- Effects beyond lowest-order seems negligible or cancel each other

Microscopic calculations in SNM and PNM

- **Soft** and **hard** interactions rely on different many-body expansions
- Lowest and higher-order contributions differ in each scheme
- Results for **soft** interaction in finite nuclei confirmed
- Momentum averaging of $m^*(k, k_F)$ reliable for **soft** interaction
- Fine tuning needed to design $m_{Sk}^*(k_F^T)$ for **hard** interaction

Outlook

Works in progress or envisioned

- Extensive study including other observables (T. Lesinski)
- Extension to deformed nuclei (T. Lesinski)
- Fine-tuned Skyrme appropriate to hard NN (T. Lesinski, A. Pastore)
- Equivalent semi-empirical DDDI functional (J. Margueron)
- Incorporate NNN (T. Lesinski)
- Construct ph part (B. Gebremariam)

Introduction
○○○○○○○○

Formalism
○○○○○

Results
○○○○○○○

Consistency of RG scales
○○

Summary

Thank you !

Self energies from many-body expansion

Expansion for soft $V_{NN} = \text{small } \Lambda$

Perturbative

1st order

$$\Sigma^{(1)} = \text{diagram with vertical line, dashed line, and circle}$$

$$\Delta^{(1)} = \text{diagram with vertical line, dashed line, and arc}$$

2nd order ($\Delta^2/\varepsilon_F \rightarrow 0$)

$$\Sigma^{(2)} = \text{diagram 1} + \text{diagram 2} + \text{diagram 3}$$

$$\Delta^{(2)} = \text{diagram 1} + \text{diagram 2}$$

Expansion for hard $V_{NN} = \text{large } \Lambda$

Hole lines/pp-irreducible vertex

1st order

$$\Sigma^{(1)} = \text{diagram with vertical line, wavy line, and circle}$$

$$\Delta^{(1)} = \text{diagram with vertical line, dashed line, and arc}$$

2nd order ($\Delta^2/\varepsilon_F \rightarrow 0$)

$$\Sigma^{(2)} = \text{diagram 1} + \text{diagram 2}$$

$$\Delta^{(2)} = \text{diagram 1} + \text{diagram 2}$$

Self energies from many-body expansion

Expansion for soft $V_{NN} = \text{small } \Lambda$

Perturbative

1st order

$$\Sigma^{(1)} = \text{diagram: a vertical line with a dashed line loop attached to its right side}$$

$$\Delta^{(1)} = \text{diagram: a vertical line with a dashed line loop attached to its bottom side}$$

Expansion for hard $V_{NN} = \text{large } \Lambda$

Hole lines/pp-irreducible vertex

1st order

$$\Sigma^{(1)} = \text{diagram: a vertical line with a wavy line loop attached to its right side}$$

$$\Delta^{(1)} = \text{diagram: a vertical line with a wavy line loop attached to its bottom side}$$

2nd order ($\Delta^2/\varepsilon_F \rightarrow 0$)2nd order ($\Delta^2/\varepsilon_F \rightarrow 0$)

- Second order extendable to all orders in screening bubbles
- Category of diagrams different depending on the starting $V_{NN}(\Lambda)$
- Work here at **1st order** for both soft and hard V_{NN}

First order - soft interaction - Hartree-Fock (HF)

Self energy

$$\text{Basic vertex} = \langle k' | V^{\tau\tau'} J_{IS} | q \rangle$$

$$\Sigma_{\tau}^{(1)}(p, \omega, k_F) = 2 \sum_{\mathbf{q}, \tau'} n(q, k_F^{\tau'}) \left\langle \frac{\mathbf{p} - \mathbf{q}}{2} \middle| V^{\tau\tau'} \middle| \frac{\mathbf{p} - \mathbf{q}}{2} \right\rangle$$

Single-particle energy

$$\varepsilon_p^{\tau} \equiv \frac{p^2}{2} + \text{Re} \Sigma_{\tau}^{(1)}(p)$$

- Non self-consistent problem

First order - hard interaction - Brueckner-Hartree-Fock (BHF)

Self energy

$$\langle k' | G^{\tau\tau'} J_{IS}(P, \omega, k_F) | k \rangle = \langle k' | V^{\tau\tau'} J_{IS} | k \rangle + \frac{2}{\pi} \int q^2 dq \langle k' | V^{\tau\tau'} J_{IS} | q \rangle \cdot \frac{\langle Q^{\tau\tau'}(P, q) \rangle}{\omega - \langle \varepsilon^{\tau\tau'}(P, q) \rangle + i\delta} \langle q | G^{\tau\tau'} J_{IS}(P, \omega) | k \rangle$$

$$\Sigma_{\tau}^{(1)}(p, \omega, k_F) = 2 \sum_{\mathbf{q}, \tau'}^{|\mathbf{q}| < k_F^{\tau'}} \left\langle \frac{\mathbf{p} - \mathbf{q}}{2} \left| G^{\tau\tau'}(|\mathbf{p} + \mathbf{q}|, \omega + \varepsilon_q^{\tau'}) \right| \frac{\mathbf{p} - \mathbf{q}}{2} \right\rangle$$

Single-particle energy

$$\varepsilon_p^{\tau} \equiv \frac{p^2}{2} + \text{Re} \Sigma_{\tau}^{(1)}(p, \varepsilon_p^{\tau})$$

- Self-consistent problem

Effective masses

Effective k-mass and e-mass

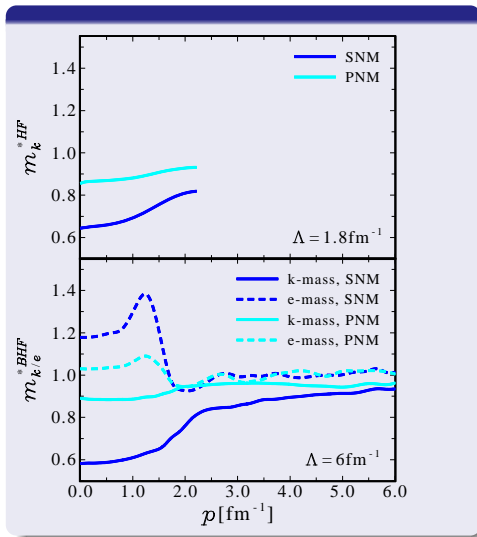
$$\frac{m_{\tau,k}^{*(1)}(p, k_F)}{m} \equiv \left[1 + \frac{1}{p} \frac{\partial \text{Re} \Sigma_{\tau}^{(1)}(\omega, p, k_F)}{\partial p} \Bigg|_{\omega=\varepsilon_p^{\tau}} \right]^{-1}$$

$$\frac{m_{\tau,e}^{*(1)}(p, k_F)}{m} \equiv 1 - \frac{1}{m} \frac{\partial \text{Re} \Sigma_{\tau}^{(1)}(\omega, p, k_F)}{\partial \omega} \Bigg|_{\omega=\varepsilon_p^{\tau}}$$

Total effective mass from $\Sigma_{\Lambda}(\mathbf{k}, \varepsilon_{\mathbf{k}})$

$$\frac{m_{\tau}^{*(1)}(p, k_F)}{m} \equiv \frac{m_{\tau,k}^{*(1)}(p, k_F)}{m} \frac{m_{\tau,e}^{*(1)}(p, k_F)}{m}$$

- $\Sigma_{\text{soft}}^{(1)}$ from soft V_{NN} provides k-mass only
- $\Sigma_{\text{hard}}^{(1)}$ from hard V_{NN} provides both k-mass and e-mass

Momentum dependence of the effective mass, $k_F^{\tau} = 1.2 \text{ fm}^{-1}$ 

Soft - HF

- Mild p -dependence
- Smaller in SNM than in PNM
- Limited to $p \lesssim \Lambda = 1.8 \text{ fm}^{-1}$

Hard - BHF

- e-mass enhancement at k_F^{τ}
- Stronger p -dependence
- Extend to $p \lesssim \Lambda = 6 \text{ fm}^{-1}$
- Larger overall

Reducing the momentum dependence

Remember

Skyrme EDF provides at best $m_{Sk}^{*\tau}(k_F^\tau)$ **independent of momentum**

Averaging procedure of $X = m_\tau^*(p, k_F^\tau)$ or $Z_\tau(p, k_F^\tau)$

- Evaluation on the Fermi surface

$$X_{pe}(k_F^\tau) \equiv X(p = k_F^\tau)$$

- Averaging around the Fermi surface

$$X_{av}(k_F^\tau) \equiv \frac{\int f(q, \Lambda) q^2 dq X(q) u(q, k_F^\tau) v(q, k_F^\tau)}{\int f(q, \Lambda) q^2 dq u(q, k_F^\tau) v(q, k_F^\tau)}$$

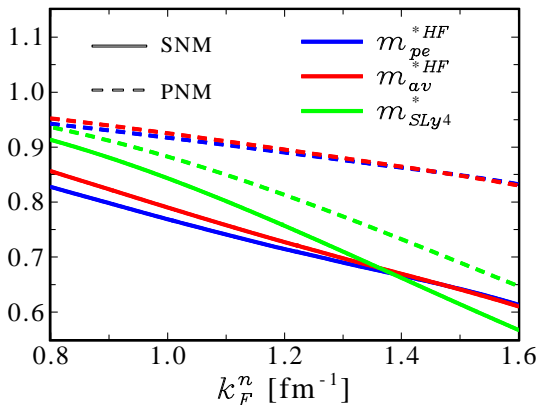
- Other variants

Questions

- Are the results sensitive to the averaging procedure?
- Is there a qualitative difference between **hard** and **soft** interactions?

Momentum independent effective masses for **soft** interaction

[K. Hebeler, T. Duguet, T. Lesinski, A. Schwenk, in preparation)]

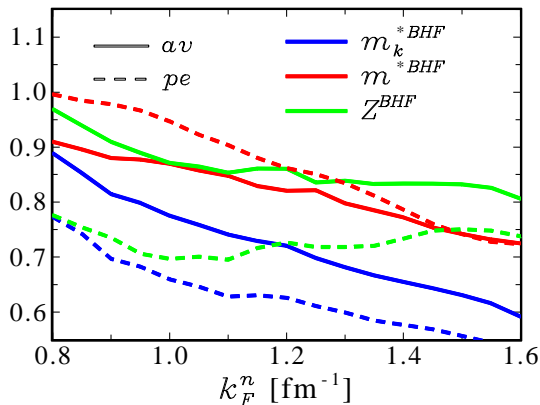


Only slight difference between av and pe at HF level

- SLy4 gives reasonable account of m_k^{*HF} ($k_F^\tau \approx 1.36 \text{ fm}^{-1}$) in SNM
- Wrong isovector m_1^* of SLy4 [T. Lesinski et al. 2006)]

Momentum independent effective masses in SNM for **hard** interaction

[K. Hebeler, T. Duguet, T. Lesinski, A. Schwenk, in preparation]

Difference between av/pe much larger than for soft cutoff interaction

- Momentum dependence stronger
- Larger momentum-space for averaging procedure

Pairing gaps

Gap equation

- After pole approximation

$$\hat{\Delta}(\mathbf{k}) = - \int \frac{d^3\mathbf{k}'}{(2\pi)^3} \frac{Z(\mathbf{k}) V_{NN}(\mathbf{k}, \mathbf{k}'; \Lambda) Z(\mathbf{k}') \hat{\Delta}(\mathbf{k}')}{2\sqrt{(\varepsilon_{\mathbf{k}'} - \mu)^2 + \hat{\Delta}^2(\mathbf{k}')}}}$$

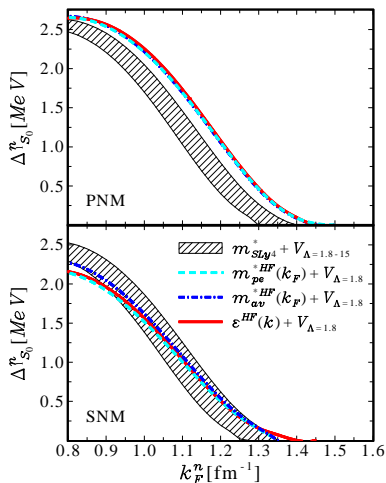
- $\hat{\Delta}(\mathbf{k}) = Z(\mathbf{k}) \Delta(\mathbf{k}) =$ physical gap of the excitation spectrum
- Effective mass approximation relates to $\varepsilon_{\mathbf{k}'} - \mu$ in the denominator

Questions of interest regarding results in finite nuclei

- Impact of using $m_{\tau}^*(k, k_F) \approx m_{\tau}^*(k_F^T)$?
- Is there a qualitative difference between **hard** and **soft** interactions?

Pairing gaps from **soft** interactions - PNM and SNM

[K. Hebeler, T. Duguet, T. Lesinski, A. Schwenk, in preparation]

Gaps from full $m_{\tau}^{*HF}(k, k_F)$

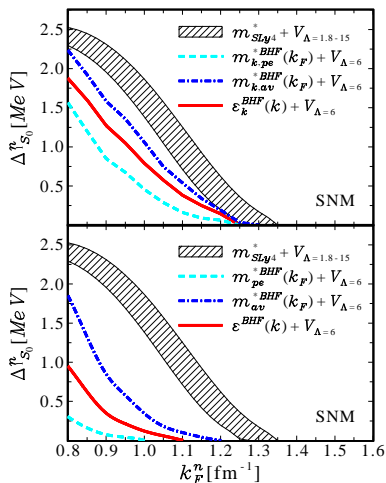
- Close to SLy4 + $V_{\Lambda=1.8}$ for $k_F^n \in [1.1, 1.4]$
- Trace of wrong m_1^* of SLy4 in PNM

Gaps from $m_{\tau}^{*HF}(k, k_F) \approx m_{\tau}^*(k_F^T)$

- Reproduce well gaps from $m_{\tau}^{*HF}(k, k_F)$
- No sensitivity to averaging procedure
- Variation \ll bandwidth of SLy4 calc.
- Could optimize $m_{Sk}^*(k_F^T)$

Pairing gaps from **hard** interactions - SNM

[K. Hebeler, T. Duguet, T. Lesinski, A. Schwenk, in preparation]

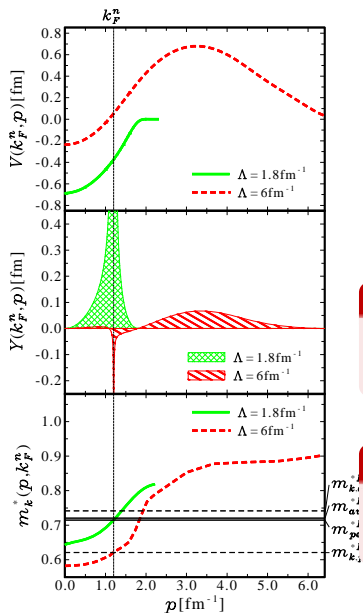
Gaps from full $m_{\tau,e/k}^{*BHF}(k, k_F)$

- Lower than SLy4 + $V_{\Lambda=15}$ for all k_F^n
- Very small if considering $Z_\tau(k, k_F)$

Gaps from $m_{\tau,e/k}^{*BHF}(k, k_F) \approx m_\tau^*(k_F^T)$

- Strong sensitivity to averaging procedure
- Variation \gtrsim bandwidth of SLy4 calc.
- Fine tuning needed to design $m_{Sk}^*(k_F^T)$
→ Collaboration with Milan group

Analysis



- Consider matrix elements $V_{\Lambda=1.8/6}(k_F^n, p)$

- Write gap equation schematically as

$$\hat{\Delta}(k_F^n) \equiv \int dq Y(k_F^n, q)$$

Gap generated

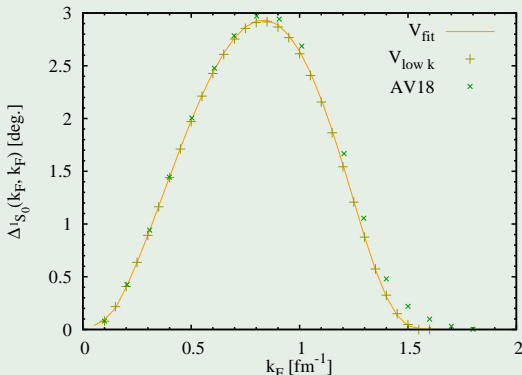
- Around the Fermi surface for **soft** Λ
- Mainly at large momenta for **hard** Λ

Effect of $m_\tau^*(k_F^\tau) = \text{constant}$ - 'pe' values here

- Good approx for **soft** Λ around k_F^n
- Bad approx for **hard** Λ at relevant $p \gtrsim 2 \text{ fm}^{-1}$

$\Delta_{1S_0}^n(k_F, k_F, \Lambda)$ in INM from $V_{\text{low } k}(\Lambda)$

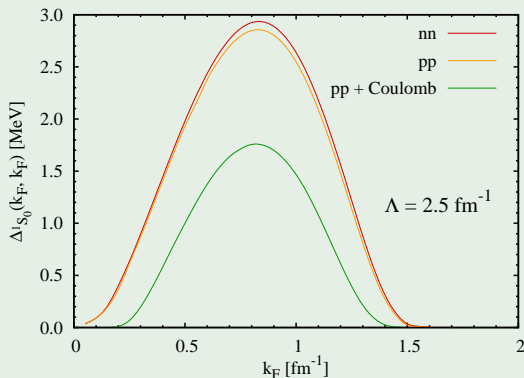
Neutron gaps with $\epsilon_k = \hbar^2 k^2 / 2m$ ($m^* = m$)



- Gaps of numerical $V_{\text{low } k}$ reproduced perfectly without direct fitting
- True for $\Delta_{1S_0}^n(k, k_F, \Lambda) \quad \forall k \lesssim \Lambda$
- $\partial_\Lambda \Delta_{1S_0}^n(k_F, k_F, \Lambda) = 0$ [K. Hebeler, A. Schwenk, B. Friman, 2007]

$$\Delta_{1S_0}^q(k_F, k_F, \Lambda) \text{ in INM } V_{\text{low } k}(\Lambda) + V_{\text{Coul}}$$

Neutron/proton gaps with $\epsilon_k = \hbar^2 k^2 / 2m$ ($m^* = m$)



- Coulomb reduces proton gaps by $\sim 40\%$
- Comparatively CSB effects are negligible

Effective masses

Effective k-mass, e-mass and total mass from $\Sigma_{\Lambda}(\mathbf{k}, \varepsilon_{\mathbf{k}})$

$$\frac{m_{\tau}^{*(1)}(p, k_F)}{m} \equiv \frac{m_{\tau, k}^{*(1)}(p, k_F)}{m} \frac{m_{\tau, e}^{*(1)}(p, k_F)}{m}$$

- $\Sigma_{\text{soft}}^{(1)}$ from **soft** V_{NN} provides **k-mass only**
- $\Sigma_{\text{hard}}^{(1)}$ from **hard** V_{NN} provides **both k-mass and e-mass**

Remember

Skyrme EDF provides at best $m_{Sk}^{*\tau}(k_F^{\tau})$ **independent of momentum**

Momentum-dependence averaging $m_{\tau}^{*}(k, k_F) \approx m_{\tau}^{*}(k_F^{\tau})$

- Result **insensitive** to procedure for **soft** V_{NN}
- Result **very sensitive** to procedure for **hard** V_{NN}

Reducing the momentum dependence

Remember

Skyrme EDF provides at best $m_{Sk}^{*T}(k_F^T)$ **independent of momentum**

Averaging procedure of $X = m_\tau^*(p, k_F^T)$ or $Z_\tau(p, k_F^T)$

- Evaluation on the Fermi surface

$$X_{pe}(k_F^T) \equiv X(p = k_F^T)$$

- Averaging around the Fermi surface

$$X_{av}(k_F^T) \equiv \frac{\int f(q, \Lambda) q^2 dq X(q) u(q, k_F^T) v(q, k_F^T)}{\int f(q, \Lambda) q^2 dq u(q, k_F^T) v(q, k_F^T)}$$

- Other variants

Results

- Provides exact same $m_\tau^*(k_F^T)$ for **soft** interaction
- Result very sensitive to procedure for **hard** interaction

Pole approximation to the gap equation

General gap equation with $\Gamma^{irr} = V$

$$\Delta(\mathbf{k}) = \int \frac{d^3\mathbf{k}'}{(2\pi)^3} \int \frac{d\omega'}{2\pi i} V(\mathbf{k}, \mathbf{k}') F(\mathbf{k}', \omega')$$

- Neglect imaginary part of Σ
- Find pole $E_{\mathbf{k}}$ by solving $F^{-1}(\mathbf{k}, \omega) = 0$
- Expand propagator around these poles
- Perform the energy integral analytically in the limit $\Delta \ll \varepsilon_F$

$$\Delta(\mathbf{k}) = - \int \frac{d^3\mathbf{k}'}{(2\pi)^3} \frac{V(\mathbf{k}, \mathbf{k}') Z(\mathbf{k}') \Delta(\mathbf{k}')}{2\sqrt{[\varepsilon_{\mathbf{k}'}^0 - \mu + \frac{1}{2} [\Sigma(\mathbf{k}', \varepsilon_{\mathbf{k}'}^0) + \Sigma(\mathbf{k}', 2\mu - \varepsilon_{\mathbf{k}'}^0)]]^2 + \Delta^2(\mathbf{k}')}}}$$

with $Z(\mathbf{k}) = m_e^{-1}(\mathbf{k})$.

Reference: Definition of Fermi momenta

$$\begin{aligned}k_{\text{F}}^p &= \left[\frac{3\pi^2}{2}(1-\beta)\rho \right]^{1/3} \\k_{\text{F}}^n &= \left[\frac{3\pi^2}{2}(1+\beta)\rho \right]^{1/3} \\ \beta &= (\rho_n - \rho_p)/\rho \\ \rho &= \rho_n + \rho_p = \frac{1}{3\pi^2} [(k_{\text{F}}^n)^3 + (k_{\text{F}}^p)^3]\end{aligned}$$

In SNM we have $k_{\text{F}}^p = k_{\text{F}}^n \equiv k_{\text{F}} = [3\pi^2 \rho_q]^{1/3}$.