

Clusters in hot and dense stellar matter

Francesca Gulminelli - LPC Caen, France

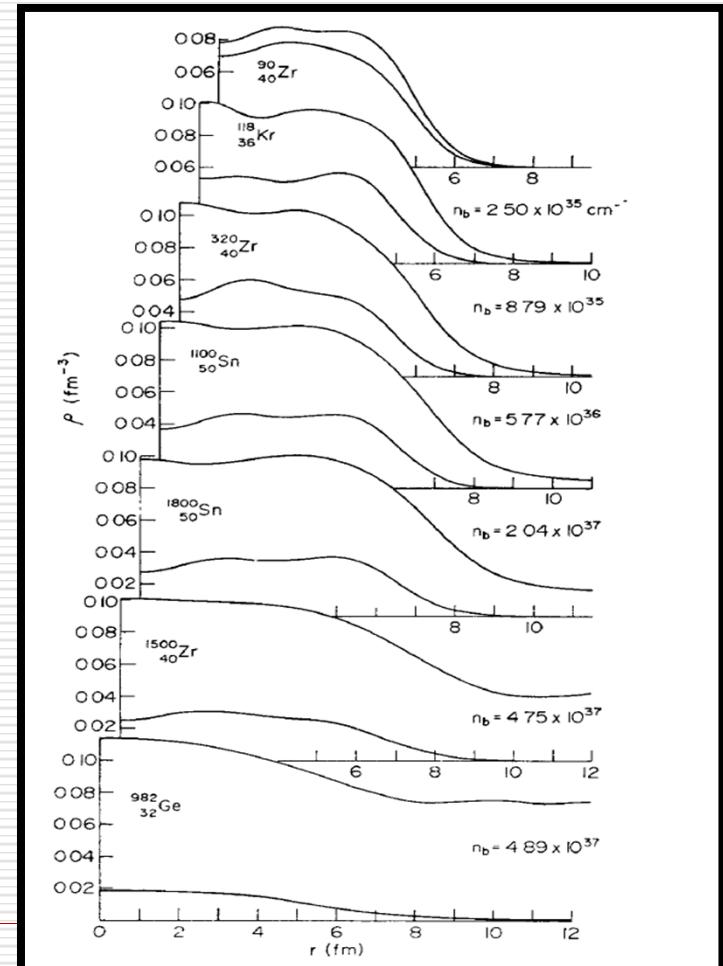
- Motivation: a realistic EoS for core collapse simulations
- A description of stellar matter with cluster dof
- Effect of the effective interaction and in-medium self energies



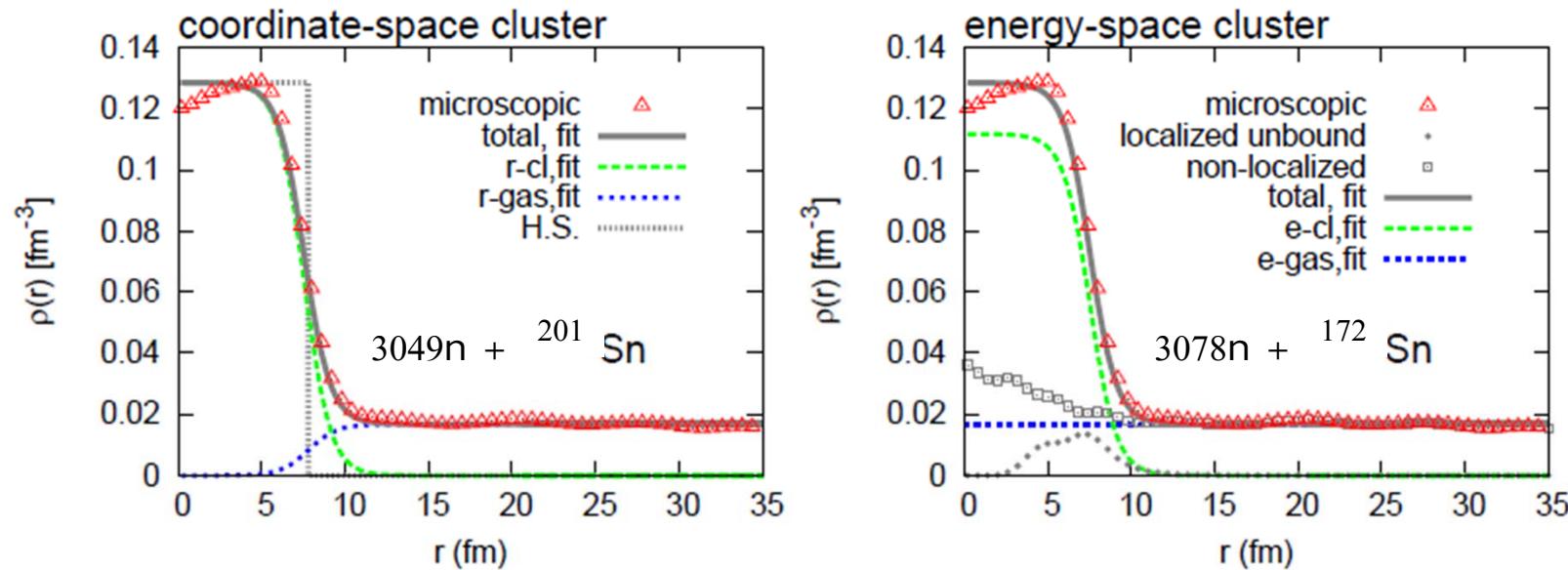
Nuclear matter at subsaturation is clusterized

J. W. Negele and D. Vautherin,
Nuclear Physics A 207, 298 (1973)

- Below nuclear saturation matter is frustrated:
- Nuclear and Coulomb forces act on comparable length scales
- The LG transition is quenched and matter is clusterized.
- A description in terms of nucleons is feasible (but heavy) in DFT
- An extension to mixed states (beyond MF and/or T>0) is (almost) hopeless
- An effective model with cluster DoF is appealing



Clusters in a dense medium



Cluster: density fluctuation

- =>excluded volume applies
- =>bulk energy \sim vacuum energy
(zero order LD approximation)

Cluster: localized wave functions

- =>excluded volume does not apply
- =>binding energy shift

A model for the energy density

- $\varepsilon_{HM}(\vec{x}_g), B_0(\vec{x}_A)$

« Best » functionals for homogeneous matter and for a cluster in the vacuum



$$\begin{aligned} V_{WS}\varepsilon_{WS}[\hat{\rho}_n, \hat{\rho}_p, \hat{\rho}_e, V_{WS}] &= B_0(\vec{x}_A) + V_{WS}\varepsilon_{HM}(\vec{x}_g) + \delta B \\ &= B_m(\vec{x}_g, \vec{x}_A) + V_{WS}\varepsilon_{HM}(\vec{x}_g) \end{aligned}$$

$B_m \approx B_0(\vec{x}_A) + V_{WS}\delta\varepsilon_c(\vec{x}_A, \rho_p) - \varepsilon_{HM}(\vec{x}_g)V(\vec{x}_A)$ *in-medium binding energy shift*

- Numerical applications:

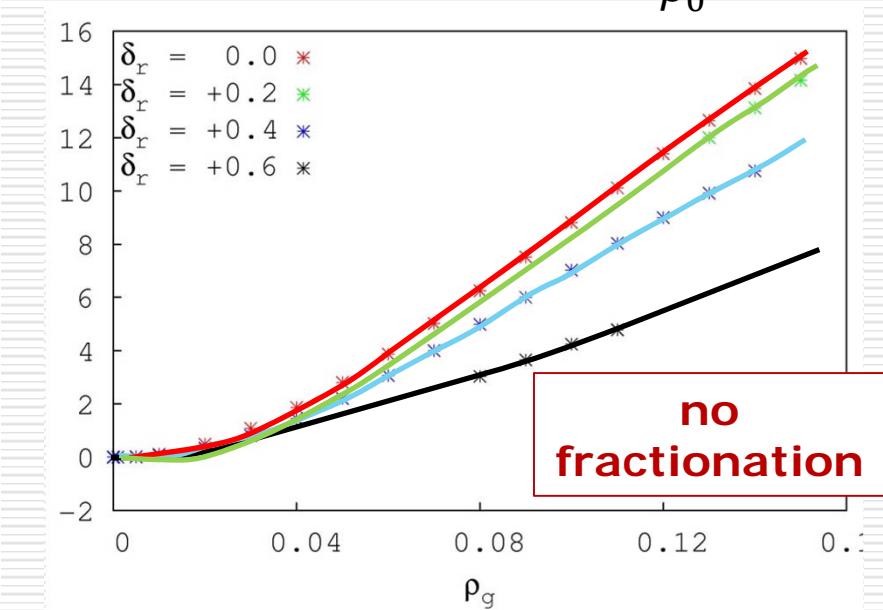
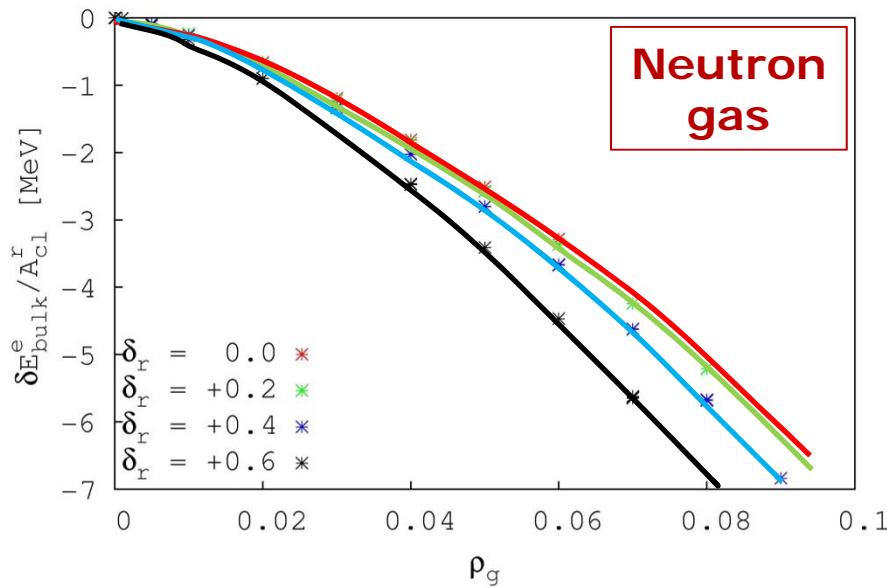
$e_{HM}(\rho_g, \rho_{Ig})$: Homogeneous matter Skyrme energy functional

$B_0(A, I)$: Mass formula from Skyrme-HF

In-medium bulk energy shift

Sly4 Skyrme functional

$$\delta = \frac{(\rho_{0n} - \rho_{0p})}{\rho_0}$$



Cluster DoF: T=0

- A model for the energy density

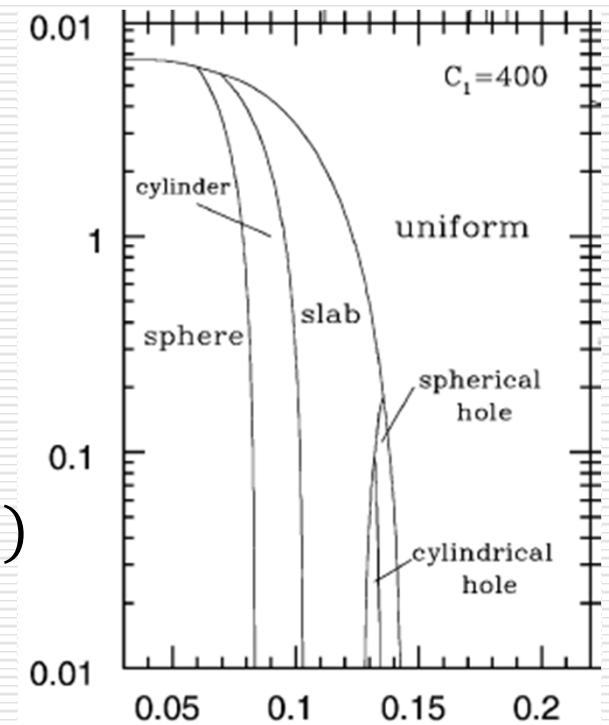
$$V_{WS}\varepsilon_{WS}[\hat{\rho}_n, \hat{\rho}_p, \hat{\rho}_e, V_{WS}] = B_m(\vec{x}_g, \vec{x}_A) + V_{WS}\varepsilon_{HM}(\vec{x}_g)$$

$$\vec{x}_A = (A, I, V_{WS} \dots) \quad \vec{x}_g = (\rho_g, \rho_{Ig} \dots)$$

Variational variables

- A variational problem for each (ρ_B, ρ_I)

$$d \left(\varepsilon_{WS}(\vec{x}_A, \vec{x}_g) - \mu \left(\frac{A_{WS}}{V_{WS}} - \rho_B \right) - \mu_I \left(\frac{I_{WS}}{V_{WS}} - \rho_I \right) \right) = 0$$



Cluster DoF: T=0

- Equilibrium equations $\vec{x}_A = (A, \delta, V_{WS})$ $\vec{x}_g = (\rho_g, \rho_{Ig})$

$$\left\{ \begin{array}{l} \frac{\partial B_m}{\partial A} = \mu_g \left(1 - \frac{\rho_g}{\rho_0} \right) - \mu_{Ig} \left(\delta - \frac{\rho_{Ig}}{\rho_0} \right) \\ \frac{1}{A} \frac{\partial B_m}{\partial \delta} = \mu_{Ig} + \frac{1}{\rho_0^2} \frac{\partial \rho_0}{\partial \delta} (\mu_g \rho_g + \mu_{Ig} \rho_{Ig}) \\ A \frac{\partial B/A}{\partial A} = \frac{\delta B}{V_{WS}} \end{array} \right.$$

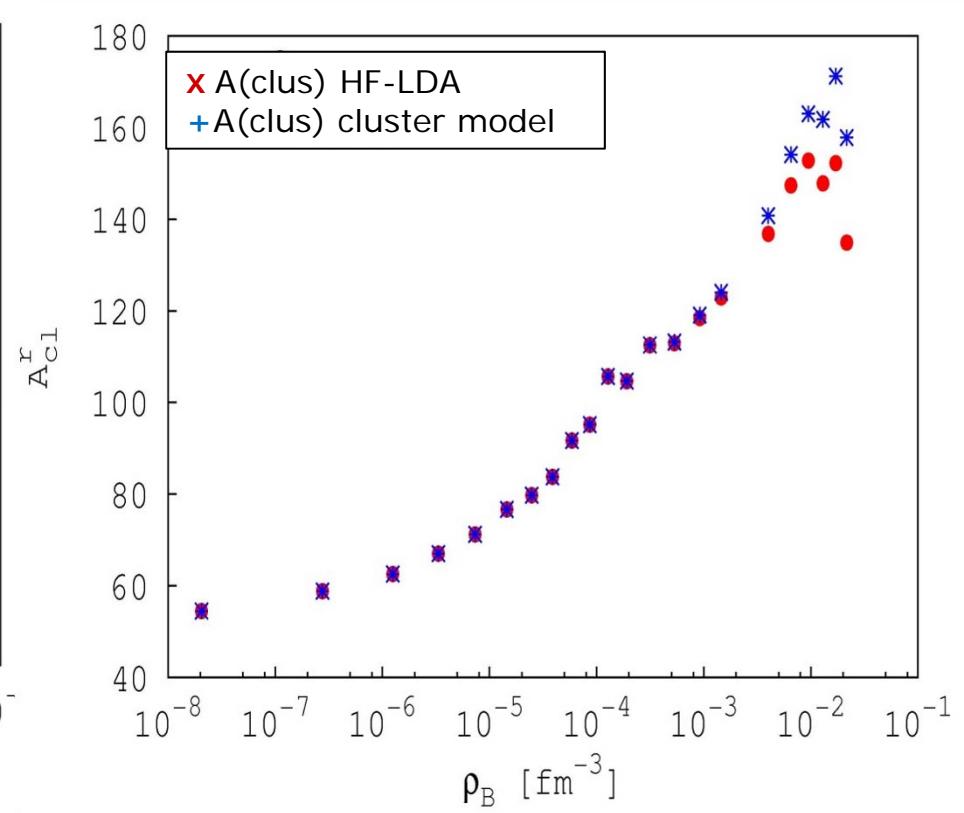
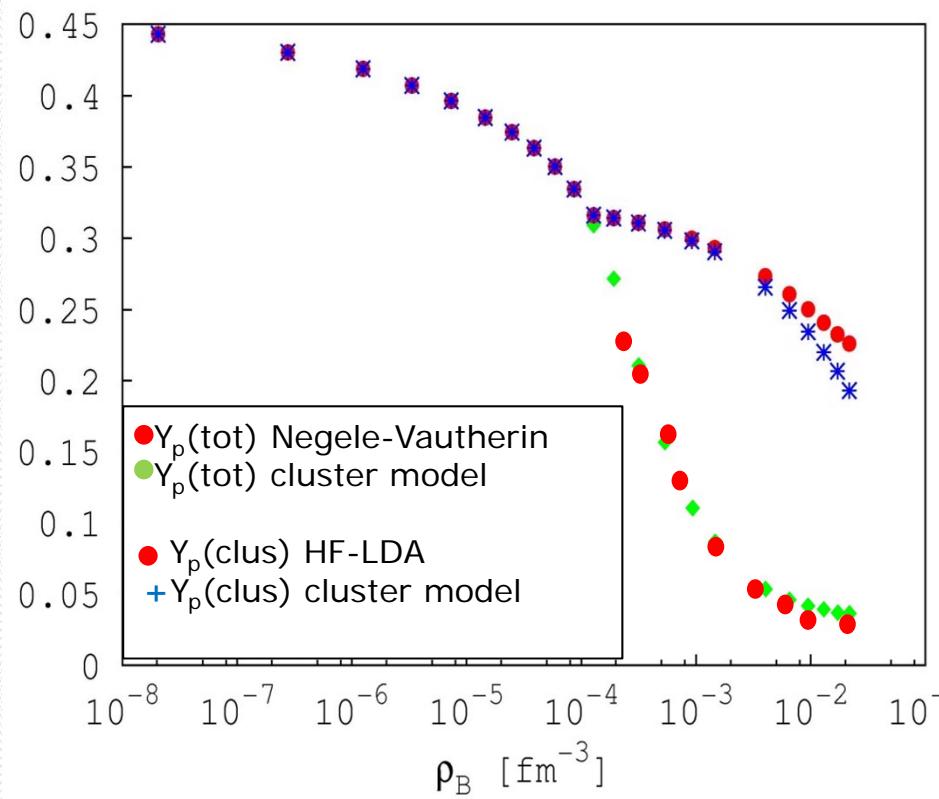
$$\mu_g = \frac{\partial \varepsilon_{HM}}{\partial \rho_g}$$

$$\mu_{Ig} = \frac{\partial \varepsilon_{HM}}{\partial \rho_{Ig}}$$

$$\delta = \frac{(\rho_{0n} - \rho_{0p})}{\rho_0}$$

Cluster DoF: T=0

Sly4 Skyrme functional



T>0: the SN approximation

- A model for the **free** energy density

$$V_{WS}f_{WS}[\hat{\rho}_n, \hat{\rho}_p, \hat{\rho}_e, V_{WS}, \mathbf{T}] = B_m(\vec{x}_g, \vec{x}_A) - \mathbf{TS}(\vec{x}_A, \vec{x}_g) + V_{WS}f_{HM}(\vec{x}_g, \mathbf{T})$$

$$\vec{x}_A = (A, I, V_{WS} \dots) \quad \vec{x}_g = (\rho_g, \rho_{Ig} \dots)$$

Variational variables

- A variational problem for each $(\rho_B, \rho_I, \mathbf{T})$

$$d \left(f_{WS}(\vec{x}_A, \vec{x}_g) - \mu \left(\frac{A_{WS}}{V_{WS}} - \rho_B \right) - \mu_I \left(\frac{I_{WS}}{V_{WS}} - \rho_I \right) \right) = 0$$

J. M. Lattimer and F. D. Swesty, NPA 535, 331 (1991).

H. Shen, H. Toki, K. Oyamatsu, and K. Sumiyoshi, NPA 637, 435 (1998).

T>0: the SN approximation

- Equilibrium equations $\vec{x}_A = (A, \delta, V_{WS})$ $\vec{x}_g = (\rho_g, \rho_{Ig})$

$$\left\{ \begin{array}{l} \frac{\partial B_m}{\partial A} = \mu_g \left(1 - \frac{\rho_g}{\rho_0} \right) - \mu_{Ig} \left(\delta - \frac{\rho_{Ig}}{\rho_0} \right) + \frac{3T}{2A} \frac{V_{WS}\rho_0}{V_{WS}\rho_0 - A} \\ \frac{1}{A} \frac{\partial B_m}{\partial \delta} = \mu_{Ig} + \frac{1}{\rho_0^2} \frac{\partial \rho}{\partial \delta} (\mu_g \rho_g + \mu_{Ig} \rho_{Ig}) + \frac{3}{2} T \frac{\rho_g}{\rho_0} \frac{\partial \rho V_{WS}\rho_0}{(V_{WS}\rho_0 - A)(\rho_0 - \rho_g)} \\ A^2 \frac{\partial B/A}{\partial A} = -f_{HM} \frac{A}{\rho_0 V_{WS}} - TV_{WS} \left(1 - \frac{3}{2} \frac{1}{V_{WS}\rho_0 - A} \right) \end{array} \right.$$

$$\mu_g = \frac{\partial f_{HM}}{\partial \rho_g}$$

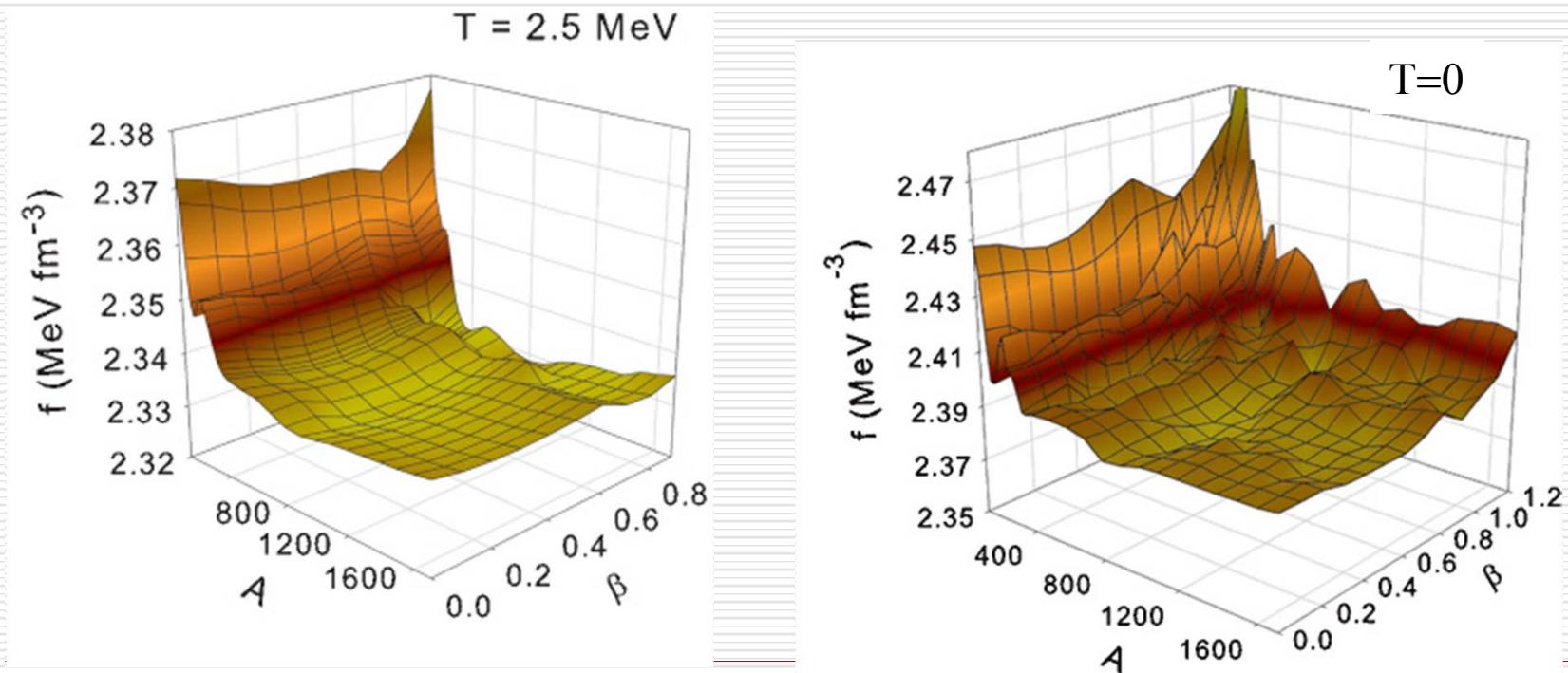
$$\mu_{Ig} = \frac{\partial f_{HM}}{\partial \rho_{Ig}}$$

$$\delta = \frac{(\rho_{0n} - \rho_{0p})}{\rho_0}$$

- These equations give the most probable cluster only and neglect mass and charge fluctuations among cells.*
-

Cluster DoF: $T > 0$

- free-energy surface in constrained HF-BCS (Sly4)



$$\rho_B = 0.08 \text{ fm}^{-3}$$

W.G.Newton J.R.Stone PHYSICAL REVIEW C 79, 055801 (2009)

T>0: beyond the SN approximation

- S_{\max} under constraints, $V \rightarrow \infty \Leftrightarrow F = \min$ in the total volume

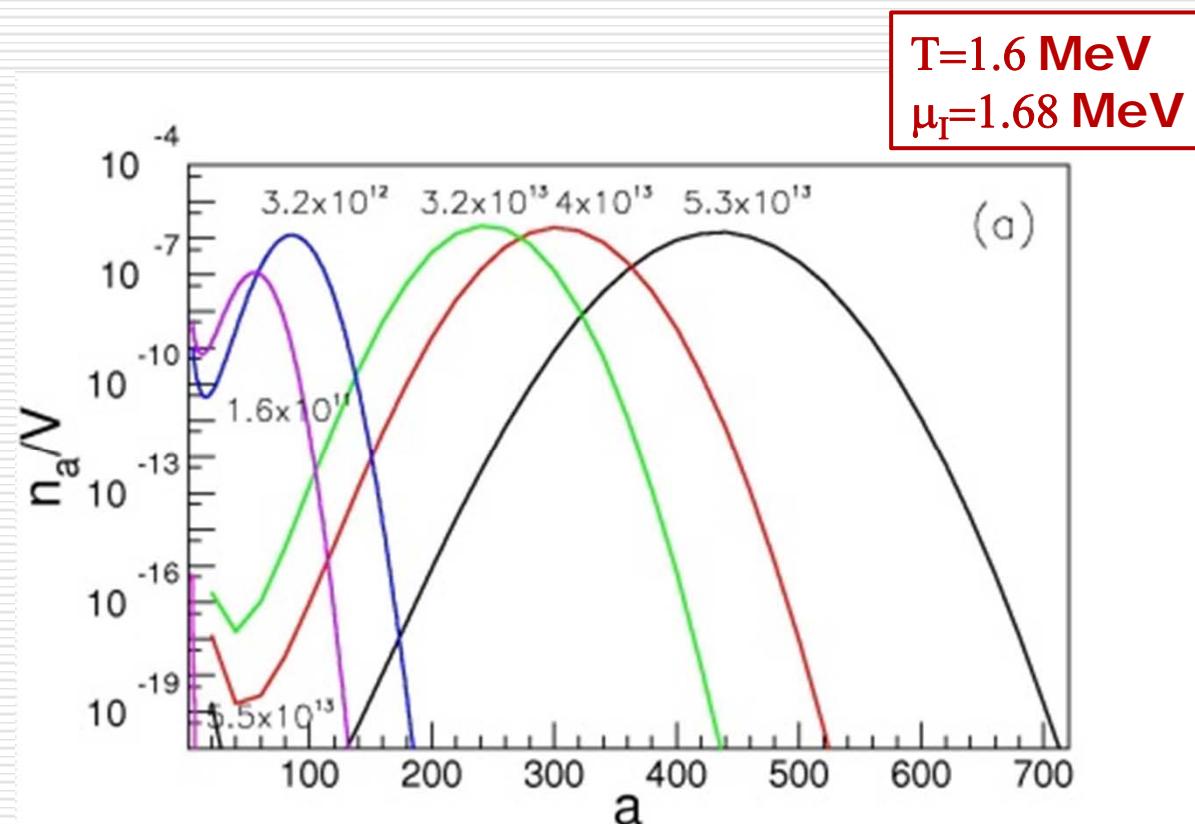
$$-\sum_k p_k \ln p_k - \beta \left(E_{tot} - \langle \hat{H} \rangle_V \right) + \beta \mu \left(A_{tot} - \langle \hat{N} \rangle_V \right) + \beta \mu_I \left(I_{tot} - \langle \hat{I} \rangle_V \right)$$

$$k = \{n_i^{(k)}, A_i, I_i\}$$

$$\bar{p}_k = \frac{1}{Z_\beta} (z_{sky})^V \prod_{A,I} \frac{1}{n_{AI}^{(k)}!} \exp - \beta n_{AI}^{(k)} (B_m - TS)$$

- All cluster A,I possible at a given (ρ_B, y_p)
 - Canonical partition sum analytically calculable via a recursion relation A.Raduta,F.G., PRC 85:025803 (2012)
-

Importance of the cluster distribution



T>0: beyond the SN approximation

□ One WS cell

$$d \left(f_{WS}(A, I, \rho_g, y_g) - \mu \left(\frac{A_{WS}}{V_{WS}} - \rho_B \right) - \mu_I \left(\frac{I_{WS}}{V_{WS}} - \rho_I \right) \right) = 0$$

A, I, ρ_g , ρ_I variational variables linked by the strict conservation law in the cell

□ Many WS cells

$$f(\rho_B, \rho_I) = f_{gas} - \frac{T}{V} \ln \sum_{(k)} \prod_{A,I} \frac{1}{n_{AI}^{(k)}!} \exp - \beta n_{AI}^{(k)} (B_m - TS)$$

$\langle n_{AZ} \rangle = \delta(A - \bar{A}) \delta(I - \bar{I})$ 1 cluster only

$$f(\rho_B, \rho_p) \approx f_{gas} + B_m(\bar{A}, \bar{I}) - TS(\bar{A}, \bar{I}) \text{ with } \mu = \partial f / \partial \rho_B, \mu_I = \partial f / \partial \rho_I$$

$$\Rightarrow d(f_{WS}(\bar{A}, \bar{I}) - \mu \rho_B - \mu_I \rho_I) = 0$$

A, I variational variables linked by the loose conservation law in the cell through the global chemical potential

❖ **Different equations at T>0**

❖ **Same ground state T=0 solution**

Most probable cluster

$$\left\{ \begin{array}{l} \frac{\partial B_m}{\partial A} = \mu_g \left(1 - \frac{\rho_g}{\rho_0} \right) - \mu_{Ig} \left(\delta - \frac{\rho_{Ig}}{\rho_0} \right) + \frac{3T}{2A} \\ \frac{\partial B_m}{\partial \delta} = \mu_{Ig} A + \frac{A}{\rho_0^2} \frac{\partial \rho}{\partial \delta} (\mu_g \rho_g + \mu_{Ig} \rho_{Ig}) + \frac{3}{2} T \frac{\rho_g}{\rho_0} \frac{1}{(\rho_0 - \rho_g)} \\ A^2 \frac{\partial B/A}{\partial A} = -f_{HM} \frac{A}{\rho_0 V_{WS}} + \frac{3}{2} T V_{WS} \end{array} \right.$$

$$\mu_g = \frac{\partial f_{HM}}{\partial \rho_g}$$

$$\mu_{Ig} = \frac{\partial f_{HM}}{\partial \rho_{Ig}}$$



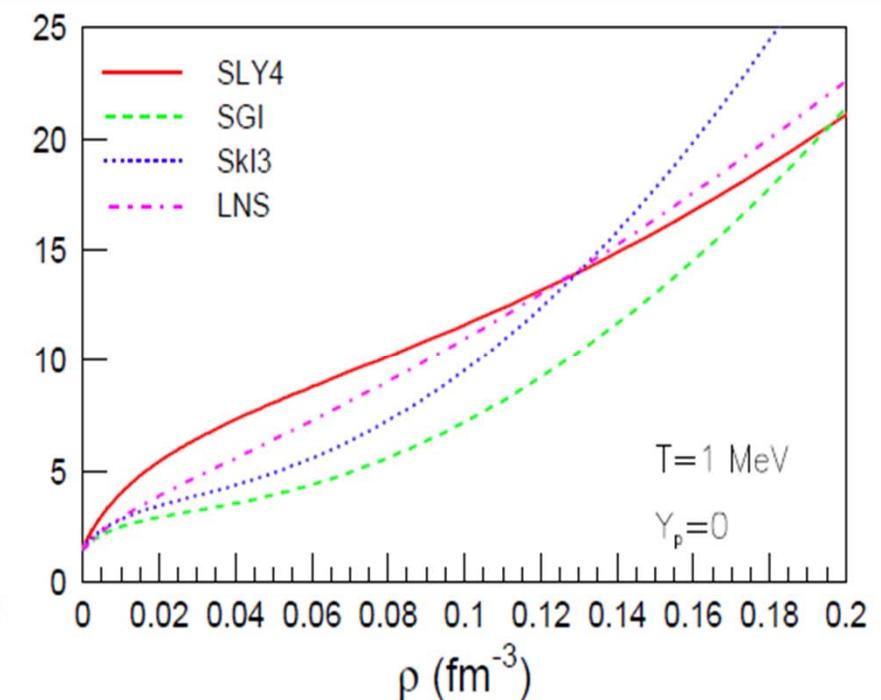
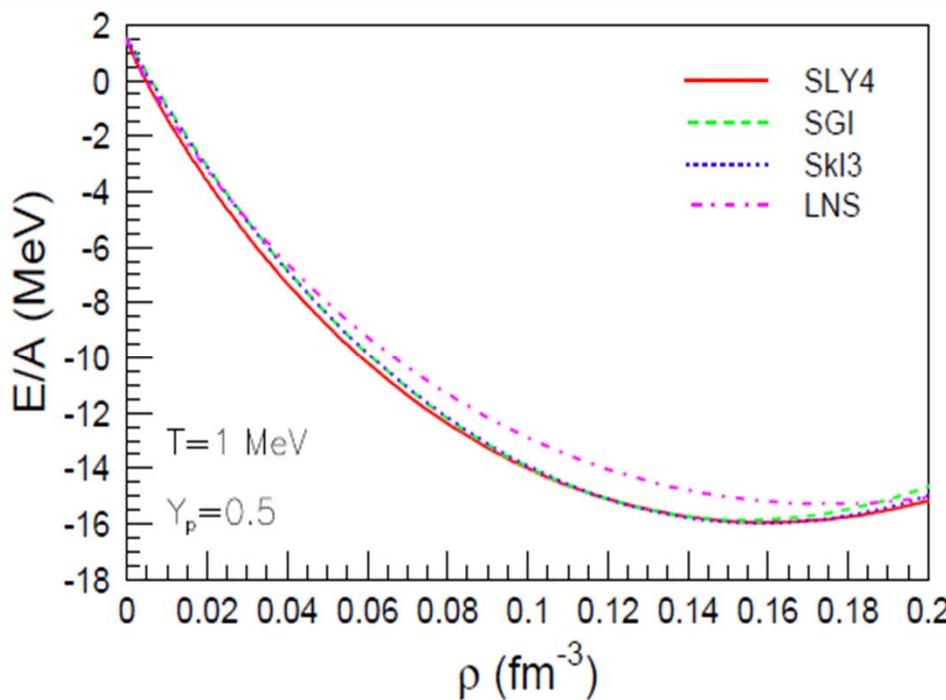
Stellar matter at beta equilibrium

RESULTS AT FINITE TEMPERATURE



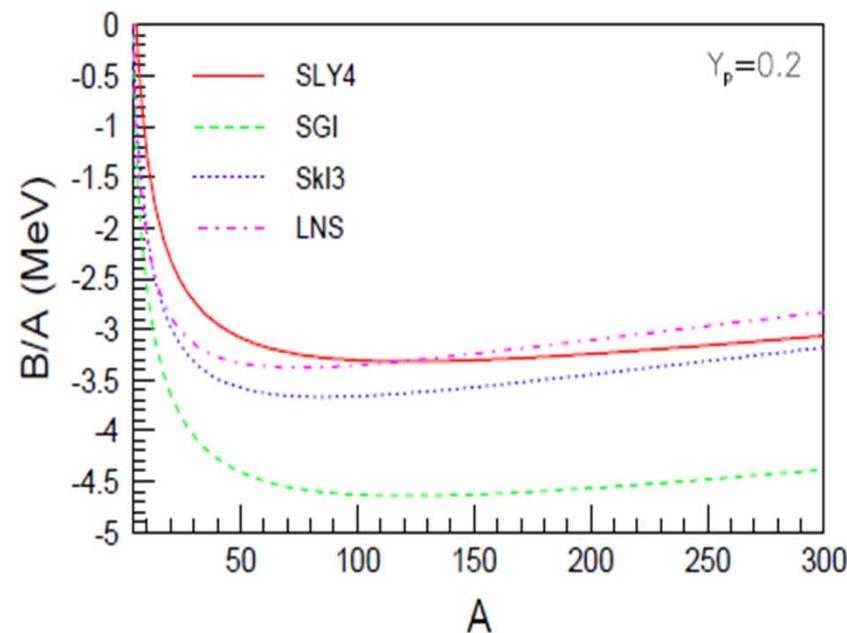
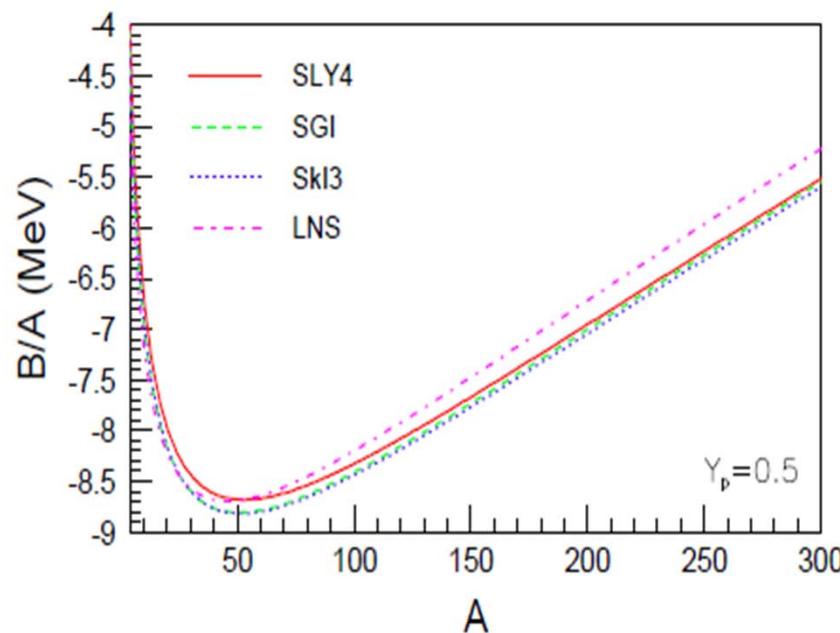
The role of the effective interaction

□ Ingredients: homogeneous matter



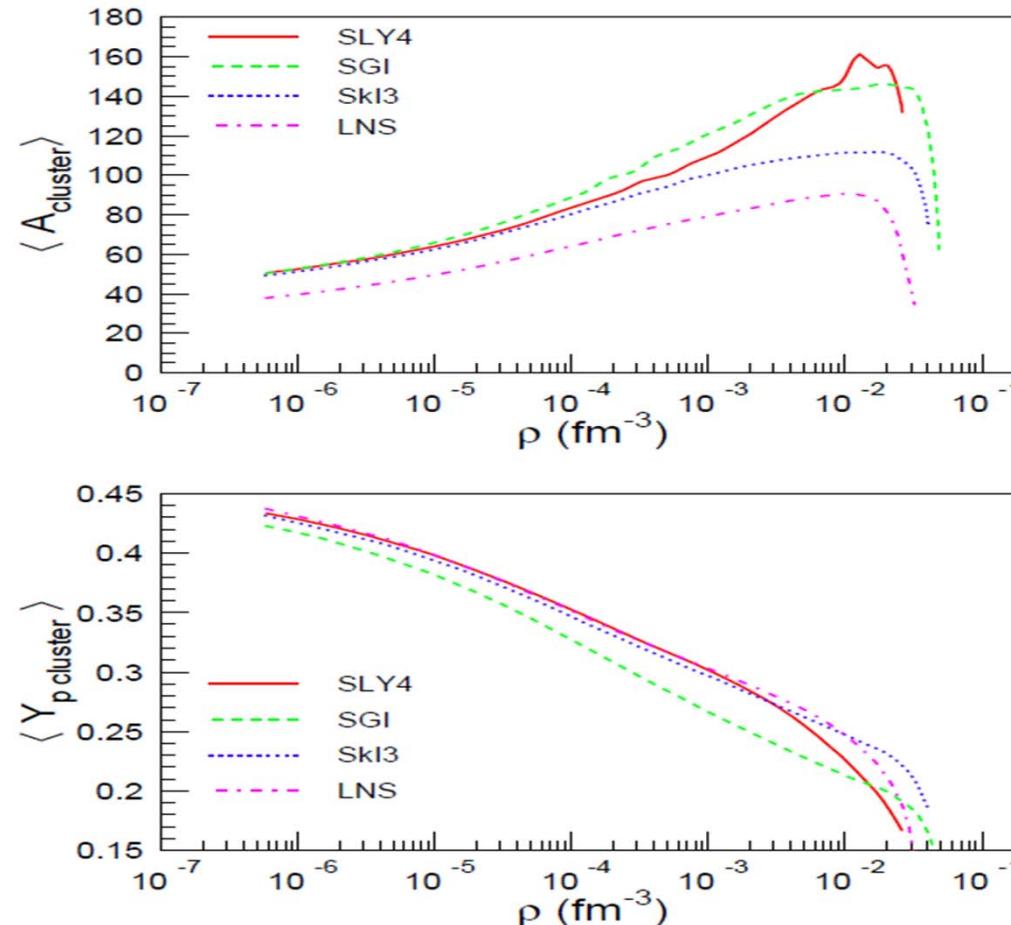
The role of the effective interaction

- ☐ Ingredients: nuclear mass (no shell, no pairing)



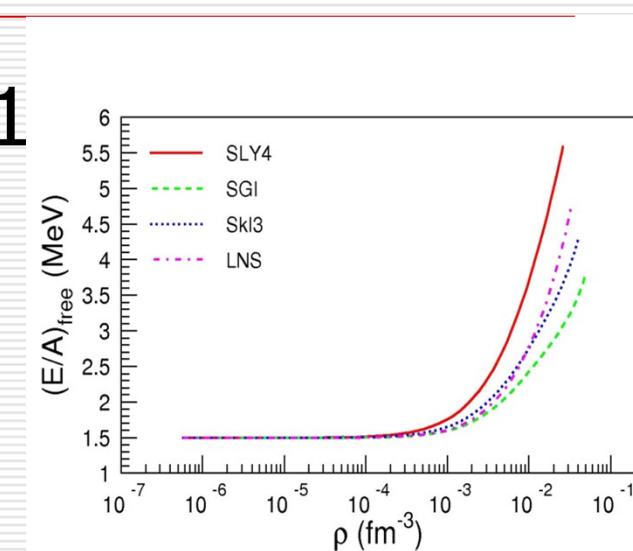
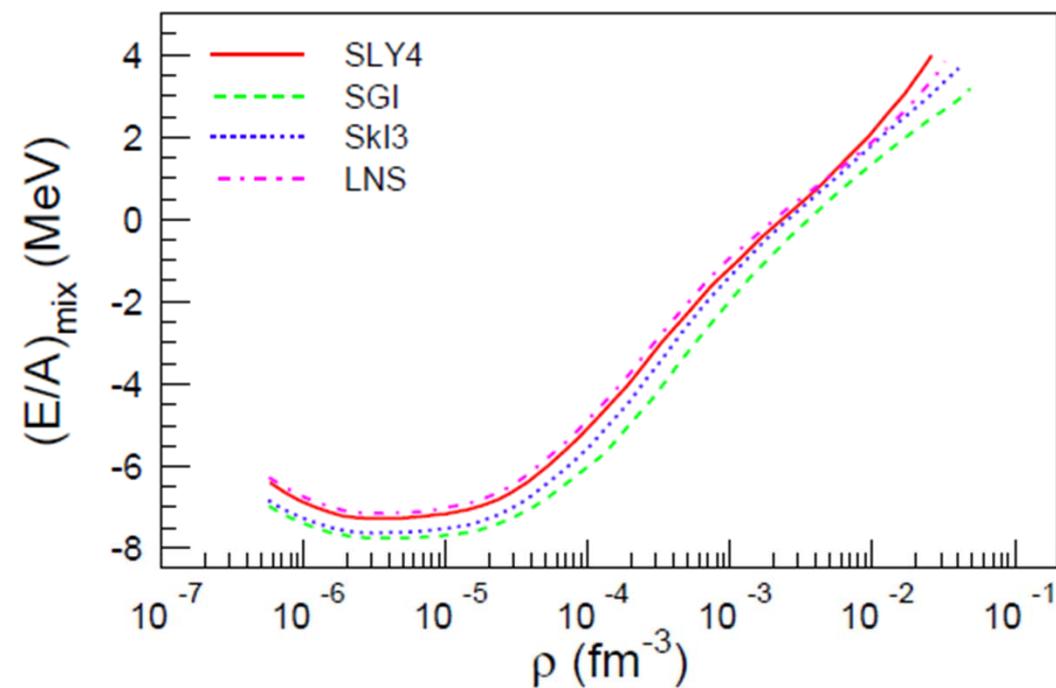
Results: T-dependent β -equilibrium

□ Average cluster mass and isospin T=1



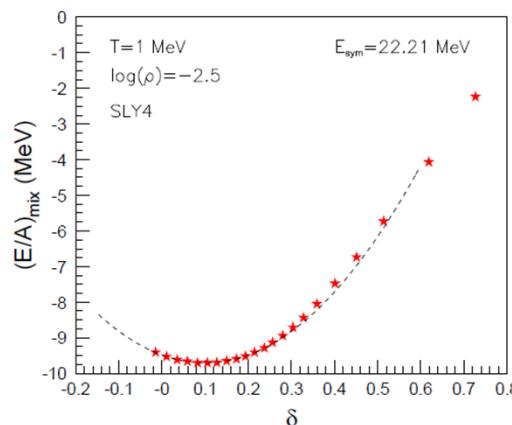
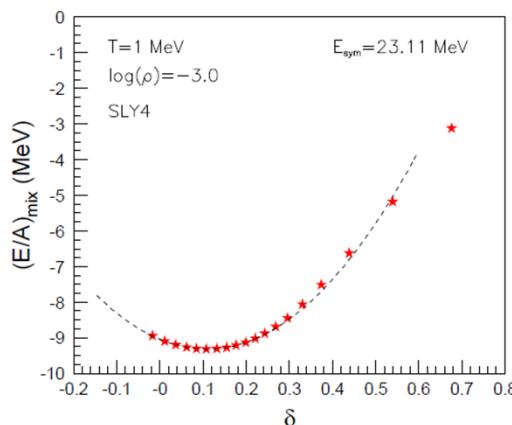
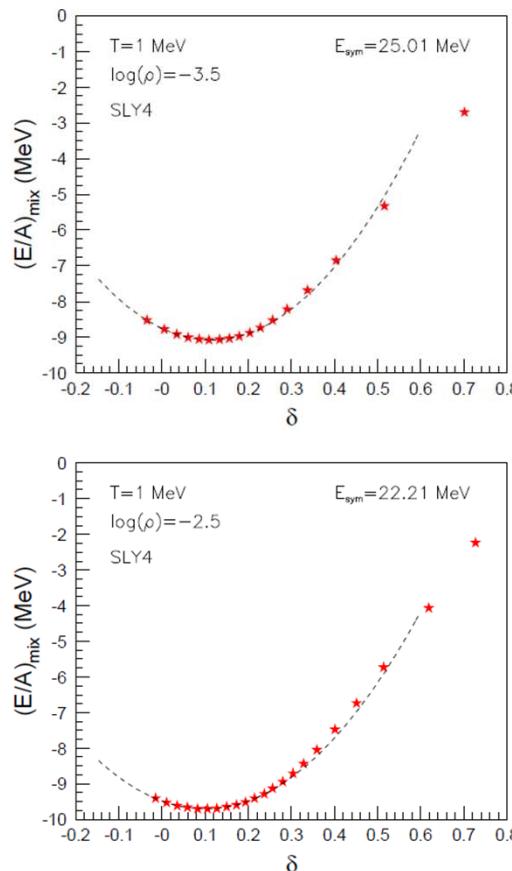
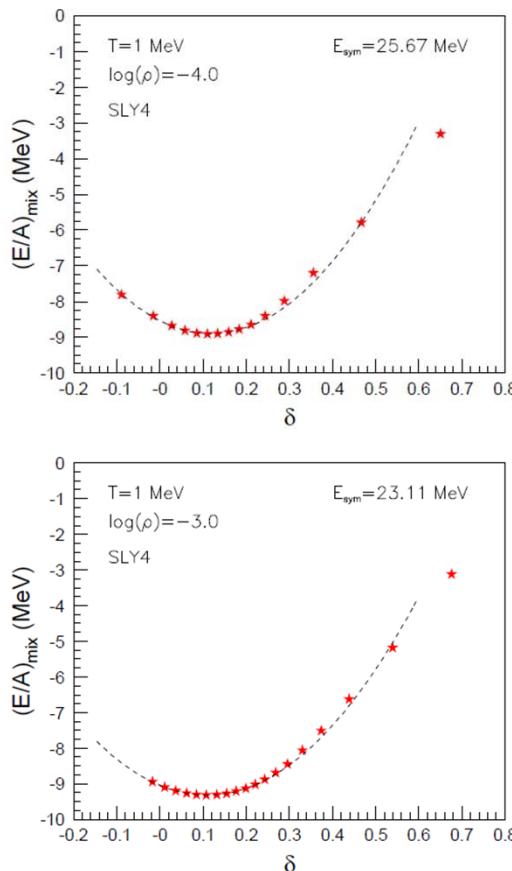
Results: T-dependent β -equilibrium

□ Inhomogeneous EoS T=1



Symmetry energy

□ $\varepsilon(\rho, \delta) = \varepsilon_0(\rho) + \varepsilon_{sym}(\rho)\delta^2$ is not correct!



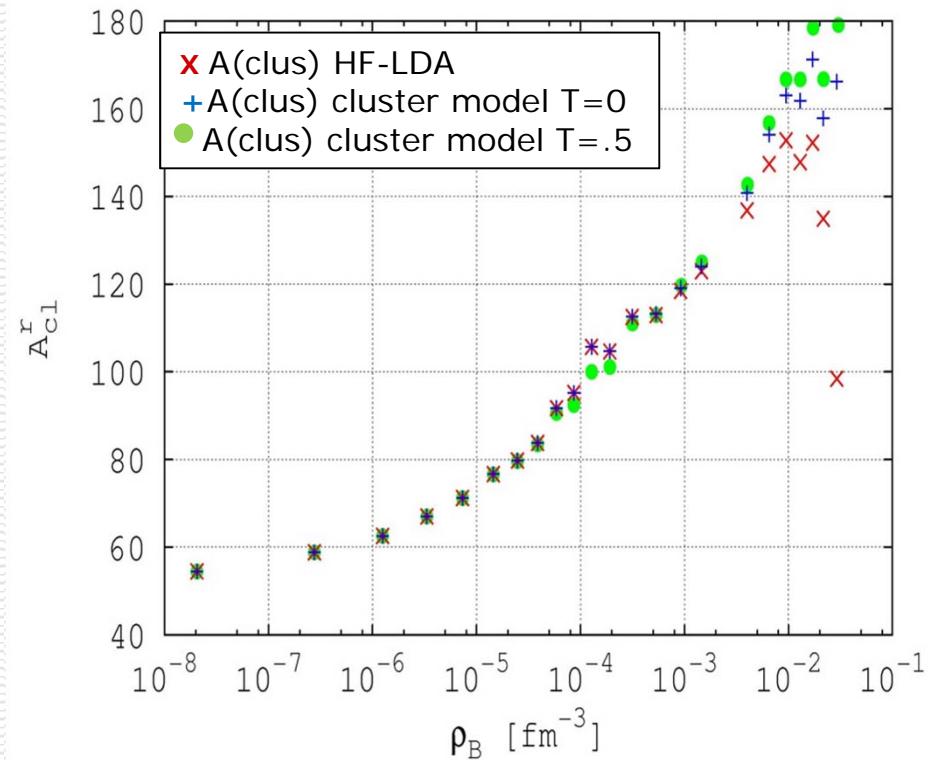


IN MEDIUM ENERGY FUNCTIONAL



Cluster DoF: T=0

- Cluster model is not identical to HF-LDA with the same effective interaction
=> In-medium self energies are not properly treated

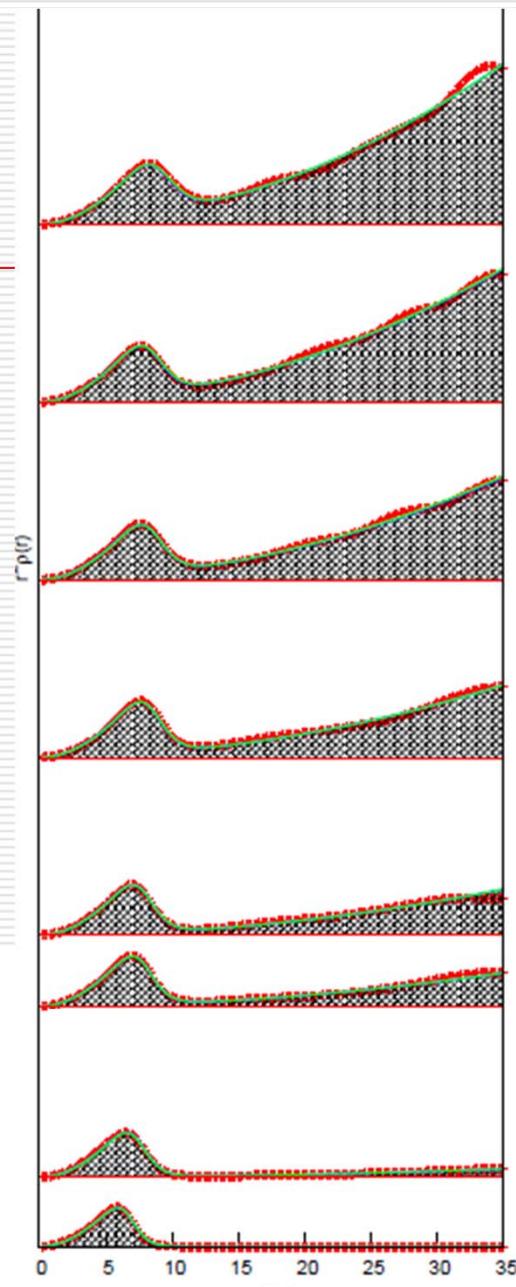
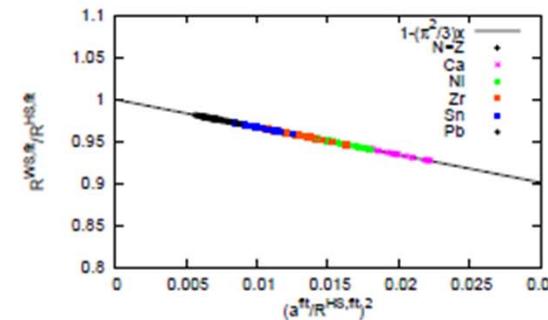
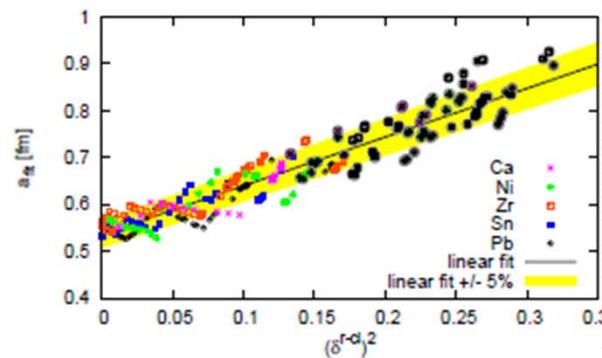
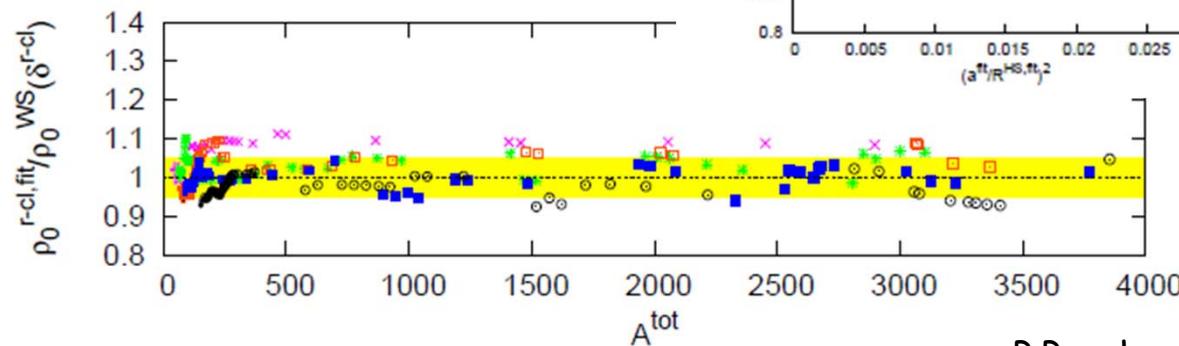


Modelling the HF density profile

$$\rho_{cl}^q(r) = \frac{\rho_0^q(\delta_{bulk})}{1 + \exp \frac{r - R^q}{a^q}}$$

$$\rho_g^q(r) = \frac{\rho_g^q}{1 + \exp \frac{R^q - r}{a^q}}$$

R^q, a^q, δ_{bulk} analytic functions of
 N, Z, ρ_g^n, ρ_g^p



An analytical in-medium cluster energy

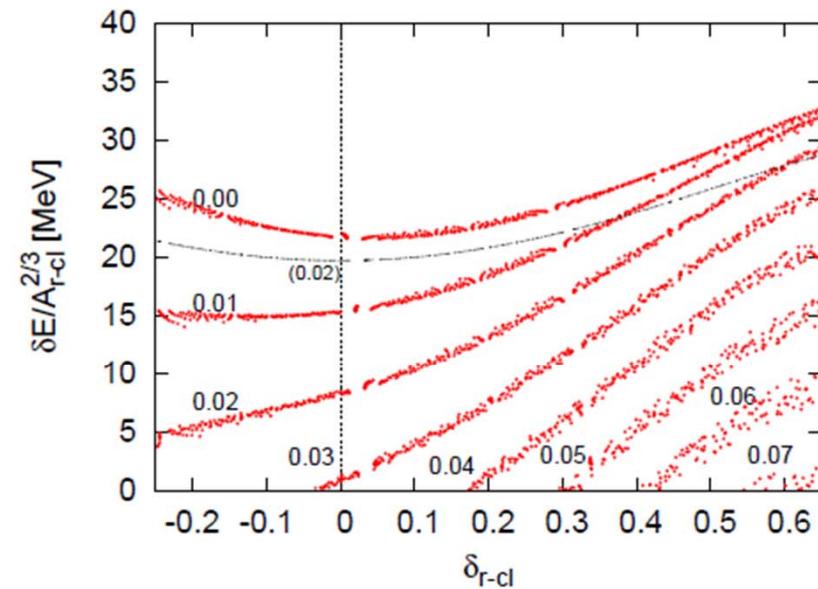
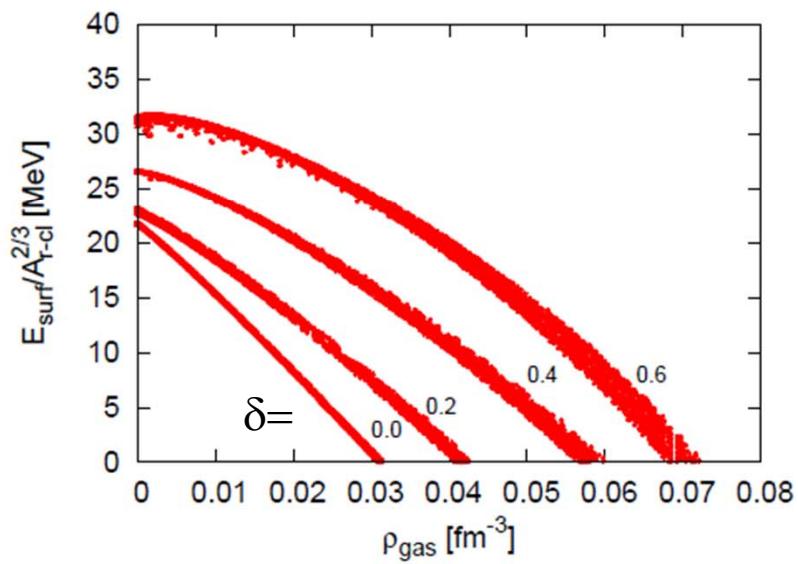
$$\rho^q(r) = \rho_{cl}^q(r) + \rho_g^q(r)$$

$$B^m(\rho_g, \rho_{Ig}, A, I) = \int d^3r \varepsilon_{HF}(\rho^n(r), \rho^p(r)) - \varepsilon_{HF}(\rho_g^n, \rho_g^p) (V_{WS} - A/\rho_0)$$

In-medium modification of the surface mass formula parameters:

$$\delta B^m = B^m - B_0(A, I) = (e_{surf}^m(\rho_g, \rho_{Ig}, I) - a_s) A^{2/3}$$

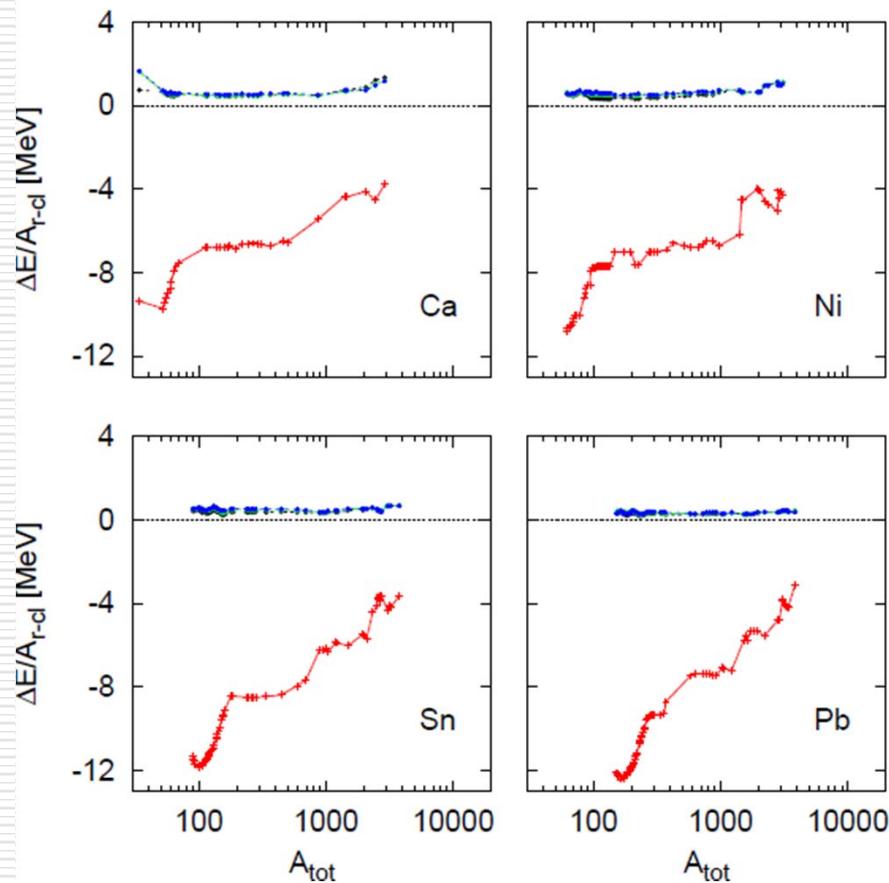
In-medium surface energies



Conclusions

- *Clusters d.o.f. essential to describe hot and dense stellar matter*
 - $\varepsilon_{WS} = B_0 + \varepsilon_{HM} + \delta\varepsilon$, $\delta\varepsilon$ *in-medium self-energy shift*
 - *A simple model: $\delta\varepsilon$ from the excluded volume mechanism+surface corrections from Skyrme-HF*
 - ⇒ *Energetics very different from the vacuum*
 - *Variational calculation at $T=0$ and $T>0$*
 - ⇒ *Quasi-analytic EoS*
 - ⇒ *Wide distributions of clusters*
 - ⇒ *Agreement with microscopic at $T=0$*

Quality of the LDA



$$E_{LDA, \text{model}}^{\text{cl}} - E_{HF}^{\text{cl}}$$

The deviation between the microscopic calculation and the LDA modelling is independent of the medium
=> $\delta e_{bulk}^m(\rho_{gas}, \delta)$,
 $\delta e_{surf}^m(\rho_{gas}, \delta)$
will be correct