

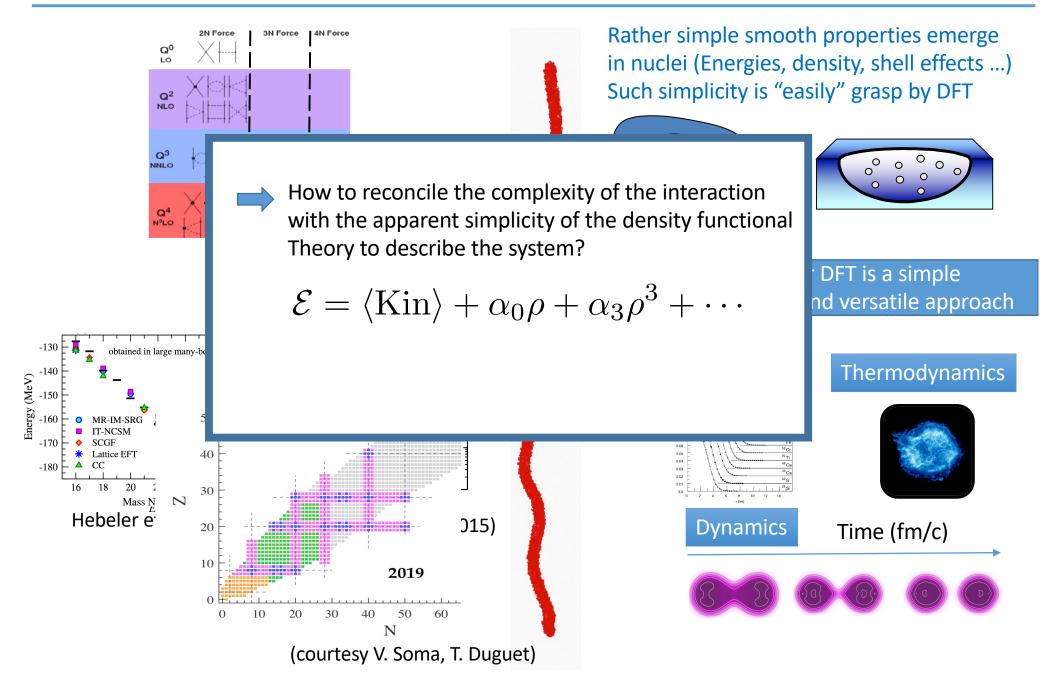
Fermi systems with large s-wave scattering length: cold atoms and nuclear matter

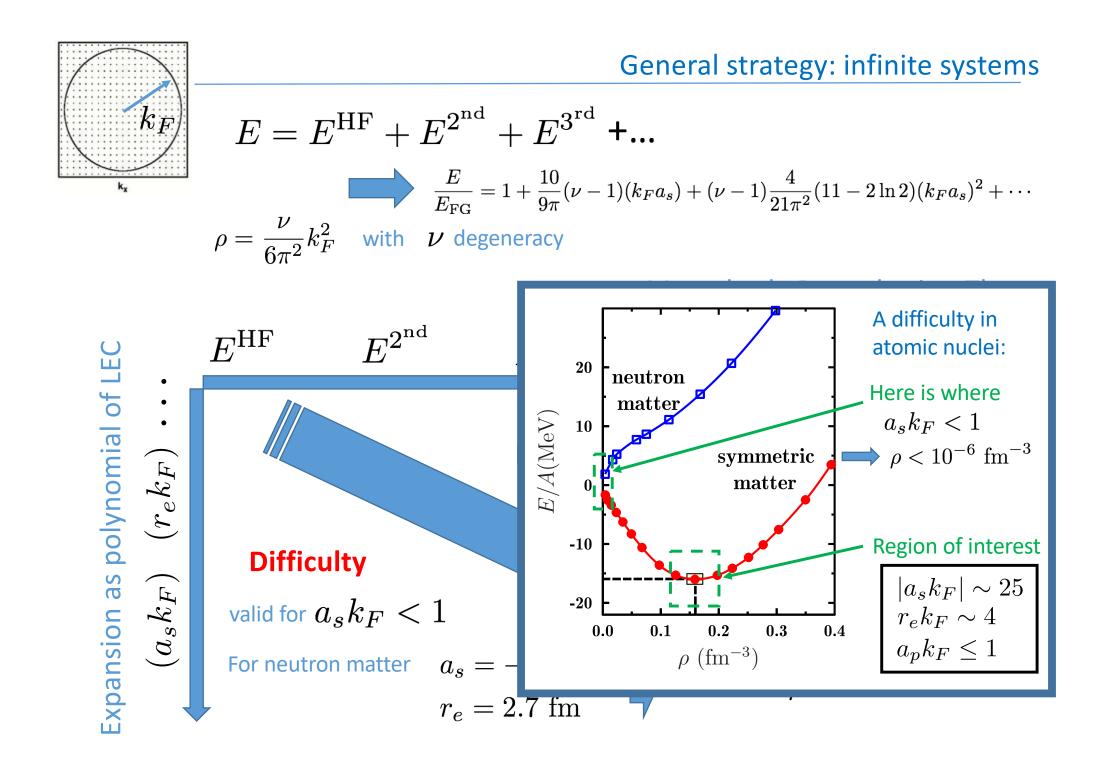
Denis Lacroix (IJCLab, Orsay, France)

Outline:

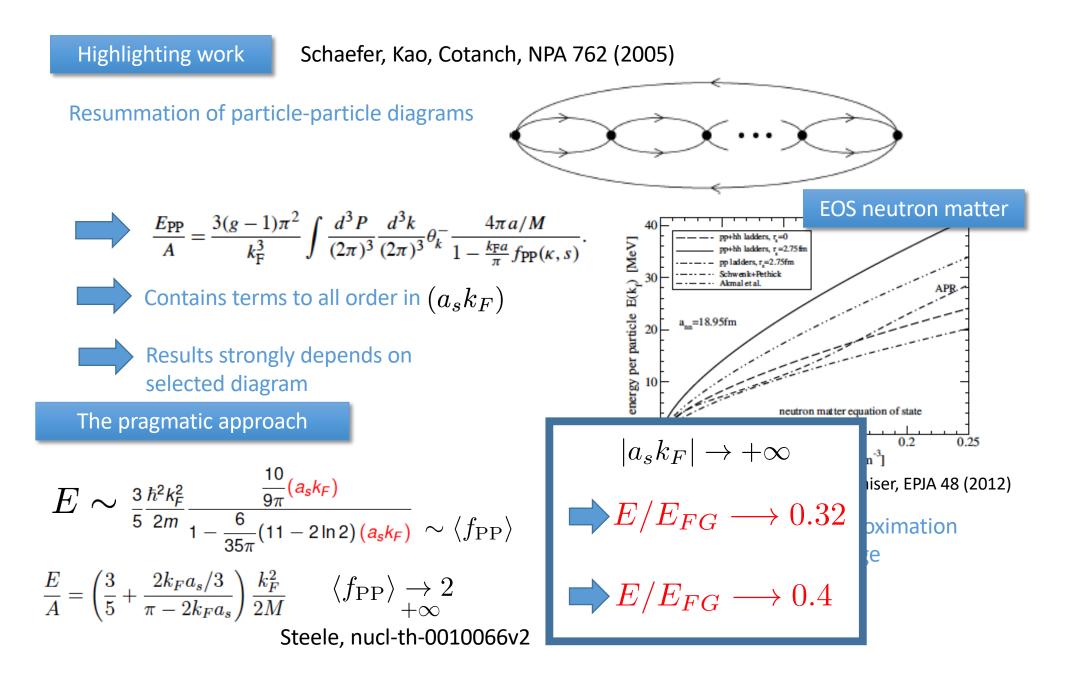
- Brief discussion on Density functional theory for nuclei
- Common DFT for nuclei and cold atoms?
- Applications: energies, static response, ...
- Self-energy: quasi-particle properties.

Bare int.+Many-Body calculation versus Density Functional theory

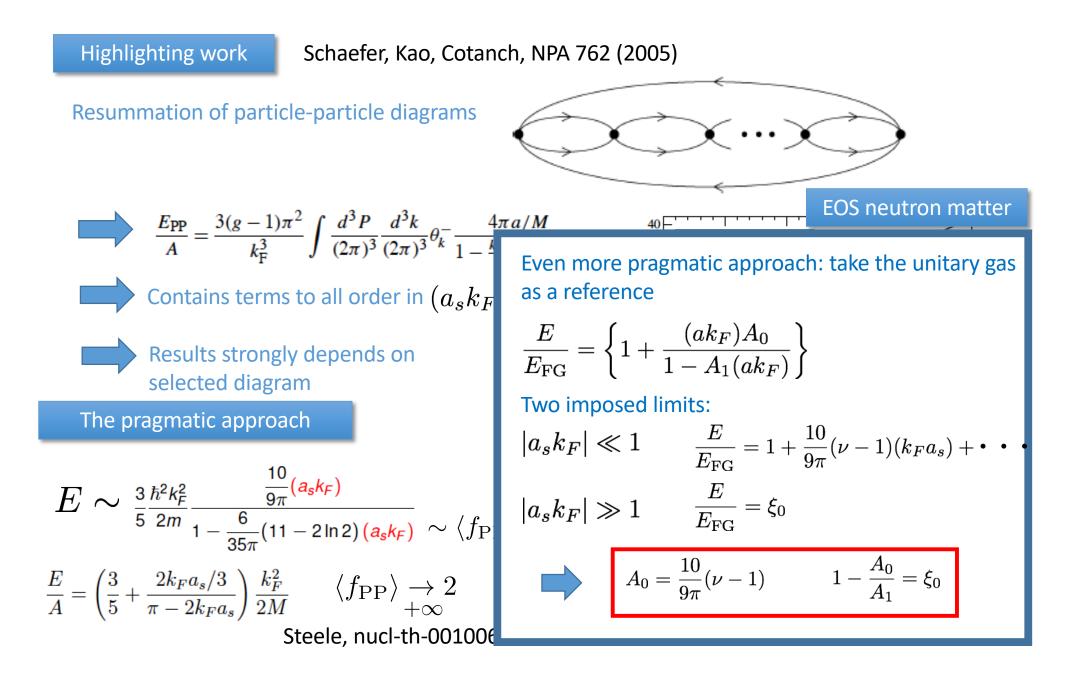




The "magic" technique: resummation and phase-space argument

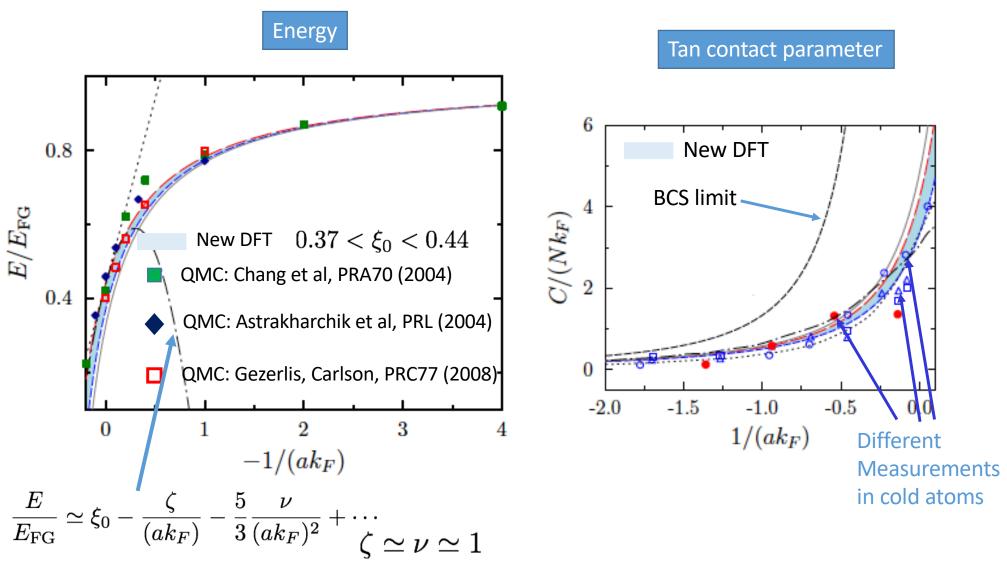


The "magic" technique: resummation and phase-space argument



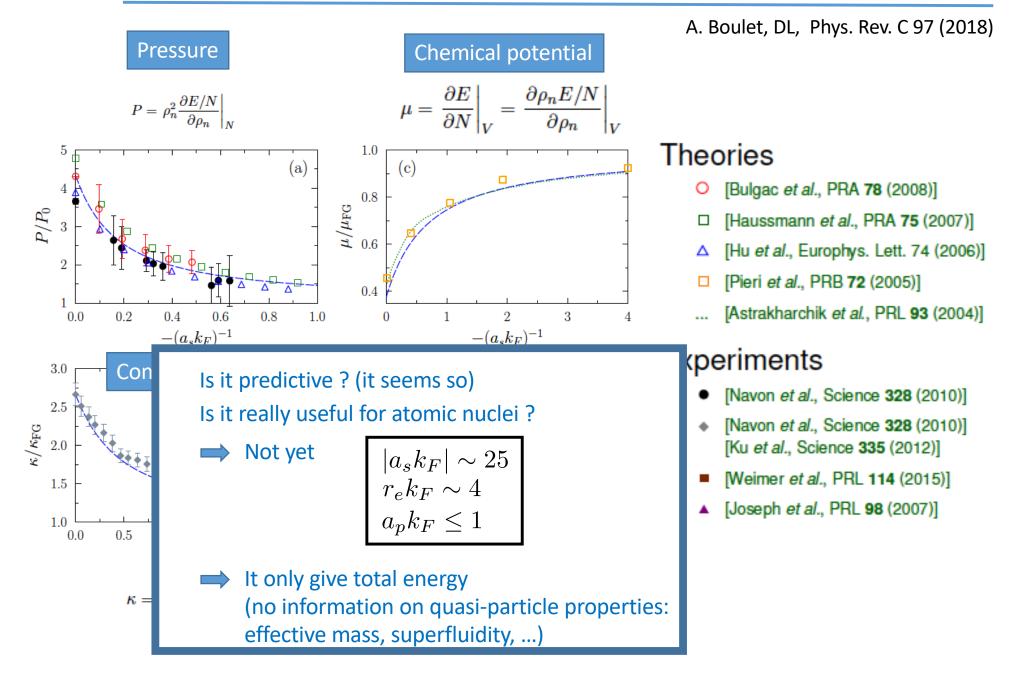
Result of the DFT for at or close to unitarity

Lacroix, PRA 94 (2016)

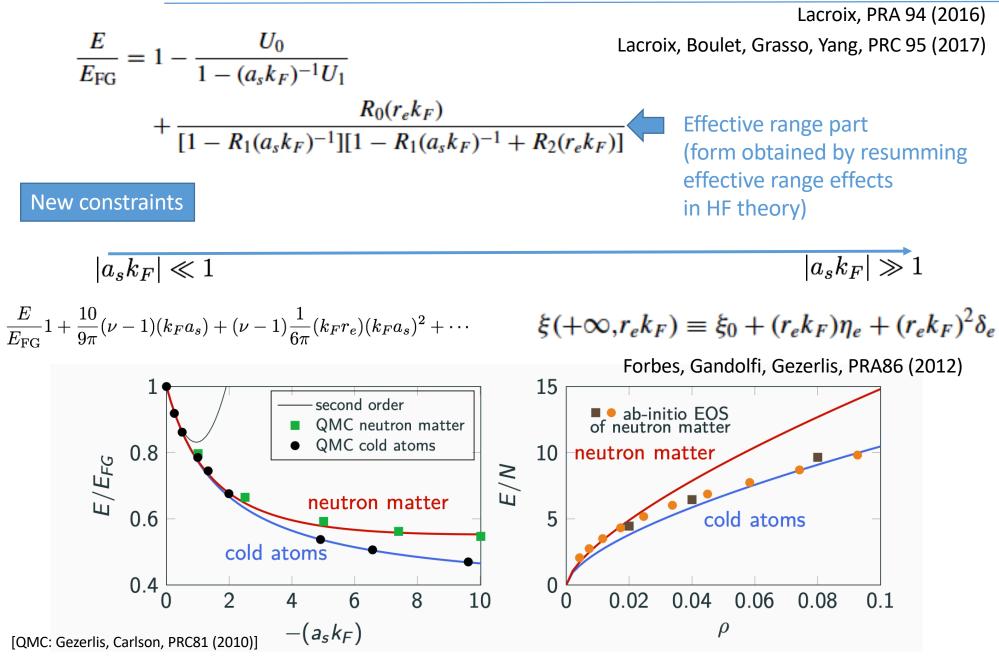


Taylor expansion in $(a_s k_F)^{-1}$: Bulgac and Bertsch, PRL 94 (2005)

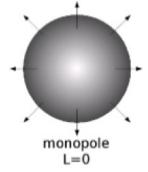
Example of applications: thermodynamical quantities around unitarity



From cold atom to neutron matter: inclusion of effective range



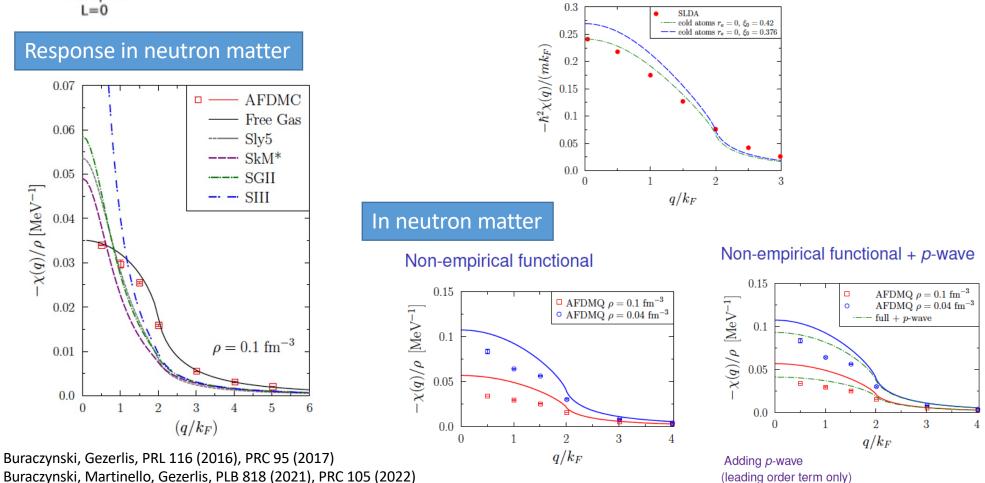
Some successful application



Static response in cold atoms and neutron matter

A. Boulet, DL, Phys. Rev. C 97 (2018) Static response In unitary gas

SLDA: [Forbes and Sharma, PRA 90 (2014)]



Buraczynski, Martinello, Gezerlis, PLB 818 (2021), PRC 105 (2022)

F. 1

$$\frac{L_p}{E_{\rm FG}} = \frac{1}{\pi} (a_p k_F)^3$$

What have we learn beyond having a unitary gas guided DFT?

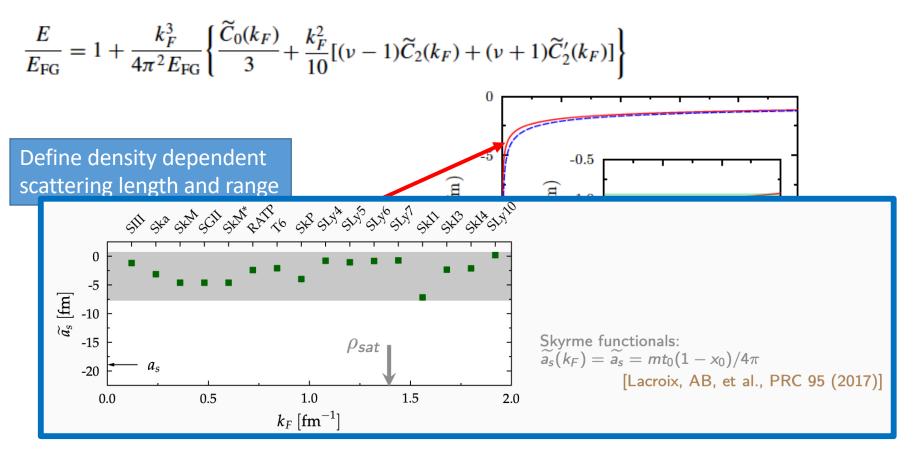
How to conceal the complexity of the interaction with the apparent simplicity of the density functional Theory to describe the system?

 $\mathcal{E} = \langle \mathrm{Kin} \rangle + \alpha_0 \rho + \alpha_3 \rho^3 + \cdots$

Restart from the functional

$$\frac{E}{E_{\rm FG}} = 1 - \frac{U_0}{1 - (a_s k_F)^{-1} U_1} + \frac{R_0(r_e k_F)}{[1 - R_1(a_s k_F)^{-1}][1 - R_1(a_s k_F)^{-1} + R_2(r_e k_F)]}$$

Rewrite it as



How to conceal the complexity of the interaction The simple DFT has guided us to design new functional theory with the apparent simplicity of the density functional Theory to describe the system? valid at low density up to densities of interest. $\mathcal{E} = \langle \mathrm{Kin} \rangle + \alpha_0 \rho + \alpha_3 \rho^3 + \cdots$ The YGLO functional The ELYO functional *YGLO : Yang Grasso Lacroix Orsay *ELYO : Extended Lee-Yang Orsay $\frac{E}{A} = K_{\beta} + \frac{B_{\beta}\rho}{1 - R_{\beta}\rho^{1/3} + C_{\beta}\rho^{2/3}} + D_{\beta}\rho^{5/3} + F_{\beta}\rho^{\alpha+1}$ $\frac{E}{N} = \frac{\hbar^2 k_F^2}{2m} \left[\frac{3}{5} + \frac{2}{3\pi} (k_F a) + \frac{4}{35\pi^2} (11 - 2\ln 2) (k_F a)^2 \right]$ $\rho(\text{fm}^{-3})$ 0.95 Neutron 0.90 0.005 0.0462 0.135 $+\frac{1}{10\pi}(k_F r_s)(k_F a)^2+0.019(k_F a)^3$, $E/\hbar\omega N^{4/3}$ drops 0.85 ····· Lee-Yang (resum.) SLv5 0.80 1.0· — · YGLO (Akmal) QMC AV4 0.75 - YGLO (FP) □ AFDMC 0.70 O HS $E_{NM}/E_{
m FG}$ 0 0.64 • FP 14 20 26 32 Akmal Bonnard, Grasso, lacroix, PRC98 (2018), PRC101 (2020) Atomic Nuclei a (fm) $-a(k_{\rm E}) k_{\rm E} = 0.5$ $- a(k_{\rm E}) k_{\rm E} = 1$ 0.4 (a) VGLO(FP) Yang 2016, Bonnard 2018 ← YGLO(FP) Yang 2016 + this work ¥YGLO(Akmal) Yang 2016 + this work 35 10 200 30 SLv5 30 $|ak_N|$ 25 % [WeV] (a) (b) 0.24 0 Yang, Grasso, Lacroix PRC94 (2016) 0.08 0.16 0.5 1.0 1.5 2.0 Density (fm⁻³) $k_{F}(fm^{-1})$ 15 Grasso, Lacroix, Yang, PRC 95 (2017) O isotopic chain

Burrello, Bonnard, Grasso, PRC103 (2021)

10 12 14 16 18 20 22

4 6

SkM*

KIDS

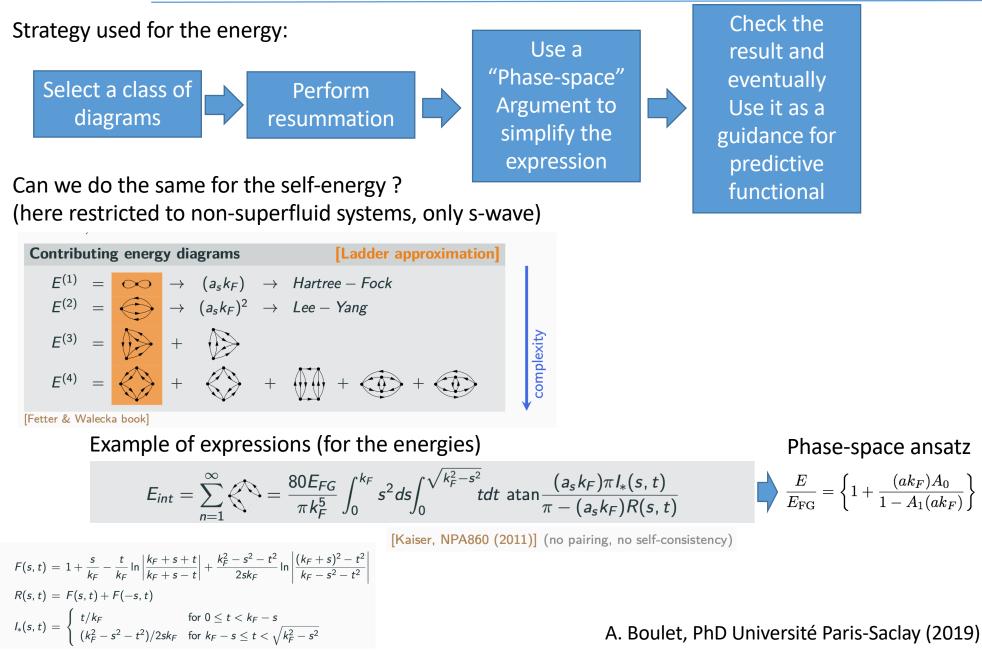
ELYO

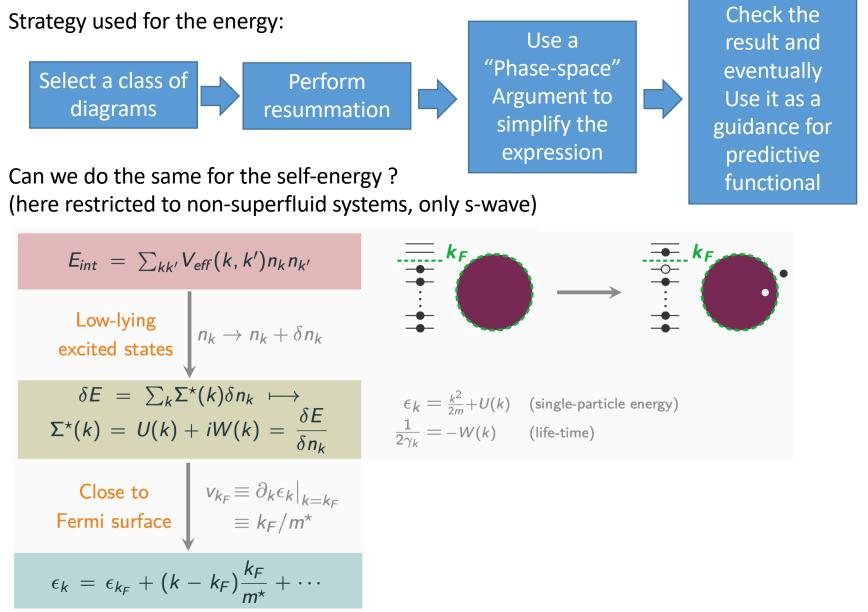
38

44 50

ah initic

UNEDF0 YGLO

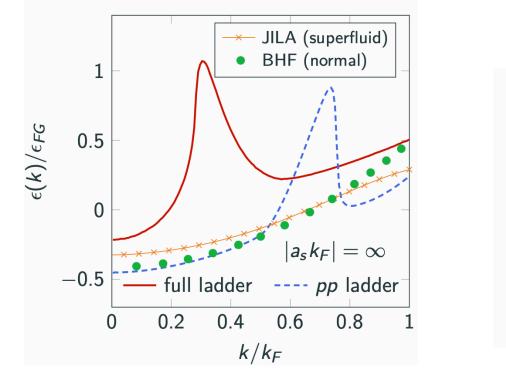




A. Boulet, PhD Université Paris-Saclay (2019)

Exact formula of the self-energy after resummation:

$$\Sigma^{\star}(k) = U(k) + iW(k) \qquad U(k < k_F) = \frac{8}{m\pi^2} \int_0^{k_F} s^2 ds \int_0^{\sqrt{k_F^2 - s^2}} t dt \, \mathcal{U}(s, t, k < k_F)$$
[Kaiser, EPJA49 (2013)]

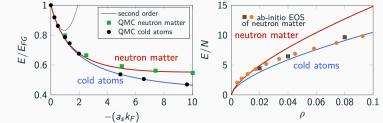




Performing a "phase-space" like average does not seem so evident to us

We used a different strategy

We have developed rather simple Density functional theory that applies to both cold atoms and nuclei

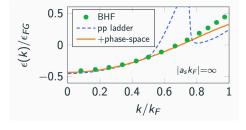


These functionals seems useful to understand the apparent simplicity of other historical successful functionals.

This has also initiated new types of functionals that respects low density limit

The problem is much harder on self-energy (without superfluidity and selected

diagrams)



Local energy density functional for superfluid Fermi gases from effective field theory

Antoine Boulet¹,¹,^{*} Gabriel Wlazłowski¹,²,[†] and Piotr Magierski¹,²,[‡] ¹Faculty of Physics, Warsaw University of Technology, Ulica Koszykowa 75, 00-662 Warsaw, Poland ²Department of Physics, University of Washington, Seattle, Washington 98195–1560, USA (Dated: January 20, 2022)

arXiv:2201.07626