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# Chapter 1

## Structure and equation of state of cold non-accreting neutron stars

This chapter deals with the structure and equation of state (EoS) of neutron stars (NS) within the framework of the “cold catalyzed matter” (CCM) hypothesis.

With typical temperature of about  $T \sim 10^8$  K  $\sim 0.01$  MeV/ $k_B$ , NS are very cold systems from the nuclear physics viewpoint, therefore the CCM hypothesis is commonly used to predict their internal composition and pressure. In this limit, thermal, nuclear, and beta equilibrium are established at zero temperature, meaning that the energy cannot be lowered by weak, strong, or electromagnetic processes, thus the matter is in its ground state. It is reasonable to expect these equilibrium conditions to be valid in any NS as far as it is not accreting matter from a neighbor. Indeed, in the accretion scenario, the typical timescale of the process are such that the matter composition is believed to be out of equilibrium.

As already discussed in the general introduction, the EoS relates, in given conditions of temperature and densities, the macroscopic quantities of the star, such as the mass density and the pressure, which determines among others the mass-radius relation of NS. The evaluation of the EoS therefore implies to know the microscopic composition at each point in the star. At subsaturation densities, the solid crust consists mainly of clusterized matter, arranged in a body-centered cubic lattice [HPY07]. The relevant degrees of freedom in the crust are the Wigner-Seitz (WS) cells, containing exactly one lattice point [WS33]. At zero temperature, WS cells are supposed to be identical, thus the so-called single nucleus approximation (SNA), which considers a unique configuration for a given thermodynamic condition of temperature and pressure ( $P, T$ ), becomes exact. The ground state of the outer crust is almost entirely characterized by experimental nuclear masses, which are available up to  $(N - Z)/A \lesssim 0.3$ . The determination of inner-crust ground state is however more challenging because the crust is permeated by free neutrons, a situation which cannot be achieved in terrestrial conditions. Therefore, different treatments, from microscopic [NV73] to classical [BBP71], can be envisaged to estimate the energy of matter, and the EoS in this region depends on the nucleon-nucleon (NN) effective interaction or energy functional. At suprasaturation densities, matter consists of a uniform plasma of neutrons, protons, electrons, and eventually muons, in both strong and weak equilibrium. The development of a unified EoS, that is such that

matter at subsaturation and supersaturation densities are treated within a unique model, is essential if one wants to make realistic predictions on NS observables [For+16].

The plan of the chapter is as follows. In Section 1.1, the ground state of the outer crust is determined by application of the method introduced by Baym, Pethick, and Sutherland (BPS) [BPS71], using experimental masses [Hua+17] supplemented by state-of-the-art microscopic theoretical mass tables. Section 1.2 is devoted to the determination of the inner-crust ground state using a compressible liquid drop model (CLDM) based on the metamodeling technique [MHG18a; CGM19a]. The phase transition from the solid crust to the liquid core, occurring at some  $\approx 1$  km from the surface of the star, is investigated. In Section 1.3, we calculate the ground state of matter in the outer core, and we address the problem of the inner-core composition. In Section 1.4, a unified metamodeling of the EoS is proposed. Finally, conclusions are given in Section 1.5.

## 1.1 Ground state of the outer crust

At zero temperature, the matter inside the outer crust corresponds to a lattice of strongly bound nuclei, immersed in an sea of electrons. The mass density at which nuclei are fully ionized and electron completely degenerated is of the order of  $\rho \gg 6AZ \sim 10^4$  g/cm<sup>3</sup> for <sup>56</sup>Fe, which is the ground state of matter at very low density. Below  $10^4$  g/cm<sup>3</sup>, some electrons are still bound to the nuclei and one must rely on the EoS calculated by Feynman, Metropolis, and Teller from  $15$  to  $10^4$  g/cm<sup>3</sup>, suitable for the envelope of neutron stars [FMT49].

This section deals with the determination of the outer-crust ground state. In 1.1.1, we detail the different terms entering in the WS cell energy, with emphasis on the relativistic electron gas energy as well as nuclear masses. The ground state of the outer crust is calculated by application of the variational BPS method [BPS71], which is presented in 1.1.2. Finally, using current knowledge on experimental masses [Hua+17] supplemented by different microscopic theoretical mass tables, we compute the equilibrium composition and EoS. Our results are presented in 1.1.3.

### 1.1.1 Wigner-Seitz cell energetics

In the outer crust, a WS cell is composed of a strongly bound nucleus at the center, immersed in a relativistic electron gas of density  $n_e$ . Charge neutrality is assured in each unit cell,  $n_e = n_p$ , with  $n_p$  the proton density inside the cell.

The energy of a WS cell in the outer crust can therefore be written as

$$E_{WS} = E_i + V_{WS}\varepsilon_e, \quad (1.1)$$

with  $E_i$  the ion energy,  $V_{WS}$  the volume of the cell, and  $\varepsilon_e$  the energy density of the relativistic electron gas. The ion energy reads

$$E_i = M'(A, Z)c^2 + E_L + E_{zp}, \quad (1.2)$$

where  $M'(A, Z)c^2$  is the nuclear mass of a nucleus with associated mass number  $A$  and charge number  $Z$ ,  $E_L$  the temperature-independent static-lattice term, and  $E_{zp}$  the



zero-point quantum vibration term given by

$$E_{zp} = \frac{3}{2} \hbar w_p u_1, \quad (1.3)$$

where  $u_1 = 0.5113875$  is a numerical constant for a body-centered cubic lattice (see Table 2.4 of [HPY07]), which is assumed to be the geometry minimizing the lattice energy. This assumption was recently confirmed in [CF16]. The ion plasma frequency  $w_p$  is given by

$$\hbar w_p = \sqrt{\frac{(\hbar c)^2 4\pi n_N (Ze)^2}{M'(A, Z) c^2}}, \quad (1.4)$$

$e$  being the elementary charge,  $c$  the speed of light,  $\hbar = h/2\pi$  the reduced Planck constant, and where the ion density  $n_N = 1/V_{WS}$  has been introduced.

The lattice energy reads

$$E_L = -C_M \frac{(Ze)^2}{a_N}, \quad (1.5)$$

with  $C_M = 0.895929255682$  the Mandelung constant for a body-centered cubic lattice (see Table 2.4 of [HPY07]), and  $a_N = (4\pi n_N/3)^{-1/3}$  the ion-sphere radius.

#### 1.1.1.1 Nuclear mass tables

An essential input for Eq. (1.2) is the nuclear mass table. When available, that is for  $I = (N-Z)/A \lesssim 0.3$ , we use experimental masses from the 2016 Atomic Mass Evaluation (AME) [Hua+17]. For more neutron-rich nuclei and until we reach the neutron drip line, the use of a model is required, thus a model dependence arises. A possibility is to rely on microscopic Hartree-Fock-Bogoliubov (HFB) theoretical mass tables [Sam+02], which are based on the nuclear energy-density functional theory that is discussed in the general introduction.

In general, atomic masses are tabulated instead of the nuclear ones which can be calculated as

$$M'(A, Z)c^2 = M(A, Z)c^2 - Zm_e c^2 + B_e(Z), \quad (1.6)$$

with  $M(A, Z)c^2 = \Delta\epsilon + Am_u c^2$  the atomic mass ( $\Delta\epsilon$  is the mass excess,  $m_u$  is atomic mass unit),  $m_e$  the electron mass, and  $B_e$  the binding energy of atomic electrons depending solely on the number of protons  $Z$  according to the approximation proposed in [LPT03] (see their Eq. (A4)),

$$B_e(Z) = 1.44381 \times 10^{-5} Z^{2.39} + 1.55468 \times 10^{-12} Z^{5.35}. \quad (1.7)$$

#### 1.1.1.2 Relativistic electron gas

In the outer crust, the mass densities are above  $\sim 10^4$  g/cm<sup>3</sup> therefore the electrons are essentially free.

The expression of the energy density of a relativistic electron gas, with rest mass energy, at zero temperature can be calculated as

$$\begin{aligned} \varepsilon_e(n_e) &= \int_0^{k_e} \frac{k^2 dk}{\pi^2} c \sqrt{\hbar^2 k^2 + m_e^2 c^2} \\ &= \frac{P_r}{8\pi^2} \left[ x_r (1 + 2x_r^2) \gamma_r - \ln(x_r + \gamma_r) \right], \end{aligned} \quad (1.8)$$

with  $P_r = m_e^4 c^5 / \hbar^3$  the relativistic unit of the electron pressure,  $x_r = \hbar k_e / (m_e c)$  the relativity parameter, and  $\gamma_r = \sqrt{1 + x_r^2}$ .  $k_e$  is the electron Fermi wave number given by

$$k_e = (3\pi^2 n_e)^{1/3}. \quad (1.9)$$

The derivation of Eq. (1.8) is given in Appendix A.

Above  $10^7 \text{ g/cm}^3$ ,  $x_r \gg 1$ , thus electrons can be considered ultrarelativistic and Eq. (1.8) becomes

$$\varepsilon_e(n_e) = \frac{3}{4} n_e m_e c^2 x_r. \quad (1.10)$$

Taking the derivative of Eq. (1.8) with respect to the electron gas density yields the pressure,

$$\begin{aligned} P_e &= -\frac{\partial(V_{WS}\varepsilon_e)}{\partial V_{WS}} = n_e \frac{\partial \varepsilon_e}{\partial n_e} - \varepsilon_e \\ &= \frac{P_r}{8\pi^2} \left[ x_r(1 + 2x_r^2)\gamma_r - \ln(x_r + \gamma_r) \right]. \end{aligned} \quad (1.11)$$

The exchanges and correlation corrections are neglected since they are known to be small in comparison to the kinetic energy of relativistic electrons. The expression of the exchange correction to the free energy density for a strongly degenerate electron gas is given in [HPY07] (see their Eq. (2.151)). We have checked that the inclusion of these corrections does not modify the results presented in this chapter.

### 1.1.2 The BPS model

The variational technique which is currently used to calculate the ground state of the outer crust was introduced by Baym, Pethick, and Sutherland in [BPS71].

The zero-temperature Gibbs free energy per nucleon is the quantity to be minimized at fixed pressure  $P$  under the condition of charge neutrality  $n_e = Zn_N$ , until the neutron drip sets in, the condition for which is  $\mu_n = m_n c^2$ , with  $\mu_n$  the neutron chemical potential, and  $m_n$  the neutron mass.

The definition of the zero-temperature Gibbs free energy per nucleon is

$$g = \frac{\varepsilon_{WS} + P}{n_B}, \quad (1.12)$$

where  $\varepsilon_{WS} = E_{WS}/V_{WS}$  is the energy density of the WS cell, and  $n_B$  is the baryon density given by  $n_B = n_N A = A/V_{WS}$ .

We can calculate the pressure as

$$P = n_B^2 \frac{\partial(\varepsilon_{WS}/n_B)}{\partial n_B} \Big|_{Z,A}, \quad (1.13)$$

yielding, using Eq. (1.1),

$$P = P_e + \frac{1}{3} E_L n_N + \frac{1}{2} E_{zp} n_N. \quad (1.14)$$

Thus the expression of the zero-temperature Gibbs free energy per nucleon can be rewritten as

$$g = \frac{M'(A, Z)c^2 + \frac{4}{3}E_L + \frac{1}{2}E_{zp} + Z\mu_e}{A}, \quad (1.15)$$

where  $\mu_e$  is the electron chemical potential given by

$$\mu_e = \frac{\partial \varepsilon_e}{\partial n_e}. \quad (1.16)$$

Essentially, one fixes the pressure  $P$ , calculates for each nucleus  $(A, Z)$  the electronic density by solving numerically Eq. (1.14), then constructs a table  $(g, A, Z)$ . The ground state of the outer crust at pressure  $P$  then corresponds to the nucleus associated to the minimum value of  $g$ .

As explained in [Pea+18], the neutron chemical potential can be calculated through

$$\mu_n = g, \quad (1.17)$$

because the beta equilibrium equation  $\mu_n = \mu_p + \mu_e$  holds throughout the star. Indeed, by exploiting the thermodynamic relation

$$\mathcal{G} = \varepsilon + P = \mu_n n_n + \mu_p n_p + \mu_e n_e, \quad (1.18)$$

together with the charge neutrality condition  $n_e = n_p$ , we obtain

$$\mathcal{G} = \mu_n n_n + (\mu_e + \mu_p) n_p, \quad (1.19)$$

where  $\mathcal{G}$  is the zero-temperature Gibbs free energy density, and  $\mu_p$  the proton chemical potential. In the case where the chemical equilibrium is established, that is if weak processes are at the equilibrium, we finally get

$$\mathcal{G} = \mu_n n_B. \quad (1.20)$$

One should stress that while minimizing the zero-temperature Gibbs free energy per nucleon at fixed pressure is less practical than simply minimizing the total energy density at constant baryon density, it makes it easier to study the transitions between layers. The pressure increases continuously with increasing depth in the star, thus a discontinuity in the density is the signature of a transition from a layer  $(A, Z)$  to another  $(A', Z')$ . Noting that the pressure in the outer crust is approximately equal to the pressure of the electron gas (the lattice and zero-point terms contribute to less than 5% to the total pressure in the bottom layers of the outer crust), we know that  $n_e = n_p$  is continuous across the transition thus

$$\Delta n_B = n'_B - n_B \simeq n_p \left( \frac{A'}{Z'} - \frac{A}{Z} \right), \quad (1.21)$$

and the fractional change in the baryon mass density results

$$\frac{\Delta \rho_B}{\rho_B} \simeq \frac{\Delta n_B}{n_B} \simeq \frac{Z/A}{Z'/A'} - 1. \quad (1.22)$$

For this reason, it is more convenient to have the pressure as the independent variable. In this way we avoid to make a Maxwell construction to estimate the pressure at which the transition from  $(A, Z)$  to  $(A', Z')$  occurs.

The BPS model is still widely used to determine the ground state of the outer crust. However, at that time the authors took the values of nuclear binding energy from the outdated phenomenological macroscopic model of Myers and Swiatecki [MS65]. Moreover, they did not take into account the zero-point vibration energy, Eq. (1.3). In the last decades, considerable efforts have been made to measure nuclear masses near the neutron drip line [LPT03], as well as theoretical developments were achieved to construct microscopic mass tables [Sam+02]. It is therefore important to reevaluate the ground state of the outer crust considering those experimental and theoretical advances. In this line, Haensel and Pichon studied the consequences of progress concerning the experimental determination of atomic masses in 1994 [HP94]. The authors found that the ground state of the outer crust can be determined exclusively by experimental masses in a fully model independent way up to  $\rho_B \approx 10^{11} \text{ g/cm}^3$ . From this density up to the neutron drip point, the phenomenological, liquid-drop based mass formula of Möller was used, the formalism of which is described in [MN88].

In the same spirit as [HP94], we calculate, in the following, the ground state of the outer crust using the present day knowledge on experimental masses of neutron rich nuclei [Hua+17; Wel+17] combined with state-of-the-art microscopic theoretical mass tables [GCP13].

### 1.1.3 Equilibrium composition and equation of state

We turn to the numerical results obtained applying the BPS method presented in 1.1.2. The ground state of the outer crust is calculated, beginning at  $P = 3 \times 10^{-11}$ , corresponding approximately to  $n_B \approx 10^{-9} \text{ fm}^{-3}$ , in order to ensure complete ionization and electron degeneracy. The pressure is increased from steps of  $0.02P$  until the neutron drip point, defined by  $\mu_n - m_n c^2 = 0$ , is reached.

The ground-state composition and EoS of the outer crust of a cold non-accreting NS is reported in Table 1.1. The upper part of the table,  $n_B < 3.84 \times 10^{-5} \text{ fm}^{-3}$ , corresponding to  $\rho_B < 6.39 \times 10^{10} \text{ g/cm}^3$ , is exclusively determined by experimental data from the AME2016 [Hua+17]. Let us notice that the maximum mass density at which experimentally studied nucleus  $^{80}\text{Zn}$  was present was found to be slightly lower in [HP94],  $\rho_B = 5.44 \times 10^{10} \text{ g/cm}^3$ . This is due to the fact that the determination of this value of density depends on the mass formula used, here HFB-24, for the neutron rich nuclides present in the bottom layers of the outer crust. From  $3.84 \times 10^{-5} \text{ fm}^{-3}$ , matter becomes so neutron rich that the nuclear masses cannot be measured experimentally for now, thus a model dependence is expected to arise because the nuclear masses have to be extrapolated from laboratory nuclei. The nuclear mass model used to calculate the ground state here is HFB-24 [GCP13], constructed from the BSk24 functional following the Hartree-Fock-Bogoliubov method [Sam+02]. The BSk24 functional is part of a family of functionals labeled BSk22 to BSk26, consisting of unconventional effective Skyrme forces with extra  $t_4$  and  $t_5$  terms that behave as density-dependent generalizations of the usual  $t_1$  and  $t_2$  terms present in traditional Skyrme forces. The functionals BSk22-

element	$Z$	$N$	$Y_p$	$n_{B,max}$ (fm <sup>-3</sup> )	$P_{max}$ (MeV/fm <sup>3</sup> )	$\mu_n - m_n c^2$ (MeV)	$\mu_e$ (MeV)
<sup>56</sup> Fe	26	30	0.4643	$4.97 \times 10^{-9}$	$3.40 \times 10^{-10}$	-8.96	0.95
<sup>62</sup> Ni	28	34	0.4516	$1.56 \times 10^{-7}$	$4.09 \times 10^{-8}$	-8.26	2.57
<sup>64</sup> Ni	28	36	0.4375	$8.07 \times 10^{-7}$	$3.60 \times 10^{-7}$	-7.52	4.34
<sup>66</sup> Ni	28	38	0.4242	$9.27 \times 10^{-7}$	$4.16 \times 10^{-7}$	-7.46	4.50
<sup>86</sup> Kr	36	50	0.4186	$1.85 \times 10^{-6}$	$1.03 \times 10^{-6}$	-7.01	5.63
<sup>84</sup> Se	34	50	0.4048	$6.85 \times 10^{-6}$	$5.64 \times 10^{-6}$	-5.87	8.59
<sup>82</sup> Ge	32	50	0.3902	$1.67 \times 10^{-5}$	$1.77 \times 10^{-5}$	-4.82	11.41
<sup>80</sup> Zn	30	50	0.3750	$3.84 \times 10^{-5}$	$5.10 \times 10^{-5}$	-3.58	14.86
<sup>78</sup> Ni	28	50	0.3590	$6.68 \times 10^{-5}$	$1.01 \times 10^{-4}$	-2.63	17.61
<sup>126</sup> Ru	44	82	0.3492	$7.52 \times 10^{-5}$	$1.12 \times 10^{-4}$	-2.47	18.15
<sup>124</sup> Mo	42	82	0.3387	$1.21 \times 10^{-4}$	$2.04 \times 10^{-4}$	-1.54	21.05
<sup>122</sup> Zr	40	82	0.3279	$1.56 \times 10^{-4}$	$2.75 \times 10^{-4}$	-1.03	22.69
<sup>121</sup> Y	39	82	0.3223	$1.63 \times 10^{-4}$	$2.84 \times 10^{-4}$	-0.98	22.86
<sup>120</sup> Sr	38	82	0.3167	$1.95 \times 10^{-4}$	$3.52 \times 10^{-4}$	-0.60	24.12
<sup>122</sup> Sr	38	84	0.3115	$2.37 \times 10^{-4}$	$4.49 \times 10^{-4}$	-0.15	25.62
<sup>124</sup> Sr	38	86	0.3065	$2.56 \times 10^{-4}$	$4.87 \times 10^{-4}$	0.00	26.14

Table 1.1: Ground state of the outer crust of a cold non-accreting neutron star. Experimental data from the 2016 Atomic Mass Evaluation [Hua+17] are used when available. Mass excesses of <sup>77–79</sup>Co are taken from [Wel+17]. Experimental masses are supplemented with masses from microscopic HFB-24 theoretical mass table [GCP13] (lower part). The last line corresponds to the neutron drip point. The following quantities are reported in the table: the element, the atomic number  $Z$ , the neutron number  $N$ , the proton fraction  $Y_p = Z/A$  ( $A$  is the number of nucleons), the maximum baryon density  $n_{B,max}$  at which the nuclide is found, the associated pressure  $P_{max}$ , the neutron chemical potential minus rest mass  $\mu_n - m_n c^2$ , and the electron chemical potential  $\mu_e$ .

26 were fitted to the 2353 experimental masses of the AME2012 [Aud+12] and differ mainly by their symmetry energy  $S$ , defined here as the difference between the energy per nucleon of pure neutron matter (PNM)  $e_{PNM}$  and the energy per nucleon of symmetric nuclear matter (SNM)  $e_{SNM}$ ,

$$S(n) = e_{PNM}(n) - e_{SNM}(n), \quad (1.23)$$

with  $n$  the density of nuclear matter. To a first approximation,  $S(n) \simeq E_{sym}$ ,  $E_{sym}$  being the symmetry energy per nucleon at the saturation density of nuclear matter  $n = n_{sat} \approx 0.16 \text{ fm}^{-3}$ . This quantity is constrained to 32, 31, 30, and 29 MeV for BSk22, BSk23, BSk24, and BSk25, respectively. For BSk26,  $E_{sym} = 30 \text{ MeV}$  as well but the functional is fitted to the APR EoS of PNM [APR98], unlike BSk22–25 that are fitted to the stiffer LS2 EoS [LS08].

The first layer of the outer crust consists of a crystal lattice of <sup>56</sup>Fe, the mass per nucleon of which is the lowest among all nuclides. The sequence of nuclides is in good

agreement with the results of [HP94]. One can note the persistence of magic numbers  $Z = 28$  and  $N = 50$  at low density, and  $Z = 82$  at higher density. In particular, in [HP94] the  $N = 82$  shell is found from  $\rho_B = 9.64 \times 10^{10} \text{ g/cm}^3$  to  $4.32 \times 10^{11} \text{ g/cm}^3$ , point of neutron drip (see Table 3.1 of [HPY07]). It is explained in [HPY07] that the apparent strong effect of the  $N = 82$  shell in the bottom layers could be an artifact of the extrapolation using the mass formula of [MN88] and that further investigations on nuclear shell structure in the vicinity of the neutron drip point were required. Here we find two additional layers of strontium ( $N = 84$  and  $N = 86$ ) with respect to [HP94] results, and the layer of  $^{118}\text{Kr}$  before the neutron drip is not observed.

At high density, a thin layer of odd-mass nuclei  $^{121}\text{Y}$  is found with the HFB-24 mass model. It is interesting since the possibility of having odd nuclei in the ground state of the outer crust was not considered in the original calculation of BPS [BPS71]. Also, it is reported in [Pea+18] that  $^{79}\text{Co}$  is favored over  $^{76}\text{Ni}$ , found for HFB-22 and HFB-25, in the case where the recent mass excess measurements of [Wel+17] are accounted for. This highlights the importance of measuring the mass of odd-nuclei. **One could expect that the presence of odd-nuclei in the outer crust of NS leads to ferromagnetic phase transitions at low temperature. In return, it could generate a magnetic field and alter the existing field, and so the electron gas.**

It can also be observed in Table 1.1 that the proton fraction  $Y_p = Z/A$  always decreases with increasing depth. The neutronization of matter can be understood by the following reason. Neglecting the static-lattice energy as well as the zero-point quantum vibrations terms, we can write the zero-temperature Gibbs energy per nucleon as

$$g \simeq \frac{M'(A, Z)c^2 + V\varepsilon_e}{n_B V} + \frac{P}{n_B}. \quad (1.24)$$

Since the lattice does not contribute to the pressure, we have  $P \simeq P_e = n_e \mu_e - \varepsilon_e$ , yielding

$$g \simeq \frac{M'(A, Z)c^2}{A} + \frac{n_e}{n_B} \mu_e. \quad (1.25)$$

The charge neutrality is ensured in the unit cell, therefore  $n_e/n_B = Z/A$  and we finally get

$$g \simeq \frac{M'(A, Z)c^2}{A} + \frac{Z}{A} \mu_e. \quad (1.26)$$

For  $\rho_B \gg 10^{-7} \text{ g/cm}^3$ , electrons are ultrarelativistic and the electron chemical potential is calculated by taking the derivative of Eq. (1.10),

$$\mu_e = \frac{\partial \varepsilon_e}{\partial n_e} = \frac{3}{4} \hbar c (3\pi^2)^{1/3} n_e^{4/3}. \quad (1.27)$$

Therefore, the electron chemical potential  $\mu_e$  scales as  $P^{1/4}$  ( $P \propto n_e^{4/3}$ ). Then it appears that, with increasing pressure, it is energetically favorable to decrease the proton fraction  $Y_p = Z/A$  in order to compensate the increase in the term  $M'(A, Z)c^2/A$ .

The neutron chemical potential monotonously increases with increasing density and at  $n_B = 2.56 \times 10^{-4} \text{ fm}^{-3}$ , corresponding to  $P = 4.87 \times 10^{-4} \text{ MeV/fm}^{-3}$ , the neutron drip is finally reached. It should be mentioned that the neutron drip density as well as pressure depend on the mass model, here chosen to be HFB-24.



Figure 1.1: Variation with baryon density  $n_B$  of the equilibrium value of atomic number  $Z$  and neutron number  $N$  in the bottom layers of the outer crust for four different models: HFB-24, HFB-26 [GCP13], HFB-14 [GSP07], and the FRDM [Mol+95].

The model dependence of the composition in the bottom layers of the outer crust (bottom part of Table 1.1) can be seen in Figure 1.1, where we represent the proton number  $Z$  and the neutron number  $N$  as a function of the baryon density  $n_B$  for four different mass models that reproduce with comparable accuracy the present experimental mass information: HFB-14 [GSP07], HFB-24, HFB-26 [GCP13], and the FRDM [Mol+95] that supplement the experimental data. The model dependence is observed from  $n_B \approx 2.5 \times 10^{-5} \text{ fm}^{-3}$ , where the experimental mass data are not available. We recover similar sequences of nuclides with the different mass models, with the strong shell effect associated to  $N = 50$  at low density and  $N = 82$  in the bottom layers. A thin layer with  $N = 52$  is also found for HFB-24 and HFB-26 just before the transition to  $N = 82$ . While this transition occurs in the vicinity of  $\approx 8 \times 10^{-5} \text{ fm}^{-3}$  for HFB-24, HFB-26, and the FRDM, it is found to happen at approximately  $4 \times 10^{-5} \text{ fm}^{-3}$  for HFB-14. Different nuclides are found close to the neutron drip, depending on the model:  $^{124}\text{Sr}$  is found for HFB-24,  $^{126}\text{Sr}$  for HFB-26,  $^{120}\text{Kr}$  for HFB-14, and  $^{122}\text{Kr}$  for the FRDM. The neutron drip density  $n_{ND}$  and pressure  $P_{ND}$  slightly depends on the model as well. These values are reported in Table 1.2.

The variation of pressure with baryon density, namely the EoS, is shown in Fig. 1.2 for the four mass models. The selected models give the same value of pressure up to  $\approx 2.5 \times 10^{-5} \text{ fm}^{-3}$ , as it is the case for the composition. At higher densities, we can perceive the small effect of the model, in particular at the transitions from a layer to

model	element	$Z$	$N$	$n_{ND}$ (fm $^{-3}$ )	$P_{ND}$ (MeV/fm $^3$ )
HFB-24	$^{124}\text{Sr}$	38	86	$2.56 \times 10^{-4}$	$4.87 \times 10^{-4}$
HFB-26	$^{126}\text{Sr}$	38	88	$2.62 \times 10^{-4}$	$4.91 \times 10^{-4}$
HFB-14	$^{120}\text{Kr}$	36	84	$2.67 \times 10^{-4}$	$5.01 \times 10^{-4}$
FRDM	$^{122}\text{Kr}$	36	82	$2.62 \times 10^{-4}$	$4.99 \times 10^{-4}$

Table 1.2: Chemical element, atomic number  $Z$ , neutron number  $N$ , density  $n_{ND}$ , and pressure  $P_{ND}$  at the neutron drip point for four different mass models: HFB-24, HFB-26 [GCP13], HFB-14 [GSP07], and the FRDM [Mol+95].



Figure 1.2: Variation with baryon density  $n_B$  of pressure  $P$  for four different models: HFB-24, HFB-26 [GCP13], HFB-14 [GSP07], and the FRDM [Mol+95].

another at which the pressure stays constant, but where a discontinuity is observed in the density.

## 1.2 Ground state of the inner crust

The neutron drip marks the onset of the inner crust, which is a system that cannot be reproduced in the laboratory since the dripped neutrons would evaporate, while they stay confined in the WS cell due the gravitational pressure. For this reason, the equilibrium composition and so the EoS in the inner crust are fully model dependent.



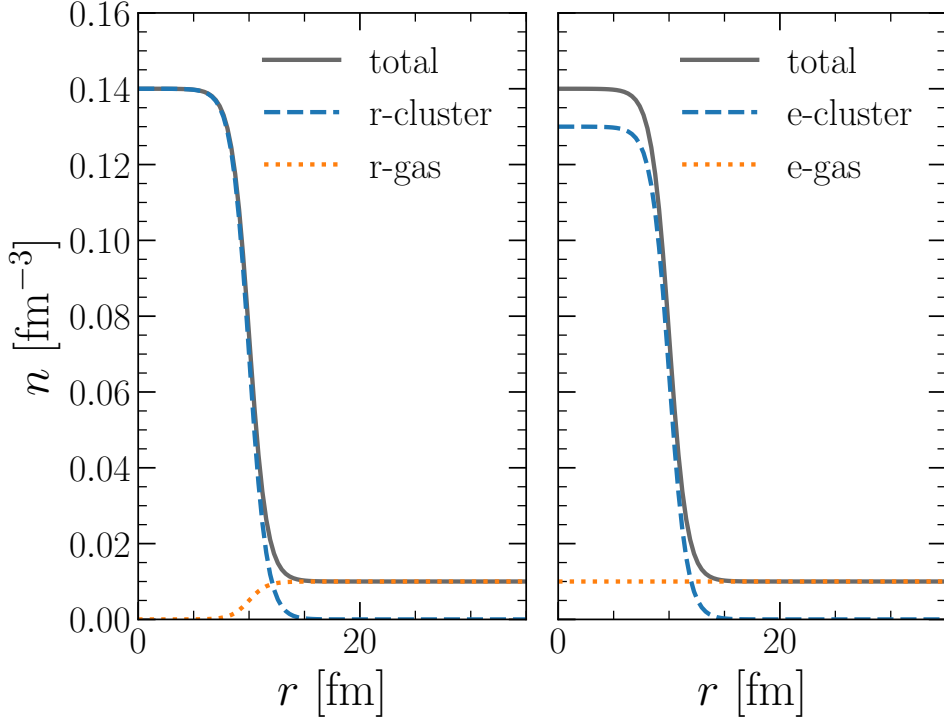


Figure 1.3: Wood-Saxon density profiles, for arbitrary values, within a WS cell in the free neutron regime. Blue dashed lines correspond to the cluster density (left, r-cluster; right, e-cluster), and orange dotted lines to the gas density. The total density profile is represented in gray solid lines.

This section deals with the determination of the ground state of the inner crust. In 1.2.1, we present the metamodeling of the infinite nuclear matter EoS here used to calculate the energy of the neutron gas, and extended to finite nuclei within the so-called compressible liquid drop (CLD) approximation so as to describe the cluster energetics. The derivation of the system of equilibrium equations for the determination of inner-crust ground state is detailed in 1.2.2. The numerical method as well as results are presented afterwards, in 1.2.3. In 1.2.4 we add Strutinsky shell corrections on top of the CLD energy as an attempt to recover magic numbers in the free neutron regime. Nonspherical pasta phases in the bottom layers of the inner crust are considered in 1.2.5. Finally, the phase transition from the solid crust to the liquid core is investigated in 1.2.6.

### 1.2.1 Modeling the nuclear energy

Once the neutron dripline is reached, neutrons start to drip out of nuclei but stay confined in the WS cell because of the gravitational pressure, whereas they would have been emitted in the laboratory. In the regime of the inner crust, we thus have, in each unit cell, a cluster immersed in an electron sea, and an ambient neutron gas.

There is an ambiguity regarding the characterization of the cluster-gas interface. One should ask whether the neutron gas penetrate the cluster or not. Hence, two classical representations can be introduced, namely the r-cluster and e-cluster representations [Pap+13], to interpret the distribution of the  $N_{WS}$  neutrons and  $Z_{WS}$  protons in the unit cell. In the former, the cluster occupy a volume at the center of the cell and is surrounded by the neutron gas. Evidently, there is a thin region where cluster and gas overlap given the fact that a sharp interface would be unrealistic. The r-cluster representation naturally emerges in the local density approximation in density functional theory. Indeed, if the energy is expressed as a function of the local density, then the cluster corresponds to the dense part and the gas to the dilute one. This interpretation is used in most of the calculations at finite temperature to model supernovae, see for example the renowned Lattimer and Swesty EoS [LS91]. In the e-cluster representation, the gas penetrate the cluster. This interpretation appears spontaneously in single-particle developments. Indeed, in very neutron-rich clusters, beyond the neutron dripline, the bound states are occupied as well as the resonant and continuum states. All the unbound single-particle states that are occupied thus represents the neutron gas, characterized by a quasihomogeneous spatial distribution, the continuum wave functions being very similar to plane waves. **The difference between the two representations is illustrated in Fig. 1.3. The cluster and gas density profiles, described by Woods-Saxon profiles with arbitrary values, are plotted along the WS cell radius in the r-cluster representation (left) and e-cluster representation (right). Woods-Saxon profiles are known to give a good description of medium-mass and heavy nuclei density profiles, and are commonly used in Thomas-Fermi (TF) and extended Thomas-Fermi (ETF) calculations [Ons+08; Pea+18]. In both representations, the total density profile (solid gray line) is the same. One can benefit from the advantage of both representations through the simple geometric relations**

$$A_e = A \left( 1 - \frac{n_g}{n_0} \right), \quad Z_e = A \left( 1 - \frac{n_{g,p}}{n_{0,p}} \right), \quad (1.28)$$

where  $A_e$  ( $Z_e$ ) is the number of nucleons (protons) in the e-cluster,  $A$  ( $Z$ ) the number of nucleons (protons) in the r-cluster,  $n_g$  ( $n_{g,p}$ ) the total (proton) gas density, and  $n_0$  ( $n_{0,p}$ ) the average total (proton) density inside the cluster. Since we do not consider the possibility of proton drip in the inner crust, we have  $n_{g,p} = 0$ , yielding  $Z_e = Z$  and implying that the total gas density is equal to that of the neutron gas. Indeed, while free protons are expected at non-zero temperature, their presence at zero temperature remains uncertain [BBP71] and depends on the nuclear model. For some models, the proton drip could set in the very bottom layers of the crust, the classical condition for which is  $\mu_p = U_p + m_p c^2$ , where  $U_p$  corresponds to the proton single-particle field at the cell surface [Pea+18]. However, if nonspherical shapes are considered then it is found that protons remain in the cluster for most models.

At zero temperature, we write the WS cell energy  $E_{WS}$  in the r-cluster representation as

$$E_{WS} = E_{cl} + V_{WS}\varepsilon_e + (V_{WS} - V_{cl})\varepsilon_g + Z_{WS}(m_p - m_n)c^2 + A_{WS}m_n c^2 \quad (1.29)$$

where  $\varepsilon_g$  represents the energy density of the surrounding neutron gas of density  $n_g$ ,  $V_{cl} = A/n_0$  the volume of the cluster,  $V_{WS}$  that of the WS cell, and  $E_{cl}$  the energy of

the cluster. The total number of protons  $Z_{WS}$  and nucleons  $A_{WS}$  inside the WS cell are given respectively by

$$Z_{WS} = A \frac{1 - I}{2} \quad \text{and} \quad A_{WS} = A \left( 1 - \frac{n_g}{n_0} \right) + n_g V_{WS}. \quad (1.30)$$

Let us notice that the number of protons inside the cell remains an integer number, as in the outer crust, unlike the number of neutrons that is noninteger because of the outside neutron gas.

As for the bottom layers of the outer crust, the determination of the ground state of the inner crust is fully model dependent. Therefore, the two necessary ingredients of the WS cell energy are the equation of state of PNM, and the energy of the cluster, which will be treated in the CLD approximation introduced by Baym, Bethe, and Pethick (BBP) in their pioneering work [BBP71].

### 1.2.1.1 Metamodeling of homogeneous nuclear matter

The ambient neutron gas in the WS cell is treated as homogeneous nuclear matter. In this limit, a large number of nucleons  $A \rightarrow \infty$  is contained in a large box. The system is entirely characterized by the neutron density  $n_n$  and the proton density  $n_p$ . Let us also introduce the total density  $n = n_n + n_p$  and the asymmetry  $\delta = (n_n - n_p)/n$ . Two limit cases can be distinguished: pure neutron matter, where  $n_p = 0$ , and symmetric nuclear matter, where  $n_n = n_p$ . Beta-equilibrated matter corresponds to the matter inside the core of neutron stars and is treated in Section 1.3. Different methods have been proposed in order to evaluate the energy per nucleon of homogeneous nuclear matter. While considerable theoretical efforts have been devoted to ab initio approaches in the recent years [Gan+14], these calculations might become unreliable at suprasaturation densities. Also, the treatment of three-body forces, which are known to play a major role in the determination of the saturation properties of nuclear matter, remains a challenge [DHS16]. Another possibility is to derive the nuclear EoS from effective forces [Cha+98]. This second approach is very practical because of translational invariance but suffers from the possible introduction of artificial correlations among the parameter space. In the following we present the metamodeling technique [MHG18a; MHG18b]. As we will see, this technique allows to parametrize in a simple analytical way the energy per nucleon obtained in the different ab initio or effective approaches, as well as interpolate continuously between them. By largely exploring the parameter space of this metamodel (or model of models), constraints coming from experimental measurements can be directly implemented. Indeed, nuclear experiments give us knowledge on the properties of nuclear matter around the saturation density of symmetric nuclear matter  $n_{sat}$ , which is roughly the density of laboratory nuclei. For instance, the isoscalar Giant Monopole Resonance energy is correlated with the empirical parameter  $K_{sat}$ , which corresponds to the isoscalar modulus. The well-known nuclear EoS empirical parameters correspond in fact to the successive density derivatives of nuclear matter energy per particle of SNM and symmetry energy at saturation density, associated to the isoscalar ( $E_{sat}$ ,  $K_{sat}$ ,  $Q_{sat}$ , and  $Z_{sat}$ ) and isovector channel ( $E_{sym}$ ,  $L_{sym}$ ,  $K_{sym}$ ,  $Q_{sym}$ , and  $Z_{sym}$ ), respectively. The

Parameter	Unit	$N$	SLy4	BSk24	BSk22	DD-ME $\delta$	Average	Uncertainty
$E_{sat}$	MeV	0	-15.97	-16.05	-16.09	-16.12	-15.8	0.3
$n_{sat}$	$\text{fm}^{-3}$	1	0.1595	0.1578	0.1578	0.1520	0.155	0.005
$K_{sat}$	MeV	2	230	246	246	219	230	20
$Q_{sat}$	MeV	3	-363	-274.5	-276	-748	300	400
$Z_{sat}$	MeV	4	1587	1184	1190	3950	-500	1000
$E_{sym}$	MeV	0	32.01	30.00	32.00	32.35	32	2
$L_{sym}$	MeV	1	46.0	46.4	68.5	52.8	60	15
$K_{sym}$	MeV	2	-120	-38	13	-118	-100	100
$Q_{sym}$	MeV	3	521	711	563	846	0	400
$Z_{sym}$	MeV	4	-3197	-4031	-3174	-3545	-500	1000
$m_{sat}^*/m$			0.69	0.80	0.80	0.69	0.75	0.1
$\Delta m_{sat}^*/m$			-0.19	0.20	0.20	-0.17	0.1	0.1

Table 1.3: Value of each of the empirical parameters, associated unit, and derivative order  $N$  for SLy4 [Cha+98], BSk24, BSk22 [GCP13], and DD-ME $\delta$  [Roc+11] functionals. Average values and uncertainties extracted from experimental analysis are taken from [MHG18a].

symmetry energy is generally defined as

$$e_{HM}^{sym}(n) = \frac{1}{2} \frac{\partial^2 e_{HM}(n, \delta)}{\partial \delta^2} \Big|_{\delta=0}, \quad (1.31)$$

where  $e_{HM}$  is the nuclear matter energy per nucleon in nuclear matter. In Table 1.3 are listed the value of each empirical parameter for the Skyrme-like SLy4 [Cha+98], BSk22, and BSk24 [GCP13], and relativistic DD-ME $\delta$  [Roc+11] functionals. Average values and associated uncertainties extracted from experimental analysis are also provided. It can be seen that the isovector parameters are in general less known with respect to the isoscalar ones for the same derivative order. For instance, the relative uncertainty on  $K_{sym}$  is about 100% while that of  $K_{sat}$  is less than 10%. It should also be stressed that high-order parameters  $Q_{sat(sym)}$  and  $Z_{sat(sym)}$  are poorly constrained by nuclear experiments.

The empirical parameters can be used in a Taylor expansion to estimate the nuclear matter EoS analytically up to  $n \approx 2 - 3n_{sat}$  [MHG18a; MHG18b]. Limiting us to derivative order  $N = 2$ , we obtain

$$e_{HM}^{N=2}(n, \delta) = E_{sat} + \frac{1}{2} K_{sat} x^2 + \delta^2 (E_{sym} + L_{sym} x + \frac{1}{2} K_{sym} x^2), \quad (1.32)$$

with  $x = (n - n_{sat})/(3n_{sat})$  a function of the density. **However, we should stress already that  $N = 2$  is not sufficient for describing the EoS at subsaturation densities, and even less for suprasaturation densities. High-order parameters  $Q_{sat(sym)}$  and  $Z_{sat(sym)}$  are essential and carry the largest uncertainty as far as the bulk part of the EoS is concerned.**

In a mean field approach, we can treat nucleons as independent particles, thus the kinetic part of the energy per particle of nuclear matter is given by the same functional form as a nonrelativistic Fermi gas (FG), that is

$$t_{HM}^{FG}(n, \delta) = \frac{1}{2} t_{sat}^{FG} (1 + 3x)^{2/3} \left[ (1 + \delta)^{5/3} \frac{m}{m_n^*} + (1 - \delta)^{5/3} \frac{m}{m_p^*} \right], \quad (1.33)$$

where  $t_{sat}^{FG} = 3\hbar^2/(10m)(3\pi^2/2)^{2/3}n_{sat}^{2/3}$  is the kinetic energy of SNM at the saturation density,  $m = (m_n + m_p)/2$  is the mean nucleon mass, and  $m_n^*$  ( $m_p^*$ ) is the neutron (proton) effective mass. The Landau effective mass is introduced in order to take into account for the in-medium nuclear interaction that alters the mass of nucleons. It can be parametrized as ( $q = n, p$ )

$$\frac{m}{m_q^*} = \sum_{\alpha=0}^1 m_{q,\alpha}(\delta) \frac{x^\alpha}{\alpha!} = 1 + (\kappa_{sat} + \tau_3 \kappa_{sym} \delta)(1 + 3x), \quad (1.34)$$

with  $\tau_3 = 1$  ( $\tau_3 = -1$ ) for neutrons (protons). Two additional parameters  $\kappa_{sat}$  and  $\kappa_{sym}$ , related to the effective mass  $m_{sat}^*$  and isospin splitting  $\Delta m_{sat}^*$  at saturation density, are then introduced. They are defined at  $n = n_{sat}$  as

$$\kappa_{sat} = \frac{m}{m_{sat}^*} - 1, \quad (1.35)$$

$$\kappa_{sym} = \frac{1}{2} \left( \frac{m}{m_n^*} - \frac{m}{m_p^*} \right). \quad (1.36)$$

In principle, it is possible to reproduce any EoS model with a Taylor expansion, considering an infinite number of parameters,  $N \rightarrow \infty$ , however the convergence would be very slow. In order to fasten the series convergence, one can add extra functional dependencies, which correspond to the true EoS in the limit of simplistic cases (independent particles, zero density limit), but which allow, by judicious choices of empirical parameters, to reproduce with precision realistic functionals. In this regard, the  $\delta^{5/3}$  dependence is added by decomposing the energy per particle into a potential and kinetic part, yielding

$$e_{HM}^N(x, \delta) = t_{HM}^{FG}(n, \delta) + v_{MM}^N(n, \delta), \quad (1.37)$$

where  $v_{MM}^N$  is the potential energy per particle expressed as a Taylor expansion in the parameter  $x$  at  $n = n_{sat}$ ,

$$v_{MM}^N(n, \delta) = \sum_{\alpha \geq 0}^N (v_\alpha^{is} + \delta^2 v_\alpha^{iv}) \frac{x^\alpha}{\alpha!}. \quad (1.38)$$

The quadratic approximation for the isospin dependence of the potential energy has been made in Eq. (1.38), as suggested by microscopic calculations in [Vid+09]. The coefficients  $v_\alpha^{is}$  and  $v_\alpha^{iv}$  are mapped to the empirical parameters following

$$v_\alpha^{is} = \left. \frac{\partial^\alpha e_{HM}}{\partial x^\alpha} \right|_{n=n_{sat}, \delta=0} \quad \text{and} \quad v_\alpha^{iv} = \left. \frac{\partial^\alpha e_{HM}^{sym}}{\partial x^\alpha} \right|_{n=n_{sat}, \delta=0}, \quad (1.39)$$

yielding the isoscalar parameters:

$$v_0^{is} = E_{sat} - t_{sat}^{FG}(1 + \kappa_{sat}), \quad (1.40)$$

$$v_1^{is} = -t_{sat}^{FG}(2 + 5\kappa_{sat}), \quad (1.41)$$

$$v_2^{is} = K_{sat} - 2t_{sat}^{FG}(-1 + 5\kappa_{sat}), \quad (1.42)$$

$$v_3^{is} = Q_{sat} - 2t_{sat}^{FG}(4 - 5\kappa_{sat}), \quad (1.43)$$

$$v_4^{is} = Z_{sat} - 8t_{sat}^{FG}(-7 + 5\kappa_{sat}), \quad (1.44)$$

and isovector parameters:

$$v_0^{iv} = E_{sym} - \frac{5}{9}t_{sat}^{FG}[1 + (\kappa_{sat} + 3\kappa_{sym})], \quad (1.45)$$

$$v_1^{iv} = L_{sym} - \frac{5}{9}t_{sat}^{FG}[2 + 5(\kappa_{sat} + 3\kappa_{sym})], \quad (1.46)$$

$$v_2^{iv} = K_{sym} - \frac{10}{9}t_{sat}^{FG}[-1 + 5(\kappa_{sat} + 3\kappa_{sym})], \quad (1.47)$$

$$v_3^{iv} = Q_{sym} - \frac{10}{9}t_{sat}^{FG}[4 - 5(\kappa_{sat} + 3\kappa_{sym})], \quad (1.48)$$

$$v_4^{iv} = Z_{sym} - \frac{40}{9}t_{sat}^{FG}[-7 + 5(\kappa_{sat} + 3\kappa_{sym})], \quad (1.49)$$

Let us notice from Eq. (1.38) that the energy per particle does not converge to zero at zero density. This artifact, arising from the nature of the series expansion, can be fixed by introducing an exponential correction  $u_\alpha^N(x)$  in the potential energy,

$$v_{MM}^N(n, \delta) = \sum_{\alpha \geq 0}^N (v_\alpha^{is} + \delta^2 v_\alpha^{iv}) \frac{x^\alpha}{\alpha!} u_\alpha^N(x), \quad (1.50)$$

the expression of which is

$$u_\alpha^N(x) = 1 - (-3x)^{N+1-\alpha} \exp(-b(1+3x)), \quad (1.51)$$

where we have introduced an extra low-density parameter  $b$ . It was shown in [Ant+19] that for most applications this parameter can be fixed to a constant,  $b = 10 \ln(2)$ .

Let us now turn to the derivatives of the energy. The expression of the symmetry energy, Eq. (1.31), reads

$$e_{HM}^{sym}(n) = \frac{5}{9}t_{sat}^{FG}(1+3x)^{2/3} [1 + (\kappa_{sat} + 3\kappa_{sym})(1+3x)] + \sum_{\alpha \geq 0}^N v_\alpha^{iv} \frac{x^\alpha}{\alpha!} u_\alpha^N(x). \quad (1.52)$$

For the chemical potential, we have

$$\mu_{HM,q}(n, \delta) = e_{HM}^N + \frac{1+3x}{3} \left( \frac{\partial e_{HM}^N}{\partial x} \right)_\delta + (\tau_3 - \delta) \left( \frac{\partial e_{HM}^N}{\partial \delta} \right)_x + m_q c^2, \quad (1.53)$$

The complete analytical expression of  $\mu_{HM,q}$  is given in Appendix B. Finally, the pressure can be calculated through

$$P_{HM} = \sum_{q=n,p} n_q (\mu_{HM,q}(n, \delta) - m_q c^2) - n e_{HM}^N(n, \delta). \quad (1.54)$$

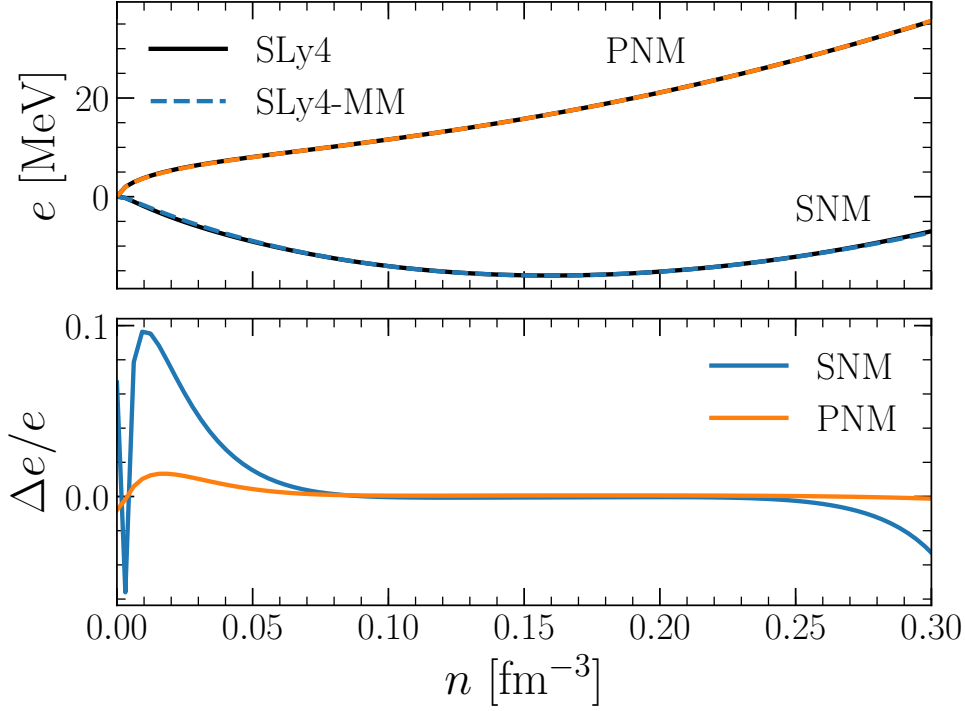


Figure 1.4: Energy per nucleon as a function of density for symmetric nuclear matter,  $\delta = 0$ , and pure neutron matter,  $\delta = 1$  (upper panel), and the associated relative error  $\Delta e/e = (e_{MM} - e)/e$  (lower panel) for the SLy4 functional, with  $N = 4$  for the metamodel. In the top panel, black solid lines (dashed lines) correspond to the calculation with the exact functional (metamodel).

There are two main advantages that come with the metamodeling technique. Firstly, It is flexible enough to reproduce the different existing energy functional of nuclear matter with very good accuracy up to  $n \approx 2 - 3n_{sat}$ , and even higher densities with a proper redefinition of the third and fourth order parameters [MHG18a]. Since very often the parameters of the model to be mimicked do not coincide with the empirical parameters, we derive the latter by calculating the derivatives of the energy per particle in nuclear matter at saturation density, following Eq. (1.39). In Fig. 1.4 is demonstrated the accuracy of the metamodeling technique for the SLy4 [Cha+98] functional with a Taylor expansion order  $N = 4$ . In the top panel, the energy per nucleon is calculated using the metamodeling technique (dashed lines) and the exact functional (black solid lines). A tiny deviation is observed at low density for the SNM, but overall a very good agreement is observed between the metamodel (labeled “SLy4-MM”) and the exact functional (labeled “SLy4”). In the lower panel the relative error is represented for SNM and PNM. We can see that the error goes up to 10% for SNM at  $n \approx 0.01 \text{ fm}^{-3}$  and rapidly drops to 0% at  $n \approx 0.07 \text{ fm}^{-3}$ . At  $n \approx 2n_{sat}$ , a deviation appears due the extrapolation far from the saturation point. The deviation at low density can be reduced

by adjusting the parameter  $b$  entering in the low-density correction, Eq. (1.51) [Ant+19]. However, this small deviation is not expected to cause any issue for the determination of the inner crust ground state. Indeed, the relative error for PNM is lower than 2% at most, and, in the inner crust, the matter inside nuclei is very neutron rich and the ambient nucleon gas is exclusively constituted of free neutrons. **This ability of mimicking existing realistic functionals will allow us to easily compare predictions of astrophysical observables of different popular models.**

Another asset of the metamodeling technique is that no correlation is assumed a priori among the empirical parameters due to the nature of the series expansion. Therefore, all density dependences of the nuclear EoS can be explored. In addition, it allows us to vary each empirical parameter independently of the others to carry out complete statistical analysis using Bayesian inference.

### 1.2.1.2 From homogeneous nuclear matter to finite nuclei in the CLD approximation

We now turn to the modeling of the energy of clusters present in the inner crust of cold neutron stars. We propose to extend the metamodeling of infinite nuclear matter to nuclei in order to take into account their surface properties, within the CLD approximation, originally introduced in [BBP71]. This approach is expected to be less accurate than microscopic Hartree-Fock (HF) [NV73] or (ETF) [Ons+08] calculations since quantum effects are neglected, though it proved its quality in the past by giving surprisingly good results, see for example [BBP71; LS91; DH00; DH01]. In particular, it appears that most of the macroscopic observables related to neutron stars, such as masses and radii, are not very sensitive to the crust composition, thus a qualitative evaluation of it is sufficient in that case. The compressible liquid drop model (CLDM) comes with multiple advantages. It is very fast from the computational point of view since the energy does not require to be integrated all along the cell volume. For this reason, it is powerful for complete statistical analyses such as the one presented in the next chapter. It is also a model that is **flexible**, giving us the opportunity to account for different geometries for example, as we consider in 1.2.5. Last but not least, because of the artificial decomposition of the nuclear cluster as a bulk term, a surface term, and a Coulomb term, we are able to identify the contributions of the different terms. For all these reasons, the CLDM is considered as a suitable alternative to semiclassical ETF and microscopic HF treatments, that additionally does not suffer from technical complications such as the definition of boundary conditions [NS09].

In the CLD approximation, the energy of the cluster reads

$$E_{cl} = e_{HM}(n_0, I)A + E_{surf} + E_{Coul}, \quad (1.55)$$

where  $E_{surf}$  represents the nuclear surface energy, and  $E_{Coul}$  is the Coulomb energy, including the lattice correction and finite-size effects, written as follows in the WS approximation,

$$E_{Coul} = \frac{3}{5} \frac{e^2}{r_0^2} \eta_{Coul}(u) \frac{Z^2}{A^{1/3}} = \frac{3}{20} \frac{e^2}{r_0^2} \eta_{Coul}(u) A^{5/3} (1 - I)^2, \quad (1.56)$$



with

$$\eta_{Coul}(u) = 1 - \frac{3}{2}u^{1/3} + \frac{1}{2}u, \quad (1.57)$$

a function of the volume fraction  $u = n_e/n_{0,p}$ , with  $n_{0,p} = n_0(1 - I)/2$  the average proton density inside cluster. Let us notice that, the factor  $3/5 \times 3/2 = 9/10$  entering in the lattice term can be replaced by the Mandelung constant for a body-centered cubic lattice,  $C_M = 0.895929255682$  (see Table 2.4 of [HPY07]).

The bulk term is taken from the metamodel, keeping the same empirical parameters as for the calculation of the neutron gas energy, thus insuring consistency between the crust and the core treatment. The bulk energy is evaluated at  $n_0$ , the average density inside the cluster, and  $I$ , its global asymmetry.

Let us now turn to the expression of the nuclear surface energy. Assuming that clusters are spherical, it is given by

$$E_{surf} = 4\pi r_0^2 A^{2/3} \sigma, \quad (1.58)$$

where  $r_0 = (4\pi n_0/3)^{-1/3}$  is related to the cluster density  $n_0$ , and  $\sigma$  is the nuclear surface tension. The simplest parametrization of the surface tension one could use is a constant,  $\sigma \approx 1.03$  MeV/fm<sup>2</sup>, as in the liquid drop model. However, clusters in the inner crust are expected to be very neutron rich and it appears evident that the surface tension of very asymmetric nuclei should be different to that of symmetric nuclei. In addition, it does not account for the interaction of the nuclear surface with the outside neutron gas, yielding  $\sigma \rightarrow 0$  as  $I \rightarrow 1$ . This leads us to consider another parametrization for  $\sigma$  which would depend on the asymmetry  $I$  and consider the interaction with the gas. It should be mentioned that the exact value of  $\sigma(I)$  is model dependent, and the uncertainty is particularly important for extreme isospin values encountered in the inner crust. Indeed, no experimental data exist on the surface energy of nuclei beyond the dripline, which is in-medium modified by the presence of the neutron gas [DHM00]. We use the expression originally proposed by Ravenhall *et al.* [RPL83] on the basis of TF calculations at extreme isospin ratios, and later employed in different works on neutron star crust and supernovae modeling within the CLD approximation [LS91; NGL12; LRP93]:

$$\sigma(I) = \sigma_0 \frac{2^{p+1} + b_s}{Y_p^{-p} + b_s + (1 - Y_p)^{-p}}, \quad (1.59)$$

with  $\mathbf{S} = \{\sigma_0, b_s, p\}$  the parameter space associated to the surface energy. The parameter  $\sigma_0 = \sigma(I = 0)$  represents the surface tension of symmetric nuclei, and  $b_s$  and  $p$  are the parameters that govern the isospin dependence of the surface tension. The proton fraction is related to the asymmetry inside the cluster via  $Y_p = (1 - I)/2$ . Following [NGL12], we add a curvature term with the aim of having a better description of the nuclear surface and so of the composition of the inner crust in the ground state. For spherical clusters, the curvature energy reads

$$E_{curv} = 8\pi r_0 A^{1/3} \sigma_c, \quad (1.60)$$

where  $\sigma_c$  is the curvature tension related to the surface tension  $\sigma$ , Eq. (1.59),

$$\sigma_c = \sigma \frac{\sigma_{0,c}}{\sigma_0} \alpha(\beta - Y_p), \quad (1.61)$$

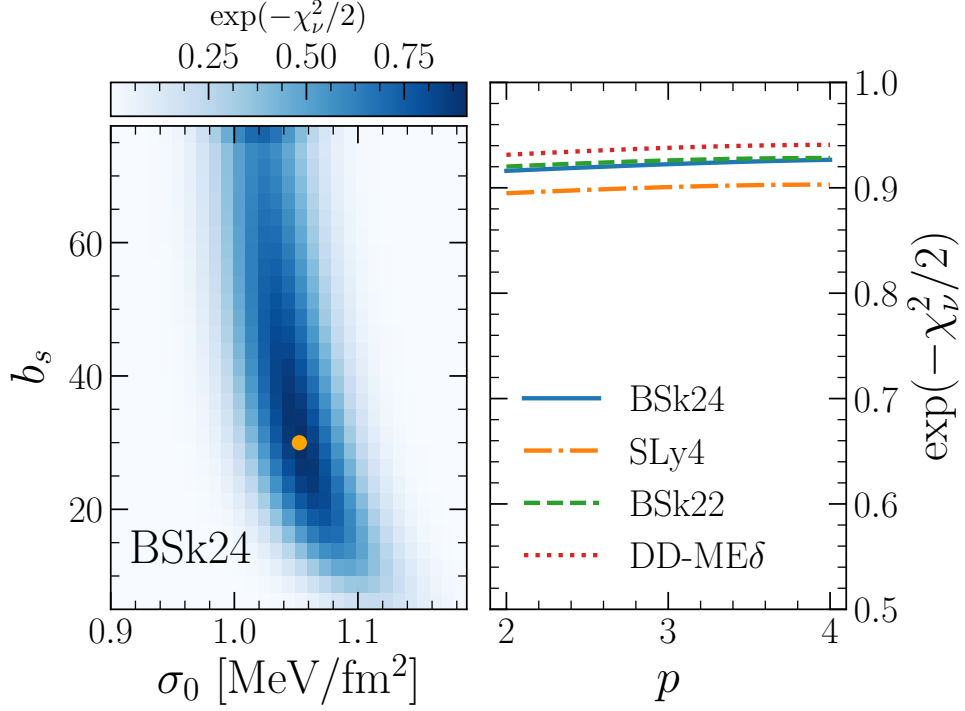


Figure 1.5: Left: Goodness of fit (color scale) as described by the reduced  $\chi^2$ , in the space of the surface parameters  $(\sigma_0, b_s)$  for the BSk24 CLDM with fixed value of the  $p$  parameter,  $p = 3$ , for optimized curvature parameters  $\sigma_{0c}$  and  $\beta$ . The orange point indicates the optimal value. Right: Weight  $\exp(-\chi^2_\nu/2)$  as a function of the parameter  $p$ , calculated for the optimal value of  $(\sigma_0, b_s, \sigma_{0c}, \beta)$ , for the selected CLDM: BSk24, SLy4, BSk22, and DD-ME $\delta$ .

with  $\alpha = 5.5$  as in [NGL12]. With the addition of the curvature energy comes two parameters added as extra dimensions to the surface parameter space,  $\mathbf{S} = \{\sigma_0, b_s, \sigma_{0c}, \beta, p\}$ . For a given model of uniform matter, that is each given set of empirical parameters, the first four parameters can be fitted to experimental masses from the AME2016 [Hua+17] for a fixed value of the parameter  $p$ . The quality of reproduction of the experimental nuclear binding energy  $E_{\text{AME2016}}$  is measured by the reduced  $\chi^2$ , defined as

$$\chi^2_\nu = \frac{\chi^2}{\nu} = \frac{1}{\nu} \sum_i \left( \frac{E_{cl}^{(i)}/A^{(i)} - E_{\text{AME2016}}^{(i)}/A^{(i)}}{\sigma^{(i)}} \right)^2, \quad (1.62)$$

where  $\nu$  is the number of degrees of freedom, and the denominator corresponds to the systematic theoretical error, 0.1 MeV/ $A$ , such as to have  $\chi^2_{\min} \approx 1$  over the parameter space sample [DNR14].

This point can be clearly understood by inspection of the left panel of Fig. 1.5, which shows the value of the weight function  $\exp(-\chi^2_\nu/2)$ , that is the goodness of fit, in the  $(\sigma_0, b_s)$  plane, for the BSk24 CLDM with  $p = 3$ , and curvature parameters already ad-



Figure 1.6: Variation with cluster asymmetry  $I$  of surface plus curvature energy per surface nucleon along the isobaric chain  $A = 200$  for SLy4 [Cha+98] and DD-ME $\delta$  [Roc+11] CLDM at three selected values of parameter  $p$ :  $p = 2$  (dashed lines),  $p = 3$  (solid lines), and  $p = 4$  (dotted lines).

justed to experimental data. It is seen that the weight function is strongly peaked to the orange point, which indicates the optimal value for the parameters  $\sigma_0$  and  $b_s$ . Conversely, we observe in the right panel of Fig. 1.5 that the error estimator  $\chi^2_\nu$  is insensitive to the choice of the parameter  $p$  in a range going from 2 to 4, for four representative functionals. This is understood from the fact that  $p$  modifies the asymmetry dependence of the surface energy only for extreme values of  $I$  where nuclear data are not available. This parameter is **therefore** not adjusted as the other parameters, in order to avoid poor extrapolations of  $\sigma$ . The value of  $p$  is selected around  $p = 3$ , which is the value used in the popular Lattimer and Swesty EoS [LS91]. We show in Fig. 1.6 the surface plus curvature energy per surface nucleon as a function of the asymmetry inside the cluster for two popular models, SLy4 [Cha+98] and DD-ME $\delta$  [Roc+11], in the CLD approximation. Different values for the parameter  $p$  are displayed in that figure. It is seen that  $p$  determines the behavior of the surface tension for extreme isospin values, and cannot be accessed from experimental nuclear physics data. **In addition, since the selected popular models contain parameters that are adjusted to the same data, we cannot consider their predictions at high isospin as reliable. For these reasons, we will keep  $p$  as a free parameter.** In particular, from  $I \approx 0.3$ , we observe that a larger value of  $p$  leads to larger surface energy at fixed asymmetry. The model dependence of the surface tension is also observed in that figure and is particularly highlighted at low asymmetry

where experimental measurements can still be performed.

Finally, we recall the final expression for the cluster binding energy in the CLD approximation,

$$\begin{aligned}
 E_{cl}(A, I, n_0, n_e) = & e_{HM}(n_0, I)A \\
 & + 4\pi r_0^2 A^{2/3} \sigma(I) + 8\pi r_0 A^{1/3} \sigma_c(I) \\
 & + \frac{3}{20} \frac{e^2}{r_0^2} \eta_{Coul}(I, n_0, n_e) A^{5/3} (1 - I)^2.
 \end{aligned} \tag{1.63}$$

Let us notice that the energy does not depend on the neutron gas density via the surface term, as in [BBP71]. This dependence is in fact implicitly accounted for in the chosen parametrization of the surface tension, suggested from microscopic calculations in the WS cell [RPL83].

### 1.2.2 Variational formalism

The variational method for the calculation of the inner crust ground state with the CLDM was introduced in the pioneering work of BBP in their classical paper [BBP71]. Since then, huge progress has been made in constraining the nucleon-nucleon (NN) effective interaction. For this reason, subsequent works using the same formalism with CLDM based on more realistic nuclear functionals followed. For instance we can mention the popular calculations of Douchin and Haensel (DH) [DH00; DHM00] with the SLy4 effective force [Cha+98]. Following those works, we use the same variational method as BBP, and we present it in the following.

Using the Lagrange multipliers method, the ground state of the inner crust is obtained by minimizing the zero-temperature free energy density of the WS cell at fixed baryon density  $n_B$  under the condition of charge neutrality,  $n_e = n_p$ . The minimization is carried out with respect to the set of variables  $A$ ,  $I$ ,  $n_0$ ,  $n_p$ , and  $n_g$ . Let us notice that, in principle, one should minimize the Gibbs free energy at constant pressure, as for the outer crust. However, it would be very heavy from the computational point of view to do so in the free neutron regime. Fortunately, it is demonstrated that the error introduced by minimizing at constant baryon density instead of pressure is negligible [Pea+12]. The auxiliary function to be minimized is defined as

$$\begin{aligned}
 \mathcal{L}(A, I, n_0, n_p, n_g) &= \frac{E_{WS}}{V_{WS}} - \mu n_B \\
 &= \frac{E_{cl}}{V_{WS}} + \varepsilon_e + \left(1 - \frac{A}{n_0 V_{WS}}\right) \varepsilon_g \\
 &\quad + n_p \Delta m_{pn} c^2 - n_B (\mu - m_n c^2),
 \end{aligned} \tag{1.64}$$

with  $\Delta m_{pn} = m_p - m_n$  the difference between the proton and neutron mass, and  $\mu$  the Lagrange multiplier.

The volume of the cell can be calculated as

$$V_{WS} = \frac{Z}{n_p} = \frac{A}{n_p} \frac{1 - I}{2}, \tag{1.65}$$

and the baryon density is given by

$$\begin{aligned} n_B &= \frac{A_{WS}}{V_{WS}} \\ &= \frac{2n_p}{1-I} \left(1 - \frac{n_g}{n_0}\right) + n_g. \end{aligned} \quad (1.66)$$

Replacing these quantities in Eq. (1.64), it results

$$\begin{aligned} \mathcal{L}(A, I, n_0, n_p, n_g) &= \frac{2n_p}{A(1-I)} E_{cl} + \varepsilon_e + \left(1 - \frac{2n_p}{n_0(1-I)}\right) \varepsilon_g + n_p \Delta m_{pn} c^2 \\ &\quad - \frac{2n_p}{1-I} \left(1 - \frac{n_g}{n_0}\right) (\mu - m_n c^2) - n_g (\mu - m_n c^2). \end{aligned} \quad (1.67)$$

Let us first minimize with respect to  $A$ :

$$\begin{aligned} \left. \frac{\partial \mathcal{L}}{\partial A} \right|_{I, n_0, n_p, n_g} &= 0 \\ \frac{2n_p}{A(1-I)} \frac{\partial E_{cl}}{\partial A} - \frac{2n_p}{A^2(1-I)} E_{cl} &= 0 \\ \frac{\partial E_{cl}}{\partial A} - \frac{E_{cl}}{A} &= 0, \end{aligned} \quad (1.68)$$

which is equivalent to

$$\frac{\partial(E_{cl}/A)}{\partial A} = 0. \quad (1.69)$$

We can go through this equation using the definition of the cluster energy, Eq. (1.63), yielding

$$\begin{aligned} -\frac{1}{3} \times 4\pi r_0^2 \sigma A^{-4/3} - \frac{2}{3} \times 8\pi r_0 \sigma_c A^{-5/3} + \frac{2}{3} \times \frac{3}{20} \frac{e^2}{r_0} \eta_{Coul} A^{-1/3} (1-I)^2 &= 0 \\ 4\pi r_0^2 \sigma A^{2/3} + 2 \times 8\pi r_0 \sigma_c A^{1/3} - 2 \times \frac{3}{20} \frac{e^2}{r_0} \eta_{Coul} A^{5/3} (1-I)^2 &= 0, \end{aligned}$$

or simply,

$$E_{surf} + 2E_{curv} = 2E_{Coul}. \quad (1.70)$$

This equation corresponds to the well-known Baym virial theorem with an additional curvature term with respect to the equation originally found in [BBP71]. It is interesting to notice that only the surface, the curvature, and the Coulomb energy are involved in this first equilibrium condition.

Minimizing with respect to the neutron gas density  $n_g$ , one obtains:

$$\begin{aligned} \left. \frac{\partial \mathcal{L}}{\partial n_g} \right|_{A, I, n_0, n_p} &= 0 \\ \left(1 - \frac{2n_p}{n_0(1-I)}\right) \frac{\partial \varepsilon_g}{\partial n_g} + \frac{2n_p}{n_0(1-I)} (\mu - m_n c^2) - (\mu - m_n c^2) &= 0, \end{aligned} \quad (1.71)$$

giving us the expression of the Lagrange multiplier, which can be identified with the chemical potential of the gas, including the rest mass energy,

$$\mu = \frac{\partial \varepsilon_g}{\partial n_g} + m_n c^2 \equiv \mu_g. \quad (1.72)$$

We now turn to the minimization with respect to average cluster density  $n_0$ :

$$\begin{aligned} \frac{\partial \mathcal{L}}{\partial n_0} \Big|_{A, I, n_p, n_g} &= 0 \\ \frac{2n_p}{A(1-I)} \frac{\partial E_{cl}}{\partial n_0} + \frac{2n_p}{n_0^2(1-I)} \varepsilon_g - \frac{2n_p n_g}{n_0^2(1-I)} (\mu - m_n c^2) &= 0 \\ \frac{n_0^2}{A} \frac{\partial E_{cl}}{\partial n_0} - n_g (\mu - m_n c^2) + \varepsilon_g &= 0. \end{aligned} \quad (1.73)$$

Using well-known thermodynamical relations, we define  $P_{cl} \equiv (n_0^2/A)(\partial E_{cl}/\partial n_0)$ , and  $P_g = n_g(\mu_g - m_n c^2) - \varepsilon_g$ . Therefore, Eq. (1.73) simply results in

$$P_{cl} = P_g, \quad (1.74)$$

which can be interpreted as a pressure equilibrium condition between the cluster and the outside neutron gas. We can derive the expression of the cluster pressure from the CLD energy, Eq. (1.63), yielding

$$P_{cl} = P_{HM}(n_0, I) - \frac{2}{3} n_0 \frac{E_{surf}}{A} - \frac{1}{3} n_0 \frac{E_{curv}}{A} + n_0 \frac{E_{coul}(u=0)}{A} \left( \frac{2}{3} + \frac{1}{2} u^{1/3} - \frac{1}{2} u \right). \quad (1.75)$$

The minimization with respect to the global asymmetry inside the cluster  $I$  yields

$$\begin{aligned} \frac{\partial \mathcal{L}}{\partial I} \Big|_{A, n_0, n_p, n_g} &= 0 \\ \frac{2n_p}{(1-I)^2} \left( \frac{E_{cl}}{A} + \frac{1-I}{A} \frac{\partial E_{cl}}{\partial I} - \frac{\varepsilon_g}{n_0} - \left( 1 - \frac{n_g}{n_0} \right) (\mu - m_n c^2) \right) &= 0 \\ \frac{E_{cl}}{A} + \frac{1-I}{A} \frac{\partial E_{cl}}{\partial I} - \frac{\varepsilon_g}{n_0} - \left( 1 - \frac{n_g}{n_0} \right) (\mu - m_n c^2) &= 0, \end{aligned} \quad (1.76)$$

resulting in

$$\frac{E_{cl}}{A} + \frac{1-I}{A} \frac{\partial E_{cl}}{\partial I} + \frac{P_g}{n_0} = \mu - m_n c^2. \quad (1.77)$$

Due to the presence of the outside neutron gas, the neutron chemical potential of the cluster is modified with respect to the usual expression in the vacuum  $\mu_n^{vac} = \partial E_{cl}/\partial N$ . We define the neutron and proton chemical potential of the cluster as

$$\mu_p^{cl} = \frac{\partial E_{cl}}{\partial Z} \Big|_N + m_p c^2 \quad \text{and} \quad \mu_n^{cl} = \frac{\partial E_{cl}}{\partial N} \Big|_Z + \frac{P_g}{n_0} + m_n c^2, \quad (1.78)$$

where the derivatives are taken at constant  $n_0$ ,  $n_p$ , and  $n_g$ . Consequently, we can work out Eq. (1.77) to find a chemical equilibrium conditions between the neutrons of the cluster and those of the gas,

$$\mu_n^{cl} = \mu_g. \quad (1.79)$$

Finally, we minimize the auxiliary function, Eq. (1.64) with respect to the proton density  $n_p$  inside the WS cell,

$$\begin{aligned} \frac{\partial \mathcal{L}}{\partial n_p} \Big|_{A, I, n_0, n_g} &= 0 \\ \frac{E_{cl}}{A} + \frac{n_p}{A} \frac{\partial E_{cl}}{\partial n_p} + \frac{1-I}{2} \left( \frac{\partial \varepsilon_e}{\partial n_p} + \Delta m_{pn} c^2 \right) - \frac{\varepsilon_g}{n_0} - \left( 1 - \frac{n_g}{n_0} \right) (\mu - m_n c^2) &= 0. \end{aligned} \quad (1.80)$$

By identifying common terms with Eq. (1.76), we recover the well-known beta equilibrium condition,

$$\frac{2}{A} \left( \frac{\partial E_{cl}}{\partial I} - \frac{n_p}{1-I} \frac{\partial E_{cl}}{\partial n_p} \right) - \Delta m_{pn} c^2 = \frac{\partial \varepsilon_e}{\partial n_p}, \quad (1.81)$$

or in terms of chemical potentials,

$$\mu_n^{cl} = \mu_p^{cl} + \mu_e + \Delta\mu, \quad (1.82)$$

where  $\Delta\mu$  represents the modification due to the excluded volume interaction with the neutron gas, and the electrostatic interaction between protons in the cluster and the background electrons. It is given by

$$\Delta\mu = \frac{P_g}{n_0} + \frac{2}{A(1-I)} n_p \frac{\partial E_{Coul}}{\partial n_p}. \quad (1.83)$$

One can remark that only the Coulomb energy enters in the derivative with respect to the proton density  $n_p$ . This derivative can be evaluated analytically from Eq. (1.56) as

$$\frac{\partial E_{Coul}}{\partial n_p} = \frac{3}{20} \frac{e^2}{r_0^2} A^{5/3} (1-I) \frac{1}{n_0} \left( 1 - \left( \frac{2n_p}{n_0(1-I)} \right)^{-2/3} \right). \quad (1.84)$$

Finally, a possible set of mechanical and chemical equilibrium equations is

$$\frac{\partial (E_{cl}/A)}{\partial A} = 0, \quad (1.85)$$

$$\frac{n_0^2}{A} \frac{\partial E_{cl}}{\partial n_0} = P_g(n_g), \quad (1.86)$$

$$\frac{E_{cl}}{A} + \frac{1-I}{A} \frac{\partial E_{cl}}{\partial I} + \frac{P_g(n_g)}{n_0} = \mu_g(n_g) - m_n c^2, \quad (1.87)$$

$$\frac{2}{A} \left( \frac{\partial E_{cl}}{\partial I} - \frac{n_p}{1-I} \frac{\partial E_{cl}}{\partial n_p} \right) - \Delta m_{pn} c^2 = \mu_e(n_p), \quad (1.88)$$

or, equivalently Eqs. (1.70), (1.74), (1.79), and (1.82), respectively. This system of four coupled differential equations can be solved numerically, as explained in 1.2.3.1, in order to obtain the equilibrium composition and so the EoS in the free neutron regime. Let us mention that taking alternatives set of variables leads to the same equilibrium equations [HPY07; Viñ+17].

### 1.2.3 Results

We now turn to the results of the minimization. The building of the numerical code is presented first and the numerical results, such as the ground-state composition and EoS of the inner crust, afterwards.

#### 1.2.3.1 Numerical code

During the thesis, an open-source C library, NSEoS, has been built in relation to the physics of NS [Car17a]. Here we focus on the numerical methods used to solve the inner crust equilibrium equations in order to estimate the composition and EoS in the free neutron regime. Those methods can be found in the module `crust` of the library.

The ground state of the inner crust at baryon density  $n_B$  is determined by solving numerically the system of four coupled differential equations, Eqs. (1.85), (1.86), (1.87), and (1.88). To do so, we use a root-finding algorithm, Broyden's method [Bro65], which is similar to the popular Newton's method. In Broyden's method, the 16 partial derivatives entering in the  $4 \times 4$  Jacobian matrix are replaced with finite differences. While the derivatives can be evaluated analytically from the CLD expressions with ease, using finite differences makes it **easier to identify the effect of the different contributions entering in the WS cell energy. This point will be covered in greater detail in Chapter 3.** The Jacobian matrix is not computed at each iteration but is only evaluated at the first iteration then simply rank-one updated. In order to avoid an unreliable estimation of the equilibrium composition  $\{A_{eq}, I_{eq}, n_{0,eq}, n_{p,eq}, n_{g,eq}\}$  due to the accumulation of errors, the maximum number of iterations is fixed to  $N_{iter}^{max} = 1000$ . In addition, the backstepping technique is employed to avoid nonphysical solutions, such as negative values for  $A$  or  $|I| > 1$ , that could result in errors when reinjected in the equilibrium equations. For instance, if at a given iteration, a spurious value of  $A$  is obtained, then we replace it by the previous value  $A_{old}$  plus a smaller step  $\Delta A' < \Delta A$ , until we get a physical value. Finally, let us mention that as for Newton's method, Broyden's method requires an initial guess for the composition.

We proceed as follows. The calculation starts at the neutron drip density  $n_{ND}$ . **A natural choice for the initial guess is to take** the last solution for the ground state of the outer crust, with reasonable values  $A_{guess} = 36$ ,  $I_{guess} = 0.3$ ,  $n_{0,guess} = 0.15 \text{ fm}^{-3}$ , and  $n_{g,guess} = 10^{-6} \text{ fm}^{-3}$ . Then, the proton density is calculated as

$$n_{p,guess} = (n_B - n_{g,guess}) \left( 1 - \frac{n_{g,guess}}{n_{0,guess}} \right) \frac{1 - I_{guess}}{2}. \quad (1.89)$$

To continue, we make use of Broyden's method described previously to evaluate the equilibrium composition  $A_{eq}$ ,  $I_{eq}$ ,  $n_{0,eq}$ , and  $n_{g,eq}$ . The process is then repeated with  $n_B$  increasing in steps of  $\Delta n_B = 10^{-4} \text{ fm}^{-3}$ . At each new value of  $n_B$ , the guess is updated to coincide with the last equilibrium composition, ensuring faster convergence of the algorithm. At a reasonable value of baryon density,  $n_B = 10^{-3} \text{ fm}^{-3}$ , we start to check whether it is beneficial or not to have a phase transition from clusterized matter to homogeneous matter. The calculation is stopped once the crust-core (CC) interface is reached, approximately for  $n_B \approx n_{sat}/2$ . The CC transition from the crust side is discussed in details in 1.2.6.





Figure 1.7: Variation with baryon density  $n_B$  of the equilibrium value of mass number  $A$ , atomic number  $Z$ , proton fraction  $Y_p$ , average cluster density  $n_0$ , neutron gas density to baryon density ratio  $n_g/n_B$ , radius  $R$ , and volume fraction  $u$  in the inner crust, for BSk24 CLDM with  $p = 3$ . Solid (dotted) lines correspond to the results for the cell (cluster). In the upper-mid panel, three values of the parameter  $p$  are selected:  $p = 2.5$ ,  $p = 3$ , and  $p = 3.5$ .

### 1.2.3.2 Equilibrium composition

Fig. 1.7 shows the evolution with baryon density  $n_B$  of the equilibrium composition of the inner crust. The empirical parameters entering in the bulk part of the CLDM energy correspond to the BSk24 effective force [GCP13], and the surface parameter  $p$  is fixed at the educated value  $p = 3$ , which appears to be a reasonable value for Skyrme-type functionals, as we will discuss in 1.2.6.

It is found that the number of nucleons inside the cluster  $A$  as well as inside the WS cell  $A_{WS}$  monotonously increases with increasing density, going up to  $\approx 700$  and  $\approx 1500$ , respectively, at high density. The number of free neutrons  $N_g = A_{WS} - A$  can also be inferred. It is obviously equal to zero at the neutron drip density, since  $A = A_{WS}$ . The behavior of  $A$  and  $A_{WS}$  is in agreement with other CLD calculations, such as the one of BBP [BBP71] and that of DH [DH00]. The later study reports a value of  $A \approx 600$  at the CC interface, using the SLy4 functional, while BBP report a much higher value of  $A = 7840$  (see their Table 1). Let us notice that the presence of such heavy spherical nuclei at the crust bottom raises the question of stability with respect to deformation and

fission, as explained in [DH00]. In particular, nuclear fission, the approximate condition for which is  $R_p/R_{WS} \gtrsim 1/2$  [BW39; PR95],  $R_p$  being the proton radius, could occur at high density. In 1.2.5, we show that considering nonspherical geometries can stabilize the large clusters. It is important to stress that the composition, in particular  $A$  and  $Z$ , is very sensitive to the isospin dependence of the surface tension at extreme isospin ratios.

The variation with baryon density of the equilibrium value of  $Z$  is also represented in Fig. 1.7. We find  $Z \approx 40$  all along the inner crust with  $p = 3$ , which is in a surprisingly good agreement with microscopic results. In particular, in the early work of Negele and Vautherin (NV) [NV73] is derived a set of nonlinear equations for the single particle wave functions of nucleons, using the density-matrix expansion for a realistic NN interaction. NV observed a predominance of  $Z = 40$  at low density, and  $Z = 50$  at high density. Subsequent works within the semiclassical ETF [Gor+05] and finite-temperature extended Thomas-Fermi plus Strutinsky integral (TETFSI) [Ons+08; Pea+18] also predict  $Z = 40$  in the inner crust. As far as CLD calculations are concerned, DH find that the equilibrium value of  $Z$  is almost constant in the vicinity of  $\approx 40$  throughout the inner crust, whereas BBP find monotonically increasing values of  $Z$ . Once again, it should be stress out that  $Z$  is very sensitive to the isospin dependence of the surface tension. In particular, it is observed by varying the parameter  $p$  that governs the behavior of the surface tension at extreme isospin values. We can see that a lower surface tension,  $p = 3.5$ , favors smaller atomic number  $Z$ , and in general lighter clusters, unlike  $p = 2.5$  that leads to heavier clusters.

As in the outer crust, we find that the proton fraction inside the cell  $Y_p$  continuously decreases in the free neutron regime, dropping from  $\approx 0.3$  at the neutron drip point to  $\approx 0.03$  at the edge of the crust. With increasing density, clusters also becomes more and more asymmetric,  $I$  going up to  $\approx 0.85$  in the bottom layers.

Clusters are found to be more and more dilute with  $n_B$  increasing. At the neutron drip density, we get  $n_0 \approx 0.145 \text{ fm}^{-3}$ , which is close to the saturation density of SNM of the BSk24 functional,  $n_{sat} = 0.1578 \text{ fm}^{-3}$ . At the CC interface, the cluster density is as low as  $n_0 \approx 0.07 \text{ fm}^{-3}$ , which is comparable to the neutron gas density, represented in the lower-mid panel. It is seen that the ratio  $n_g/n_B$  rapidly reaches  $\approx 75\%$  at  $n_B = 2 \times 10^{-3} \text{ fm}^{-3}$  then stays approximately constant at higher densities.

The equilibrium radius of the spherical WS cell  $R_{WS}$  and of the spherical cluster  $R_{cl}$  are also displayed in Fig. 1.7. As DH, we observe that  $R_{WS}$  monotonically decreases unlike  $R_{cl}$  that slowly increases with increasing depth. Therefore, nuclei becomes closer and closer, eventually becoming close enough to touch each other at some point, creating a very large cluster, ultimately leading to homogeneous matter. We can also understand this by calculating the volume fraction  $u = (R_{cl}/R_{WS})^3 = 2n_p/(n_0(1 - I))$  (orange dashed line). We find that the cluster fills up to 60% of the WS cell volume at the CC interface. This feature combined with the virial theorem, Eq. (1.70), outlines the fact that lattice and finite-size contributions to the Coulomb energy are crucial for the determination of the ground state of the inner crust.



Figure 1.8: Left: Equilibrium value of the energy per nucleon of the WS cell  $e_{WS} = E_{WS}/A_{WS}$  (upper panel), and pressure  $P$  (lower panel) as a function of the baryon density  $n_B$  in the inner crust for four selected CLDM: BSk24, SLy4, BSk26, and DD-ME $\delta$ . Right: Variation of symmetry energy  $e_{HM}^{sym}$ , Eq. (1.52), with density  $n$ , for the four models.

### 1.2.3.3 Equation of state

In Fig. 1.8, the equilibrium energy per nucleon of the WS cell  $e_{WS} = E_{WS}/A_{WS}$  is plotted as a function of the baryon density  $n_B$  for different CLDM, based on the relativistic model DD-ME $\delta$ , and on three Skyrme-type interactions: SLy4, BSk22, and BSk24. It is observed that  $e_{WS}$  depends on the nuclear model. In particular, it is known to be correlated with the symmetry energy, with higher symmetry leading to higher  $e_{WS}$ , at subnuclear densities [Pea+18]. This trend can be verified by looking at the right panel of the figure, that shows the symmetry energy, Eq. (1.52), as a function of density for the four models considered here.

Fig. 1.8 also shows the variation with the baryon density of pressure  $P$ , commonly referred to as the EoS, in the inner crust. We demonstrate in Appendix C that the total pressure, defined by the first law of thermodynamics as

$$P = n_B^2 \frac{de_{WS}}{dn_B}, \quad (1.90)$$

can be expressed as

$$P = P_g + P_e + P_L, \quad (1.91)$$

at the equilibrium composition, where  $P_L$  is the lattice contribution to the total pressure. Using this expression rather than finite differences to estimate the derivative, Eq. (1.90), is more reliable and has the advantage to be faster from the computational point of view. In addition, it allows identifying the different contributions to the pressure. In the inner crust, the pressure is no longer dominated by the relativistic electron gas but by the ambient neutron gas [Car17b]. At subnuclear densities, we observe that the EoS is also correlated with the symmetry energy. In particular, the higher the empirical parameter  $E_{sym}$ , which is the symmetry energy at  $n = n_{sat}$ , the stiffer is the EoS. The value of  $E_{sym}$  for each model is reported in Table 1.3.

#### 1.2.4 Strutinsky shell corrections to the CLD energy

In the CLD approximation, shell effects, which are known since the pioneering work of BPS [BPS71] to be essential to correctly evaluate the outer-crust composition at zero temperature, are lost. In this same limit microscopic calculations have shown that the neutron shell effects become vanishingly small beyond the neutron drip point [Cha06; Cha+07], but proton shell effects persist in the inner crust.

There are different ways to include shell corrections to the CLD energy. The simplest approach is to use an empirical formula, such as the one proposed by Myers and Swiatecki [MS66] which is accurate for reproducing masses of laboratory nuclei. However, a limitation of their method is that it requires a priori a list of magic numbers, which are obviously not known in the regime of the inner crust. Let us recall that NV find a persistence of  $Z = 40$  while it is not considered as a magic number in ordinary nuclei. A more suitable alternative is to calculate shell corrections using the Strutinsky method [Ons+08]. Fortunately, Strutinsky shell corrections were calculated for modern BSk functionals, and are tabulated in [Pea+18] (see Supplementary data). We therefore add these corrections on top of the CLD energy for the corresponding models.

We now turn to the numerical method, proceeding as follows. The energy density is the quantity to be minimized at constant baryon density for a fixed number of protons inside the WS cell. We derive the system of equilibrium equations using the Lagrange multipliers method, as in 1.2.2, yielding

$$\frac{\partial E_{cl}}{\partial A} - \frac{E_{cl}}{A} = \frac{1 - I}{2} \left( \mu_e(n_p) + \frac{2np}{A(1 - I)} \frac{\partial E_{cl}}{\partial n_p} - \frac{2}{A} \frac{\partial E_{cl}}{\partial I} + \Delta m_{pn} c^2 \right), \quad (1.92)$$

$$\frac{\partial E_{cl}}{\partial A} + \frac{1 - I}{A} \frac{\partial E_{cl}}{\partial I} + \frac{P_g(n_g)}{n_0} = \mu_g(n_g) - m_n c^2, \quad (1.93)$$

$$\frac{n_0^2}{A} \frac{\partial E_{cl}}{\partial n_0} = P_g(n_g). \quad (1.94)$$

Let us notice that beta equilibrium is not imposed anymore since  $Z$  is fixed, causing a modification of Eq. (1.92) with respect to the virial theorem with curvature term, Eq. (1.85). For a given  $n_B$ , the minimization is carried out for each tabulated  $Z$  using

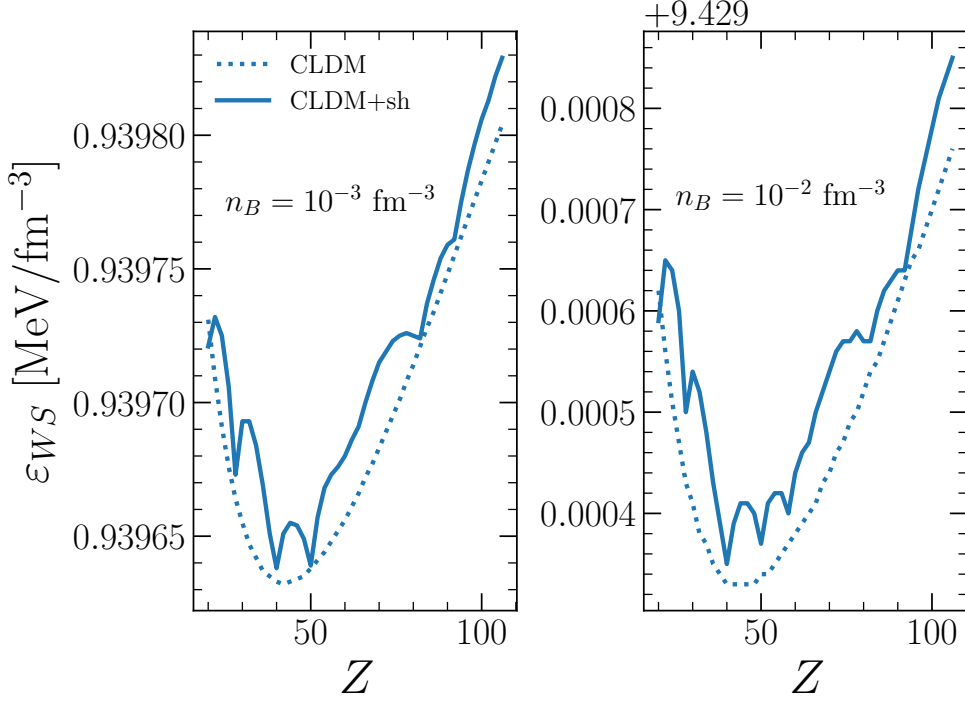


Figure 1.9: WS cell energy density as a function of number of protons  $Z$  at two different densities in the inner crust for BSk24 CLDM with (solid lines) and without (dotted lines) Strutinsky shell corrections on top of the CLD energy. **In the right panel, the constant value  $9.429 \text{ MeV/fm}^3$  is subtracted from the WS cell energy density.**

Broyden's method. The proton Strutinsky shell corrections are then added perturbatively to the CLD energy by interpolation among values of the table, which gives the shell energy per nucleon for a given  $(n_B, Z)$ . The WS energy density including shell corrections thus reads

$$\varepsilon_{WS}(n_B, Z) = \varepsilon_{WS}^{CLD}(n_B, Z) + n_B \Delta e_{sh}(n_B, Z), \quad (1.95)$$

where  $\varepsilon_{WS}^{CLD}$  is the smooth part of the WS energy density at the equilibrium, and  $\Delta e_{sh}$  is the interpolated shell energy per nucleon. The value of  $Z$  that minimizes the energy density of matter is naturally defined as the equilibrium value.

Fig. 1.9 shows the variation of the WS energy density as a function of  $Z$  at two selected values of baryon density  $n_B$ , for BSk24 CLDM. As expected, the pure CLD results, represented in dotted lines, are close to those including Strutinsky shell corrections, represented in solid lines, for closed-shell configurations, while remarkable differences exist for all other values of  $Z$ . This feature confirms the importance of a proper account of the shell structure. At  $n_B = 10^{-3} \text{ fm}^{-3}$ , a competition between  $Z = 40$  and  $Z = 50$  can be observed, with  $Z = 40$  being slightly favored. However, it should be stressed that the difference in energy density is about  $10^{-6} \text{ MeV/fm}^{-3}$ , which is expected to be lower

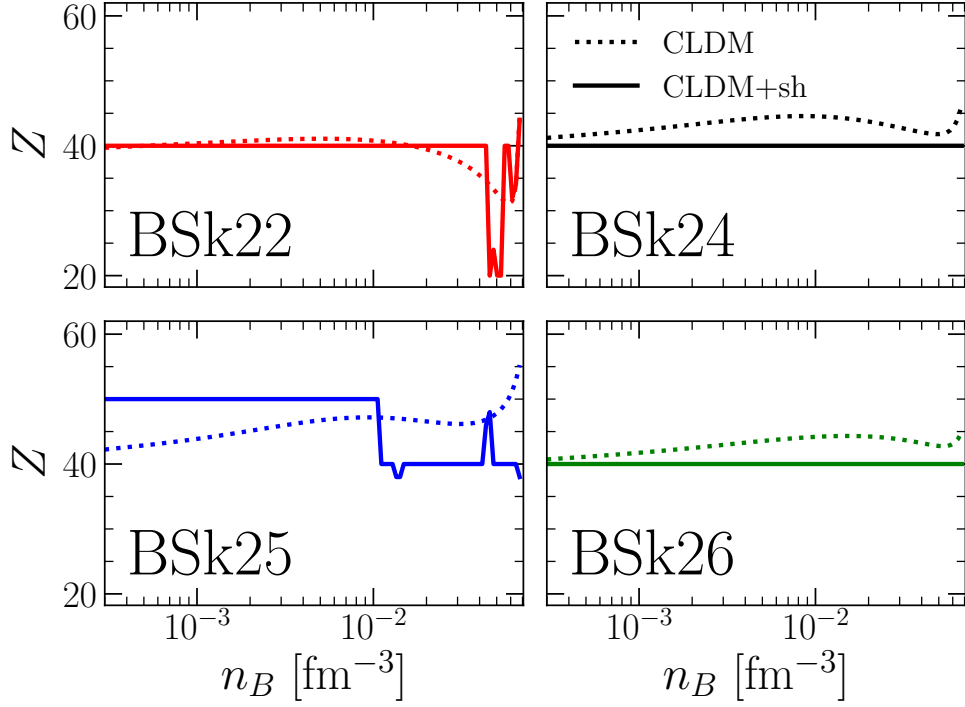


Figure 1.10: Equilibrium value of  $Z$  as a function of baryon density  $n_B$  in the inner crust for BSk22, BSk24, BSk25, and BSk26 CLDM with (solid lines) and without (dotted lines) Strutinsky shell corrections on top of the CLD energy.

than the precision reached by our modeling. It is observed that the shell closure  $Z = 40$  is also favored over  $Z = 50$  and  $Z = 58$  at  $n_B = 10^{-2} \text{ fm}^{-3}$ .

Our results for the equilibrium value of  $Z$  in the inner crust using BSk22, BSk24, BSk25, and BSk26 CLDM with (solid lines) and without (dotted lines) Strutinsky shell corrections can be seen in Fig. 1.10. It can be observed that magic numbers are recovered when shell corrections are added on top of the CLD energy. In particular, remarkable stability at  $Z = 40$  is seen for BSk24 and BSk26, as well as  $Z = 20$  and  $Z = 40$  for BSk22, in a very good agreement with Fig. 12 of [Pea+18]. A small difference only appears for the BSk25 model, reflecting the limitations of the CLD approach. For this functional, after the plateau at  $Z = 50$ , also obtained with extended Thomas-Fermi plus Strutinsky integral (ETFSI) calculations of [Pea+18] for the same functional, the equilibrium number of protons drops to  $Z = 40$  in our case, instead of increasing as in [Pea+18]. However, we stress that the authors find a second minimum at  $Z = 40$ , and the energy difference between the two minima is of the order of  $10^{-3} \text{ MeV}$  [Pea+19].

Let us recall that surface and curvature parameters  $\sigma_0$ ,  $b_s$ ,  $\sigma_{0c}$ , and  $\beta$  are so far fitted to experimental masses of nearly symmetric nuclei from the AME2016 [Hua+17] with a fixed value of the parameter  $p$  that governs the behavior of the surface tension at extreme isospin values. Therefore the shell energy is implicitly accounted for in

the surface tension by underestimating the value of  $\sigma_0$ , that is the surface tension of symmetric nuclei. This results in predicting slightly overbound nuclei with the CLD approach, but overall in a better description of spherical nuclei. However, we face an issue when adding Strutinsky shell corrections to the CLD energy, since we do not want to double-count the shell energy. To address this problem, we fit the surface and curvature parameters to the ETF results for each functional [Pea19], keeping the parameter  $p$  fixed to the value  $p = 3$  [Car+20], which gives an accurate reproduction of the CC transition points obtained in [Pea+19].

### 1.2.5 Nonspherical pasta phases

We have assumed so far that the inner crust only consists of spherical clusters. However, at high density just before the CC transition, nuclei are almost close enough to touch each other, and thus the lattice energy significantly reduces the Coulomb energy. Consequently, matter could in principle arrange itself into exotic structures, sometimes referred to as nuclear “pasta”. In particular, from the density at which the volume fraction exceeds  $1/2$ , it is expected that nuclei turn inside out, therefore forming neutron bubbles immersed in a proton-rich phase (the former cluster) [BBP71]. Furthermore, molecular dynamics simulations, in which a large number of interacting nucleons are let evolving in a very large box, predict the appearance of complex structures for  $Y_p \approx 0.1$ , which is comparable to the typical value in the bottom layers of the inner crust of neutron stars [Wat+03]. It should be stressed that these exotic phases of matter could constitute half of the crust mass [LRP93]. It is therefore straightforward to imagine that various astrophysical phenomena, such as the cooling process by neutrino emission, can be affected by their existence [WM11].

One of the virtue of the CLD approach is that different geometries for nuclear clusters can be considered in a simple way, allowing for the study of the pasta phases [RPL83; LS91; LRP93; NGL12]. Let us first recall the virial theorem with curvature term, here expressed in term of the surface, curvature, and Coulomb energy densities,

$$\varepsilon_{surf} + 2\varepsilon_{curv} = 2\varepsilon_{Coul}. \quad (1.96)$$

Following [RPL83; NGL12], the general expression for the surface energy density is

$$\varepsilon_{surf} = \frac{ud\sigma}{r}, \quad (1.97)$$

and that of the curvature energy density is

$$\varepsilon_{curv} = \frac{ud(d-1)\sigma_c}{r^2}, \quad (1.98)$$

where the surface tension  $\sigma$  and curvature tension  $\sigma_c$  are independent of the dimensionality, and therefore given by Eqs. (1.59) and (1.61), respectively. The Coulomb energy density reads

$$\varepsilon_{Coul} = 2\pi(eY_p n_0 r)^2 u \eta_{Coul,d}(u), \quad (1.99)$$

with  $\eta_{Coul,2} = \frac{1}{4} \left[ \ln\left(\frac{1}{u}\right) + u - 1 \right]$  for  $d = 2$ , and  $\eta_{Coul,d}(u) = \frac{1}{d+2} \left[ \frac{2}{d-2} \left( 1 - \frac{du^{1-2/d}}{2} \right) + u \right]$  otherwise.

In the previous expressions,  $r$  represents the radius (half-width in the case of planar geometry) of clusters or holes,  $u$  the volume fraction occupied by the cluster (hole),  $n_0$  the density of the dense phase, and  $Y_p$  its associated proton fraction. The expression of the volume fraction is different depending on whether we have clusters or holes:

$$u = \begin{cases} (n_B - n_g)/(n_0 - n_g) & \text{for clusters,} \\ (n_0 - n_B)/(n_0 - n_g) & \text{for holes.} \end{cases} \quad (1.100)$$

The parameter  $d = \{3, 2, 1\}$  is related to the dimensionality, with  $d = 3$  for spheres and bubbles,  $d = 2$  for cylinders and tubes, and  $d = 1$  for plates. In the former case, we naturally recover the expressions Eqs. (1.58), (1.60), and (1.56). Also, let us note that nuclei and bubble phases are identical as far as plates are concerned.

Replacing these quantities in Eq. (1.96), we obtain an equation for the cluster (hole) radius/half-width,

$$4\pi(eY_p n_0)^2 \eta_{Coul,d}(u) r^4 - d\sigma r - 2d(d-1)\sigma_c = 0, \quad (1.101)$$

which has to be solved numerically. Let us notice that, neglecting the curvature term, the virial theorem reduces to  $\varepsilon_{surf} = 2\varepsilon_{Coul}$ , thus the expression of the  $r$  is analytical,

$$r = \left( \frac{d\sigma}{4\pi(eY_p n_0)^2 \eta_{Coul,d}(u)} \right)^{1/3}. \quad (1.102)$$

In order to find the most stable phase at a given baryon density, we proceed as follows. The composition is first calculated at a given  $n_B$  in the spherical nuclei case ( $d = 3$ ) following 1.2.2. Then, keeping the same composition, the radius/half-width  $r$  is evaluated for the five different phases by solving numerically Eq. (1.101). Finally, since the bulk and electron contributions to the WS cell are independent of the dimensionality, the equilibrium phase is the one that minimizes  $\varepsilon_{surf} + \varepsilon_{curv} + \varepsilon_{Coul}$ . **Let us notice that this method has to be considered as an approximation since we do not consider the dimensionality dependence of the surface plus curvature energy in the minimization. This is similar to the so-called “coexisting phase method” employed in several works [Ava+08; Ava+10; Gra+17].**

In the left panel of Fig. 1.11, we represent the WS energy density as a function of the baryon density for the five phases, using our CLDM to calculate the surface and curvature energy for the BSk24 functional. It is clearly seen that the spheres dominates up to  $n_B \approx 0.05 \text{ fm}^{-3}$ , which is in good agreement with the semiclassical ETF results for this force [PCP20]. From this density, we find cylinders up to  $n_B \approx 0.065 \text{ fm}^{-3}$ , before observing a transition to plates. Then, the differences in energy density become so small that the possible transition to another phase cannot be distinguished in the figure. However, one could expect the usual sequence, that is a transition to tubes eventually followed by one to bubbles before the final transition to homogeneous matter [RPL83], here marked by the black dashed lines.

The right panel of Fig. 1.11 shows the equilibrium phase as a function of  $n_B$  for four CLDM: BSk24, SLy4, BSk22, and DD-ME $\delta$ . While model dependence is observed, it is interesting to notice that the sequence spheres  $\rightarrow$  cylinders  $\rightarrow$  plates  $\rightarrow$  tubes is





Figure 1.11: Left panel: energy density of matter as a function of baryon density for the five phases, using BSk24 CLDM. The black dashed lines mark the CC transition point. Right panel: equilibrium phase as a function of baryon density for four selected CLDM: BSk24, SLy4, BSk22, and DD-ME $\delta$ . For these calculations, the surface parameter  $p$  is fixed to the value  $p = 3$ .

recovered for every model. Another shared feature is that the transition to homogeneous matter is found before the appearance of bubbles. Let us remark that pasta phases are observed for SLy4 here, while previous CLD and TF calculations for this functional have predicted that the sphere is the most favorable shape at all densities up to the CC transition point [DH00; Viñ+17]. This contradiction can be understood by the fact that we are dealing with very tiny differences in energy density (all the shapes are very close), in a region where the matter is extremely neutron rich, thus where the nuclear surface tension is poorly constrained. In particular, we observed that varying the value of the surface parameter  $p$  can affect the sequence of pasta phases.

### 1.2.6 Crust-core transition from the crust side

It is well known that a phase transition from a solid crust to a liquid core takes place at some  $\approx 1$  km from the surface of the star, corresponding to subsaturation densities. A precise estimation of this transition point is necessary in order to evaluate the crust mass, thickness and moment of inertia, and so to understand phenomena involving NS,

such as glitches, that are irregularities in their rotational motion [Esp+11]. In that sense, a Bayesian analysis of the CC transition has been performed [CGM19a] and is presented in details in Chapter 2.

The simplest approach for computing the phase transition is from the core side, in which the transition is defined as the density point where homogeneous nuclear matter (NM) at beta equilibrium becomes unstable with respect to density fluctuations. Two versions of this technique can be distinguished: the so-called thermodynamical [Gon+17] and dynamical [PRL95; Ant+19] methods.

The former method consists in evaluating the thermodynamical spinodal, that is the instability point of NM with respect to the liquid-gas (LG) phase transition. The only necessary nuclear physics input for this method is the energy functional of homogeneous NM, which can be well constrained through nuclear experiments and/or ab initio calculations in the vicinity of the nuclear saturation density  $n_{sat}$ . While it appears as a substantial advantage, it is known that the dynamics of the CC transition is very different from the LG one [DCG07; Duc+07; DMC08], thus applying this method results in an overestimation of the CC transition density and pressure. Indeed, the CC transition is expected to take place at approximately  $\approx n_{sat}/2$ , where the energy density of clusterized matter overcomes that of uniform NM [BBP71]. Since Coulomb, surface, and curvature terms, that are crucial for determining the equilibrium composition in the inner crust, vanish in homogeneous NM, they do not contribute to the determination of the LG thermodynamical spinodal.

In the dynamical method, finite-size density fluctuations are added, thus it is expected to give a better estimation of the CC transition and pressure. It should be stressed that the isovector gradient terms, which are needed in addition to the EoS with respect to the thermodynamical method, play a nonnegligible role for the determination of the dynamical spinodal, the location of which also depends on the many-body formalism adopted [DMC08].

While the dynamical method gives a better estimation of the CC transition point in comparison with the thermodynamical one, one should remark that the spinodal decomposition scenario is not compatible with the equilibrated crust in which matter is not uniform but composed of clusters immersed in a sea of electrons and neutrons. For this reason, the CC transition point is here computed from the crust side, by comparing the energy density of the two competing phases in beta equilibrium. The equation for the CC transition density  $n_t$  is thus

$$\varepsilon_{WS}^{crust}(n_t) = \varepsilon_{npe}(n_t), \quad (1.103)$$

where  $\varepsilon_{WS}^{crust}$  is the equilibrium energy density in the WS cell in the inner crust, and  $\varepsilon_{npe}$  is the equilibrium energy density of matter in the outer core, which consists of homogeneous *npe* matter, studied in Section 1.3. The CC transition point is computed in the following for several nuclear models, using the metamodeling technique extended to finite nuclei in the CLD approximation, Eq. (1.63), to calculate the WS cell energy density in the inner crust, and keeping the same empirical matter for the description of uniform *npe* matter. One should stress the importance of having a unified EoS for the crust and the core [DH01] to estimate correctly the CC transition point. Indeed, it was shown that matching a crust and core EoS based on different nuclear models induce



Figure 1.12: Crust-core transition density as a function of the surface parameter  $p$  with truncation order  $N = 2$ ,  $N = 3$ , and  $N = 4$ , for SLy4 CLDM. The black horizontal bars represent different evaluations of  $n_t$  for the SLy4 functional found in the literature:  $n_t = 0.072 \text{ fm}^{-3}$  [Viñ+17] (dashed line) and  $n_t = 0.077 \text{ fm}^{-3}$  [DH00] (solid line) in the CLDM approximation,  $n_t = 0.076 \text{ fm}^{-3}$  [Viñ+17] (dotted line) in the TF approximation, and  $n_t = 0.080 \text{ fm}^{-3}$  [Viñ+17] (dashdotted line) with the dynamical method.

an uncertainty on the determination of the CC transition point, consequently inducing an uncertainty in the crustal observables, which can be as large as 30% for the crust thickness [For+16].

Let us mention that pasta phases are not considered here, since we have checked that their presence does not affect the estimation of the CC transition density and pressure.

In Fig. 1.12, we show the variation with surface parameter  $p$  of  $n_t$  for SLy4 CLDM. For each value of  $p$ , the surface and curvature parameters are fitted to experimental masses from the AME2016. The CC transition density is estimated for different values of the truncation order  $N$  in the density development, Eq. (1.50). A convergence feature is observed from  $N = 2$  to  $N = 4$ , almost achieved at  $N = 3$ . We can see that truncating at  $N = 2$  leads to a substantial underestimation of about  $\approx 20\%$  of the transition density. This observation highlights the importance of keeping high-order parameters beyond  $(Q_{sat}, Q_{sym})$  ( $N = 3$ ), and  $(Z_{sat}, Z_{sym})$  ( $N = 4$ ) in the Taylor expansion.

A strong correlation is seen between the isovector surface parameter  $p$  and  $n_t$ : the higher the value of  $p$  the higher the transition density. This can be understood from the fact that a low value of  $p$  leads to a higher surface tension at extreme isospin values, thus it makes the cluster less bound and consequently the uniform  $npe$  matter becomes the

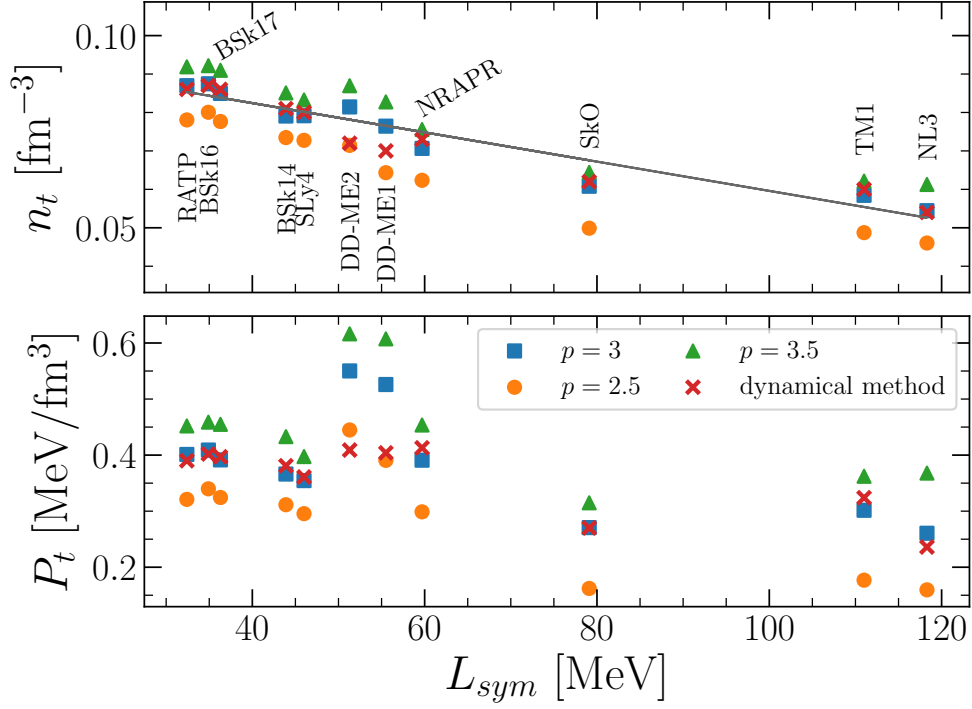


Figure 1.13: Crust-core transition density  $n_t$  (upper panel) and pressure  $P_t$  (lower panel) as a function of  $L_{sym}$  for several interactions. The crosses are the transition points calculated in [Duc+11] using the dynamical spinodal. The circles, squares, and triangles correspond respectively to our estimation of the transition points with  $p = 2.5$ ,  $p = 3$ , and  $p = 3.5$ . The gray line is the linear regression curve fitted to the  $p = 3$  results.

favorable phase at lower density.

Different estimations of the CC transition density for the SLy4 functional found in the literature are reported in Fig. 1.12 (black lines), with  $n_t$  ranging from  $0.072 \text{ fm}^{-3}$  [Viñ+17], in the CLD approximation, to  $0.089 \text{ fm}^{-3}$  [Duc+11] (not represented in the figure), with the thermodynamical method. It is interesting to observe that the **two different CLDM calculations of the literature based on SLy4 functional, reported in the figure**, give a different estimation of  $n_t$ , with  $0.072 \text{ fm}^{-3}$  in [Viñ+17], and  $0.077 \text{ fm}^{-3}$  in [DH00]. This stresses once again the importance of the treatment of the isovector surface tension for the determination of the CC transition point.

The effect of the varying the low-density parameter  $b$  on  $n_t$  is also evaluated in [CGM19a]. It is found that while it does not significantly affect the transition density in comparison with  $N$  and  $p$ , it should be kept as an additional EoS parameter in any statistical analysis in order to make a quantitative prediction of the CC transition point. **The introduction of the parameter  $b$  in the parameter space is, in fact, an effective manner to account for the effect of orders  $N > 4$ , which is not important.**

Fig. 1.13 shows our estimations of the transition density  $n_t$  and pressure  $P_t$  as a func-

tion of the slope of the symmetry energy  $L_{sym}$  for a set of relativistic and nonrelativistic functionals, in comparison with the dynamical spinodal calculations of [Duc+11]. It is seen that the value  $p = 3$  is in good agreement with the dynamical results for most of the models. For DD-ME2, the lower value needed for  $p$  is supported by the TF calculations in [GPA12].

As in [Viñ+17], an anticorrelation is observed between  $L_{sym}$  and  $n_t$ . Using a linear regression to model the relationship between  $L_{sym}$  and our estimation of  $n_t$  with  $p = 3$ , we find  $n_t = -3.807 \times 10^{-4} L_{sym} + 0.098 \text{ fm}^{-3}$ , for a root mean square error of  $3.094 \times 10^{-3} \text{ fm}^{-3}$ . This linear equation for  $n_t$  is in very good agreement with Eq. (17) of [Duc+11]. This downwards tendency with increasing values of  $L_{sym}$  is not clearly observed for the transition pressure. **We will see in Chapter 2 that the relation between those two quantities is less explicit because  $P_t$  rather depends on  $K_{sym}$ , which is logical because the pressure requires the calculation of one more derivative in comparison to the energy density that defines  $n_t$ .**

### 1.3 Matter of the core

In the very bottom layers of the inner crust, clusters are so large and asymmetric that the CLD energy per nucleon, Eq. (1.63), becomes equal to that of infinite nuclear matter, marking the CC interface. While the composition and structure of the core close to the saturation density are well known, various scenarios have been proposed in the literature [Oer+17] concerning the relevant degrees of freedom above  $2 - 3n_{sat}$ .

This section deals with the matter of the core. In 1.3.1 we derive the system of variational equations for the ground state of the outer core. The model dependence of the composition and EoS is then explored. In particular, we focus on apparent correlations with the symmetry energy. The matter inside the inner core is finally discussed in 1.3.2.

#### 1.3.1 Outer core: homogeneous $npe\mu$ matter

From the CC transition point, previously discussed in 1.2.6, up to  $n_B \approx 2n_{sat}$ , the matter of the core consists of a uniform plasma of neutrons, protons, electrons, and eventually muons, hereafter referred to as  $npe\mu$  matter (or  $npe$  if muons are not present). This region of the star corresponds to the outer core. The energy density of  $npe\mu$  matter reads

$$\varepsilon_{npe\mu}(n_B, \delta, n_e, n_\mu) = \varepsilon_{HM}(n_B, \delta) + \varepsilon_e(n_e) + \varepsilon_\mu(n_\mu) + n_B \frac{1 - \delta}{2} \Delta m_{pn} c^2 + n_B m_n c^2, \quad (1.104)$$

with  $\varepsilon_{HM} = n\varepsilon_{HM}$  the energy density of nuclear matter given by the metamodel, Eq. (1.37). The expression of energy density of the relativistic muon gas  $\varepsilon_\mu(n_\mu)$  is the same as that of the relativistic electron gas, and it is derived in Appendix A.

##### 1.3.1.1 Variational equations

Following 1.2.2, the system of differential equations determining the equilibrium composition and EoS in the outer core is determined by minimizing the energy density of

matter, Eq. (1.104), at constant baryon density under the condition of overall charge neutrality, which now reads

$$n_B \frac{1 - \delta}{2} = n_e + n_\mu. \quad (1.105)$$

since muons may exist in this region of the star. Muons are present in the outer core if the electron chemical potential exceeds the muon rest-mass energy, that is  $\mu_e \gtrsim m_\mu c^2$ . As far as variational variables are concerned, a reasonable choice appears to be the asymmetry  $\delta$  and the muon density  $n_\mu$ . Let us first minimize the energy density of matter with respect to  $\delta$ ,

$$\begin{aligned} \left. \frac{\partial \varepsilon_{npe\mu}}{\partial \delta} \right|_{n_\mu} &= 0 \\ 2 \frac{\partial e_{HM}}{\partial \delta} - \Delta m_{pn} c^2 - \mu_e &= 0. \end{aligned} \quad (1.106)$$

We can easily identify this equation with the well-known beta equilibrium condition. Indeed, using the definition of the neutron and proton chemical potentials, Eq. (1.53), we obtain

$$\mu_n = \mu_p + \mu_e. \quad (1.107)$$

We now turn to minimization with respect to the muon density,

$$\begin{aligned} \left. \frac{\partial \varepsilon_{npe\mu}}{\partial n_\mu} \right|_\delta &= 0 \\ \mu_\mu - \mu_e &= 0, \end{aligned} \quad (1.108)$$

where we have introduced the muon chemical potential  $\mu_\mu = \partial \varepsilon_\mu / \partial n_\mu$ . This equation corresponds to the chemical equilibrium between muons and electrons, and has to be satisfied if muons are present. These two equations, Eq. (1.107) and (1.108), express the equilibrium with respect to the weak interaction processes. **Let us remark that each time we mention electrons and muons, we are always referring to their net number, that is particles minus antiparticles.**

From the numerical point of view, we use Broyden's method to solve the beta equilibrium equation, Eq. (1.107), starting from the transition density  $n_t$ . The initial guess for the asymmetry  $\delta_{guess}$  is taken to be the value of the global asymmetry inside the WS cell at the CC transition point, which is approximately  $\approx 0.9$ . At each new step of baryon density, the value of  $\delta_{guess}$  is updated with the last equilibrium solution, and we check if  $\mu_e \gtrsim m_\mu c^2$ . In the case of  $npe\mu$  being energetically favorable, Eq. (1.108) is added as a supplementary equation to be solved, setting the initial guess for the muon density to  $n_{\mu, guess} = 10^{-5} \text{ fm}^{-3}$ . The calculation is stopped when the baryon density reach  $\approx 2n_{sat}$ , from which the composition of matter becomes less certain.

### 1.3.1.2 Equilibrium composition and equation of state

In Fig. 1.14, the proton fraction  $Y_p = (1 - \delta)/2$  and muon fraction  $Y_\mu = n_\mu/n_B$  are plotted as a function of the baryon density for the four nuclear models BSk24, SLy4, BSk22, and DD-ME $\delta$ , the empirical parameters of which are listed in Table 1.3. The variation

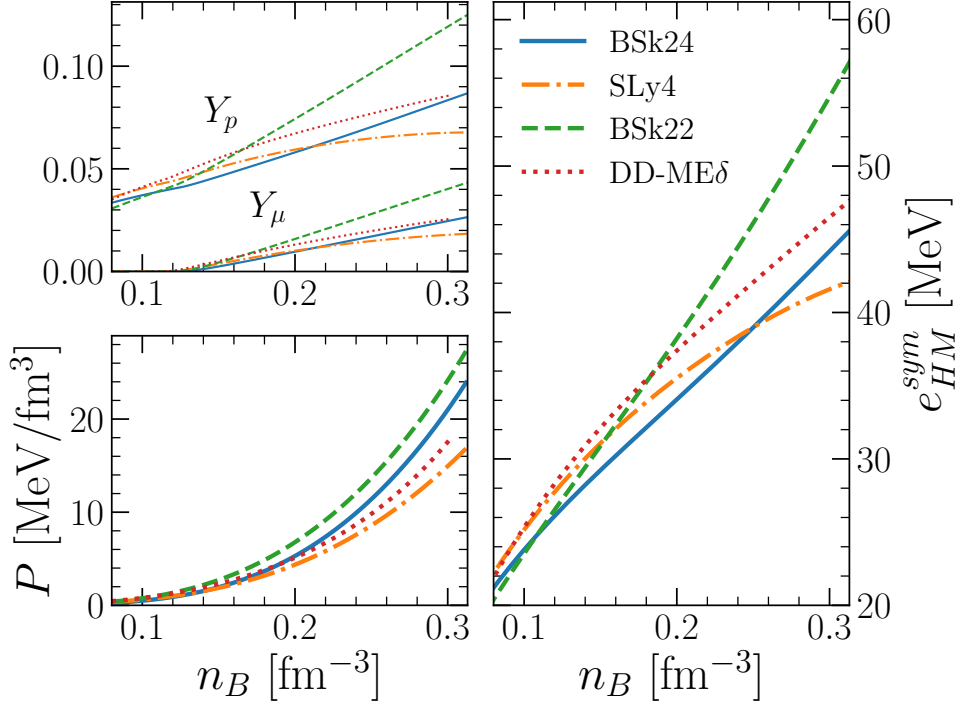


Figure 1.14: Left: Proton fraction  $Y_p$  and muon fraction  $Y_\mu$  (upper panel), and pressure  $P$  (lower panel) as a function of the baryon density  $n_B$  in the outer core for four selected models: BSk24, SLy4, BSk22, and DD-ME $\delta$ . Right: Variation with baryon density of the symmetry energy  $e_{HM}^{sym}$ , Eq. (1.52), for the four models.

$n_B$  of the electron fraction  $Y_e$  can be deduced from the charge neutrality condition, Eq. (1.105). Surprisingly, it is observed that the proton fraction rises with  $n_B$  in the outer core, while it monotonously decreases in the crust. A strong correlation between the proton fraction and the symmetry energy, represented in the right panel for the four models, is seen. One can relate the two quantities using the beta equilibrium condition. Indeed, we can easily go through Eq. (1.107) assuming a quadratic expansion for the NM energy per nucleon,

$$e_{HM}(n_B, \delta) \simeq e_{HM}(n_B, \delta = 0) + \delta^2 e_{HM}^{sym}(n_B). \quad (1.109)$$

In that case, the beta equilibrium condition reads

$$4\delta e_{HM}^{sym}(n_B) - \Delta m_{pn} c^2 = \mu_e(n_e), \quad (1.110)$$

and we can thus express the proton equilibrium proton fraction as a function of the symmetry energy,

$$Y_p = \frac{1}{2} - \frac{\mu_e(n_e) + \Delta m_{pn} c^2}{8e_{HM}^{sym}(n_B)}. \quad (1.111)$$

From this equation it is clear that the higher the symmetry energy, the higher the proton fraction, which is the trend observed in the figure.

The density from which muons are present is also shown in the upper left panel. It appears that we are dealing with *npe* matter up to  $n_B \approx 0.12 \text{ fm}^{-3}$ , and it is seen that the threshold density for the appearance of muons, when  $\mu_e \gtrsim m_\mu c^2$ , is not very sensitive to the nuclear model.

Fig. 1.14 also shows the pressure as a function of the density (lower left panel). In the outer core, we can write the pressure as

$$P(n_B) = P_{HM}(n_B, \delta) + P_e(n_e) + P_\mu(n_\mu), \quad (1.112)$$

where  $\delta$ ,  $n_e$ , and  $n_\mu$  are the equilibrium values at  $n_B$ . The expression of the nuclear contribution  $P_{HM}$  is given by Eq. (1.54). The electron and muon contributions to the total pressure, respectively  $P_e$  and  $P_\mu$ , share the same expression, Eq. (1.11), which is that of a relativistic Fermi gas. A strong model dependence is observed for the pressure at suprasaturation densities, reflecting the large uncertainties associated to the isovector parameters beyond  $E_{sym}$ , as reported in the last column of Table 1.3.

### 1.3.2 Inner core

From the crust-core interface up to  $n_B \approx 2n_{sat}$ , it is commonly accepted that matter is exclusively composed of neutrons, protons, electrons, and muons. At higher densities however, the composition of matter becomes uncertain and different scenarios have been considered in the literature for the structure and composition of the inner core [Oer+17]. The presence of hyperons at suprasaturation densities has been considered in many studies, and its effect on the EoS and subsequently on NS observables has been evaluated [Bed+12; For+15]. In particular, the presence of hyperons softens the high-density EoS and thus lowers the maximum mass of neutron stars. As a consequence, many models fail to satisfy the NS mass constraint of  $\approx 2M_\odot$  [Dem+10; Ant+13], leading to the so-called “hyperon puzzle” [ZH13]. In fact, the measurement of  $\approx 2M_\odot$  NS also rules out most of the models that include a phase transition to quark matter or a boson condensate in the inner core, because the EoS is softened by the introduction of any additional degree of freedom without an interaction. Let us however mention a recently proposed EoS with a quark-hadron crossover at high density that allows a maximum mass of  $2.35M_\odot$  [Bay+19] and agrees with the constraints on NS properties inferred from the GW170817 event [Abb+18; Abb+19].

We propose here to extrapolate the *npe $\mu$*  matter at higher densities, in the inner core. The different studies that consider hyperons and satisfy the maximum mass constraint show that the hyperon-nucleon and hyperon-hyperon are such that, if hyperons exist, they are not abundant and can be neglected as a first approximation, at least in the EoS energetics (see [Oer+17] and references therein). Since the metamodeling does not make any assumption regarding the degrees of freedom, but that the functional is based on a Taylor expansion, we can assume that the technique can be safely extended at high density in the aim of calculating static properties, such as the EoS or the mass-radius relation. The only thing that could invalidate it would be the presence of a large first-order phase transition, that is a large discontinuity in the EoS. Thus,



the only hypothesis that is made here is considering that there is no such transition. Actually, even a transition to quark matter could be smooth. Indeed, several works consider pasta phases in the hadron-quark phase transition, with the effect of washing out discontinuities [MTC09; YMT09; YMT11; Yas+14]. Finally, this is equivalent to a null hypothesis: since the metamodeling allows to control uncertainties, a disagreement with astrophysical observations would be the sign of new physics, indicating either the existence of a first-order phase transition or an alternative theory of gravity. Let us recall that the metamodeling technique gives an accurate reproduction of existing nuclear models up to  $\approx 2 - 3n_{sat}$  (metamodeling ELFc in [MHG18a]), as we have verified in Fig. 1.4. A solution for achieving faster convergence at high density, the metamodeling ELFd, is proposed in [MHG18a]. It is based on a reevaluation of the high-order empirical parameters  $Q_{sat}$ ,  $Q_{sym}$ , and  $Z_{sat}$ ,  $Z_{sym}$ , by imposing a high-density point at which the energy per nucleon and the pressure are known for a given functional. This is encouraged by the fact that the impact of these parameters is very small in the vicinity of the saturation density. We will use this technique when we will want to reproduce existing hadronic models.

## 1.4 Unified metamodeling of the equation of state

As previously discussed, matching a crust and core based on different nuclear models can lead to large uncertainties associated with the CC transition density and pressure, which are the quantities for the determination of the crust observables. In particular, it was shown in [For+16] that it can lead to an error as large as  $\approx 30\%$  for the crust thickness and  $\approx 4\%$  for the radius. The main justification for working with nonunified EoS is that the core is responsible of most of the NS mass, thus one could simply match any crust EoS, for instance that of BPS in the outer crust [BPS71] plus that of BBP in the inner crust [BBP71], with its core EoS based on a different nuclear interaction. This is indeed sufficient if one is interested in observables such as the maximum mass. However, it obviously leaves a large freedom in the matching procedure. In addition, this amounts to consider that the crust segment is not model dependent, while we have clearly seen in 1.2.3.3 that the crust EoS is sensitive to the nuclear model, and is in particular correlated with the symmetry energy. It is therefore preferable to describe the crust and core in a unified manner.

Several unified EoS for cold non-accreting NS have been proposed in the literature. For instance, the well-known DH EoS [DH01] is based on the Skyrme SLy4 effective interaction [Cha+98]. In the inner crust, the authors use a version of the CLDM to calculate the energy of finite nuclei, the bulk term of which is given by the SLy4 functional in the infinite NM limit, which is also used to describe the outside neutron gas, and at high density to calculate the  $npe\mu$  matter energetics in the core. Let us notice that the EoS of the outer-crust ground state is taken from [HP94] in this work.

More recently, a series of unified EoS based on modern BSk functionals [GCP13] has been proposed [Pea+18]. In this work, the ground state of the outer crust is determined by application of the microscopic HFB method [Sam+02], and the ground state of the inner crust is calculated within the ETFSI approximation [Ons+08]. Also, a set of unified EoS were built within a relativistic mean-field approach (RMF) in [For+16].

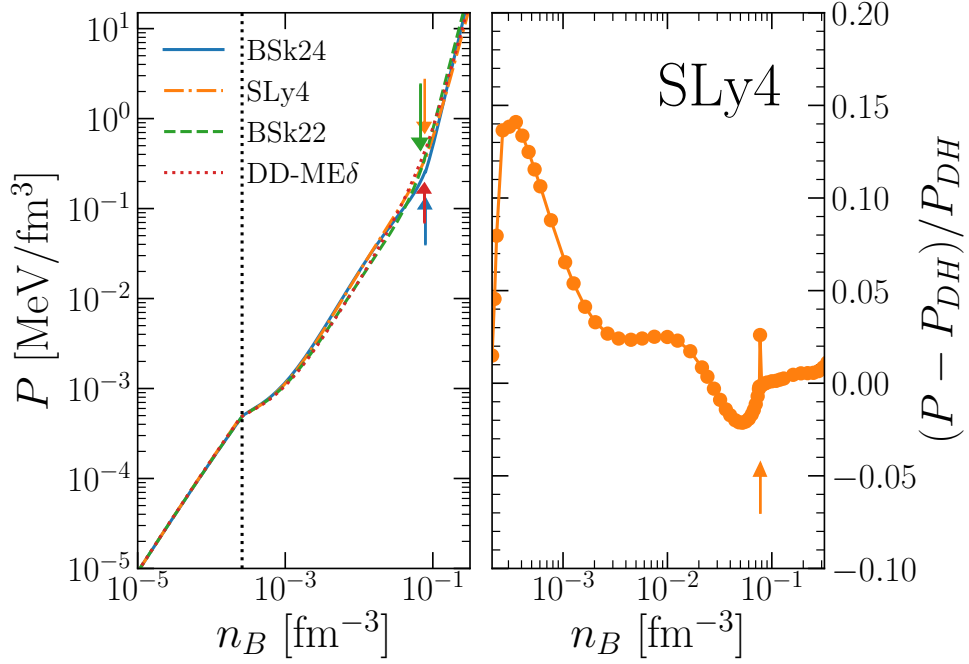


Figure 1.15: Left: Pressure as a function of baryon density in the different regions of a NS for BSk24, SLy4, BSk22, and DD-ME $\delta$ . Crust and core are described in a unified manner, with the metamodeling technique. The black dotted line marks the onset of the inner crust, and arrows indicate the CC transition density for the different models. Right: Relative difference  $(P - P_{DH})/P_{DH}$  with DH unified EoS for SLy4 [DH01] as a function of  $n_B$ . The arrow indicates our estimation of the CC transition density for SLy4.

One obvious reason explaining why crust and core matter are not consistently treated using the same microscopic interaction is that it is less trivial to evaluate the ground state of the inner crust in comparison to the ground state of matter in the core. Indeed, one can evaluate the inner-crust ground state within different many-body approaches, either using the CLDM, the semiclassical ETFSI method, or full HF calculations. While the two last methods are expected to be more precise with respect to the CLD approximation as far as the composition is concerned, they can become computationally heavy if one wants to compute the EoS for many nuclear models. **In addition, these more microscopic methods can only be extended to finite temperature in the SNA, whereas the CLD approximation allows distributions of clusters, that exist at  $T > 0$ , to be considered.**

In view to provide a unified and thermodynamically consistent treatment of the crust and core of cold non-accreting NS, we propose to compute the EoS within the metamodeling approach that we have presented throughout this chapter. For a given nuclear model, in other words a given set of empirical parameters  $\mathbf{P}_\alpha$ , we proceed as follows. The EoS in the outer crust is evaluated by application of the BPS method presented in 1.1.2, using

the present day knowledge on experimental masses [Hua+17] supplemented by state-of-the-art microscopic HFB theoretical mass tables [GCP13] up to the neutron-drip point. The EoS in the inner crust is then calculated by solving the system of four differential equations, Eqs. (1.85), (1.86), (1.87), and (1.88). The metamodeling technique, discussed in 1.2.1.1, is used to calculate the neutron gas energy. The concept of metamodeling of homogeneous NM is extended to finite nuclei in the CLD approximation in order to evaluate the cluster energy, the bulk term of which is calculated with the same empirical parameters as the neutron gas. Finally, from the CC transition point, discussed in 1.2.6, the core EoS is computed by solving the two chemical equilibrium equations Eqs. (1.107) and (1.108). Again, the empirical parameters are the same as those used in the inner crust. Within such an approach, one can compute the unified EoS for any nuclear model. In addition, this method is not expensive from the computational point of view, which will allow us to make comprehensive statistical evaluations of the model uncertainties in Chapter 2.

The left panel of Fig. 1.15 shows the resulting unified EoS for the four nuclear models considered in this chapter: BSk24, SLy4, BSk22, and DD-ME $\delta$ . The neutron-drip point is represented by the vertical dotted line, and arrows indicate the CC transition density for each model.

The relative difference between the DH EoS [DH01] and our EoS for SLy4 is represented as a function of the baryon density in the right panel of Fig. 1.15. It is seen that the overall difference is smaller than 10% in the inner crust. The slight disagreement at low density comes from the fact that the onset of the inner crust is marked at  $n_{ND} = 2.0905 \times 10^{-4} \text{ fm}^{-3}$  in the DH EoS (calculated in [HP94]), which is low in comparison with our estimation of the neutron drip density,  $n_{ND} = 2.503 \times 10^{-4} \text{ fm}^{-3}$ . Let us notice that the parametrization of the surface tension in both studies is different, therefore it might explain the difference between the two EoS at higher densities in the inner crust, in addition to the slight error introduced by the metamodel. In the core, we can observe that the relative difference is smaller than 1%.

## 1.5 Conclusion

In this chapter, we have first evaluated the ground state of the outer crust, which is obtained by the application of the BPS method, using recent experimental masses [Hua+17; Wel+17] supplemented by state-of-the-art microscopic HFB theoretical mass tables. We have shown that the outer-crust composition and so the EoS are entirely determined by the present day knowledge on experimental masses up to  $n_B \approx 3 \times 10^{-5} \text{ fm}^{-3}$ . At higher densities, we have observed the persistence of  $N = 82$  for each of the four mass models considered.

We have proposed a version of the CLDM based on the metamodeling technique [MHG18a; MHG18b]. We have seen that the metamodel offers the possibility to reproduce any functional of nuclear matter very precisely, **with a unique functional form**. The parametrization of surface tension was suggested from microscopic calculations in the free neutron regime [RPL83], and surface and curvature parameters are fitted to experimental masses. We have used the CLDM to calculate the ground state of the inner crust, which is obtained by minimizing the energy density of matter at constant baryon

density under the condition of overall charge neutrality. We have therefore derived a system of four coupled differential equations, corresponding to chemical and mechanical equilibrium conditions, that we have solved numerically using Broyden's method. The ground-state composition of the inner crust for BSk24 empirical parameters was presented, and a very good agreement with more microscopic approaches was observed concerning the value of  $Z \approx 40$ . We have also seen that the proton fraction is continuously decreasing, as in the outer crust. The inner crust EoS was computed and a positive correlation with the empirical parameter  $E_{sym}$  was revealed. We have shown that magic numbers, that vanish within the CLD approach, can be recovered by adding perturbatively Strutinsky shell corrections on top of the CLD energy. In this way, a very good agreement with ETFSI calculations was observed for recent BSk functionals [Pea+18]. With the CLDM, we have explored the possible presence of nonspherical pasta phases in the bottom layers of the inner crust, finding the sequence spheres  $\rightarrow$  cylinders  $\rightarrow$  plates  $\rightarrow$  tubes for each of the four model considered: BSk24, SLy4, BSk22, and DD-ME $\delta$ . We have shown that the transition point to homogeneous matter is very sensitive to the surface tension at extreme values of isospin, the behavior of which is governed by the isovector surface parameter  $p$ . Unfortunately, this parameter cannot be accessed from empirical nuclear physics data, which are limited to values around  $I \lesssim 0.3$ . We have seen that the CC transition density and pressure results of the literature for the dynamical spinodal [Duc+11] are globally nicely reproduced by our calculation from the crust with  $p \approx 3$ . We have also confirmed the correlation of the transition density  $n_t$  with the empirical parameter  $L_{sym}$  already observed in previous works.

We have derived the equilibrium equations characterizing the ground state of matter in the core, which consists of  $npe\mu$  matter up to  $n_B \approx 2n_{sat}$ . The strong correlation between the symmetry energy and the proton fraction was explained. We have found that muons appear in the vicinity of  $n_B \approx 0.12 \text{ fm}^{-3}$  for each model. As in previous works [WFF88; DH01], we have extrapolated the  $npe\mu$  model to higher densities, since the hyperon-hyperon and hyperon-nucleon interactions remain currently poorly constrained.

Finally, we have stressed that building unified EoS is essential to properly estimate crustal observables. In that sense we have proposed a metamodeling of the EoS of cold non-accreting NS, where the crust and core are treated in a uniform manner, that is with the same empirical parameters. Using the SLy4 empirical parameters, we have shown that the relative difference with the DH EoS is smaller than 10% in the inner crust and 1% in the core.

In Chapter 2, we will exploit the second advantage of the metamodeling technique, namely the fact that no artificial correlations are introduced a priori among the empirical parameters, to carry out complete statistical analyses that will allow us to settle the model dependence of the results.

## Chapter 2

# Bayesian inference of neutron star observables

### 2.1 From the equation of state to neutron star observables

The stellar EoS, the determination of which was studied thoroughly in Chapter 1, is the fundamental ingredient to build NS models. It enters into the basic equation for calculating macroscopic observables relevant to NS, that is so-called Tolman-Oppenheimer-Volkoff (TOV) equation, which describes the hydrostatic equilibrium for a spherically symmetric star in the framework of general relativity [Tol39; OV39].

This section deals with the estimation of NS observables for several popular EoS calculated within the metamodeling technique introduced in Chapter 1. In 2.1.1, we first solve the hydrostatic equilibrium equations so as to determine the mass-radius relation. The thickness and mass of the crust are also evaluated. The calculation of the NS moment of inertia in the slow rotation approximation is performed in 2.1.2. In particular, we show how the fraction of crust moment of inertia is connected to the observation of pulsar glitches. In 2.1.3, we finally compute the tidal deformability, which describes how much the NS is deformed by tidal forces, arising, for instance, when two NS orbit each other.

#### 2.1.1 Masses and radii

From the observational point of view, it is easier for astronomers to measure the mass of a NS entering into a binary system. There are several types of binaries: X-ray binaries, double NS binaries, radio pulsar–white dwarf binaries, and radio pulsar–nondegenerate star binaries. Depending on the type of binary, different techniques are used to infer the NS mass [HPY07]. For example, in a pulsar binary system, various phenomena, such as the relativistic Shapiro delay [Sha64], can be measured via pulsar timing, which consists of the regular monitoring of the rotation of a pulsar over long periods (years to decades). The relativistic Shapiro delay is a phenomenon from which precise masses for both a millisecond pulsar and its companion can be inferred [Dem+10; Cro+20]. Let us notice however that it is only observed in a small subset of high-precision, highly inclined binary pulsar systems.

Measuring NS radii with high precision is a more challenging task, and it constitutes one of the goals of the ambitious NICER mission [GAO12]. This observable can be extracted from the analysis of thermal emission from neutron star surfaces, or X-ray emission from accreting neutron stars in binaries.

In the following, we turn to the theoretical prediction of NS masses and radii. The hydrostatic equilibrium equations in general relativity are first solved in order to represent the mass-radius relation for several popular EoS, which are then confronted to NS mass measurements. The estimation of the crustal thickness and mass is discussed afterwards.

### 2.1.1.1 Mass-radius relation

Theoretically, the NS masses and radii correspond to the solution of the hydrostatic equilibrium equations, which are, in general relativity for spherical and nonrotating stars [Tol39; OV39],

$$\frac{dP}{dr} = -\frac{G\rho_B m}{r^2} \left(1 + \frac{P}{\rho_B c^2}\right) \left(1 + \frac{4\pi P r^3}{mc^2}\right) \left(1 - \frac{2Gm}{rc^2}\right)^{-1}, \quad (2.1)$$

$$\frac{dm}{dr} = 4\pi r^2 \rho_B, \quad (2.2)$$

$$\frac{d\Phi}{dr} = -\frac{1}{\rho_B c^2} \frac{dP}{dr} \left(1 + \frac{P}{\rho_B c^2}\right)^{-1}, \quad (2.3)$$

where  $G$  is the gravitational constant,  $P$  the pressure,  $\rho_B c^2$  the energy density, and  $m(r)$  the is the gravitational mass inside the sphere of radius  $r$ , defined within the Schwarzschild metric  $ds^2 = c^2 dt^2 e^{2\Phi} - e^{2\lambda} dr^2 - r^2(d\theta^2 + \sin^2\theta d\phi^2)$ . The function  $\Phi(r)$  corresponds to the gravitational potential, and  $\lambda(r)$  is related to the enclosed mass  $m(r)$  through

$$e^{-\lambda} = \sqrt{1 - \frac{2Gm}{rc^2}}. \quad (2.4)$$

Eq. (2.1) is the TOV equation of hydrostatic equilibrium. The integration of Eq. (2.2) from  $r = 0$  to the boundary  $r = R$ ,  $R$  being the NS radius, gives the total gravitational mass  $M = m(R)$  of the star. Eq. (2.3) is a relativistic equation for the metric function  $\Phi(r)$ . Here, we focus on solving Eqs. (2.1) and (2.2), which gives the profiles  $P(r)$ ,  $\rho_B(r)$ , and  $m(r)$  for a given EoS,  $P(\rho_B)$ , the determination of which was discussed in detail in Chapter 1.

In view of solving the hydrostatic equilibrium equations for a given tabulated EoS, one first have to choose an arbitrary value for the central density  $\rho_{B,c}$  and interpolate the central pressure  $P_c = P(\rho_{B,c})$ , corresponding to  $r = 0$ , with the associated boundary condition  $m(r = 0) = 0$ . Then, using a Runge-Kutta method, one integrate up to  $r = R$ , defined as  $P(r = R) = 0$  (our numerical condition is  $P < 5 \times 10^{-4}$  MeV/fm<sup>3</sup>, which is a sufficiently low value of the pressure). At each new step of pressure, one has to interpolate the mass density  $\rho_B$  in the table. In the end, for a given value of central density  $\rho_{B,c}$ , one obtains the profiles  $P(r)$ ,  $\rho_B(r)$ , and  $m(r)$ , and therefore a value for the NS mass  $M$  and radius  $R$ . The typical values of these observables are  $M = 1.4M_\odot$

Parameter	Unit	$N$	BSk14	PKDD	TM1
$E_{sat}$	MeV	0	-15.85	-16.27	-16.26
$n_{sat}$	$\text{fm}^{-3}$	1	0.1586	0.1495	0.1450
$K_{sat}$	MeV	2	239	261	281
$Q_{sat}$	MeV	3	-359	-119	-285
$Z_{sat}$	MeV	4	1435	4213	2014
$E_{sym}$	MeV	0	30.00	31.19	36.94
$L_{sym}$	MeV	1	43.9	79.5	111.0
$K_{sym}$	MeV	2	-152	-50	34
$Q_{sym}$	MeV	3	389	-28	-67
$Z_{sym}$	MeV	4	-2191	-1315	-1546
$m_{sat}^*/m$			0.80	0.65	0.71
$\Delta m_{sat}^*/m$			0.03	-0.08	-0.09

Table 2.1: Value of each of the empirical parameters ,associated unit, and derivative order  $N$  for BSk14 [GSP07], PKDD [Lon+04], and TM1 [SKT95] functionals.

and  $R = 10$  km [HPY07].

In fact, the central density inside a specific NS is not known, and is expected to range from  $\approx 4.6 \times 10^{14}$  g/cm<sup>3</sup> to  $\approx 4 \times 10^{15}$  g/cm<sup>3</sup>. For this reason, we are rather interested in the mass-radius relation, which can be obtained by calculating  $M$  and  $R$ , following the numerical method previously explained, for this range of central mass densities.

Let us recall that one has to redefine the high-order parameters  $Q_{sat(sym)}$  and  $Z_{sat(sym)}$  in order to reproduce existing functionals at densities greater than  $2n_{sat}$  with the meta-modeling technique. This corresponds to the ELFd technique introduced in [MHG18a].

The left panel of Fig. 2.1 shows the mass-radius relation for several popular EoS based on Skyrme-type functionals BSk24, SLy4, and BSk14, and relativistic models PKDD and TM1, calculated within the metamodeling technique. A strong model dependence is observed, with radii of canonical NS ranging from  $R_{1.4} = 11.5$  km for BSk14 to  $R_{1.4} = 14.5$  km for TM1. More specifically, it is seen that the radius is positively correlated to the slope of the symmetry energy  $L_{sym}$ . The value of the empirical parameters associated to BSk14, TM1, and PKDD are reported in Table 2.1, whereas those of BSk24 and SLy4 can be found in Table 1.3.

The bands correspond to the measured masses of the two most massive pulsars observed to the present day. The mass of PSR J0348+0432 was precisely estimated to  $(2.01 \pm 0.04)M_{\odot}$  [Ant+13], and the very recent relativistic Shapiro delay measurements of PSR J0740+6620 led to  $2.14_{-0.09}^{+0.10}M_{\odot}$  (68.3% credibility interval) [Cro+20]. Naturally, if an EoS cannot insure such high masses, it should not be considered as reliable, in particular at high density, since the maximum mass is determined by the stellar core EoS. In particular, it is seen that the maximum mass corresponding to the BSk14 functional,  $1.8M_{\odot}$ , is lower than the measured mass of PSR J0348+0432. Also, while the SLy4 EoS satisfy this constraint, the maximum mass for this EoS,  $2.05M_{\odot}$ , is lower than that of PSR J0740+6620. This outline the fact that measuring the mass of pulsars is important,



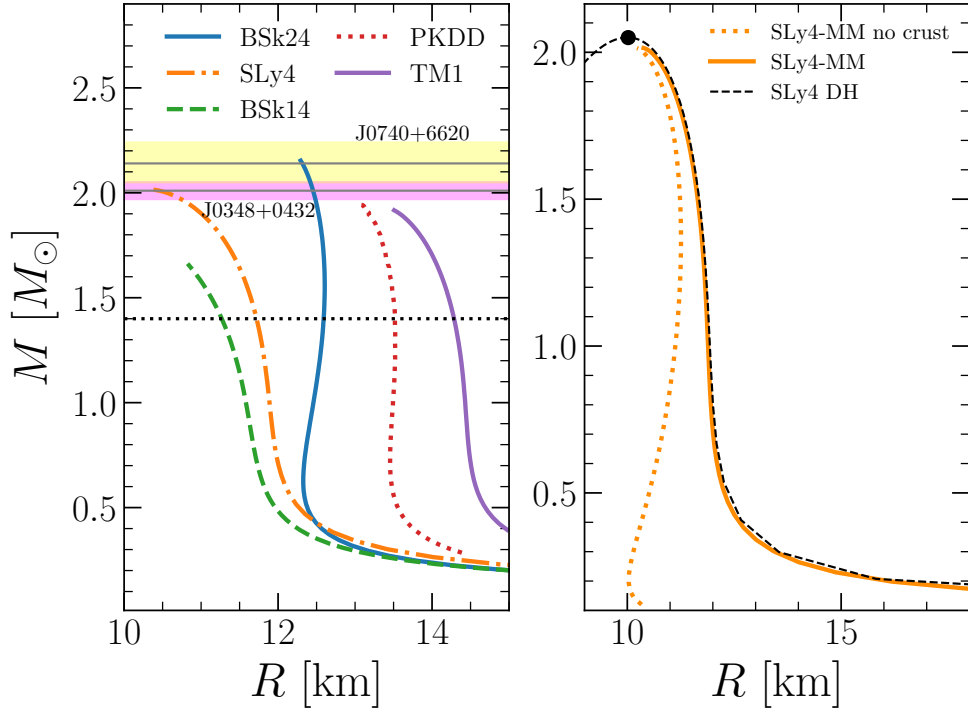


Figure 2.1: Left: Mass-radius relation for several popular EoS calculated with the meta-modeling technique. The black dotted line marks the NS canonical mass, that is  $1.4M_{\odot}$ . The magenta band represents the measured mass of PSR J0348+0432,  $(2.01 \pm 0.04)M_{\odot}$  [Ant+13], and the yellow band that of PSR J0740+6620,  $2.14^{+0.10}_{-0.09}M_{\odot}$  (68.3% credibility interval) [Cro+20]. Right: Mass-radius relation for the DH EoS for SLy4 functional (black dashed line), and the SLy4 EoS calculated with the meta-modeling technique (solid orange line). The black circle indicates the maximum mass.

because it provides a strong constraint on the stellar matter EoS. Similarly, the NICER telescope is expected to provide, for years to come, a constraint on the NS radii with a precision of 5%. Ultimately, in the ideal case where one could accurately measure the mass and radius of one neutron star, it would be possible to determine the EoS by positioning the point in the  $M - R$  diagram.

The mass-radius relation for the SLy4 EoS is represented in the right panel of Fig. 2.1. The solid orange line corresponds to  $M(R)$  for the SLy4 EoS calculated within the meta-modeling technique, and the dashed black line is the calculation based on the DH EoS for SLy4 functional. A perfect agreement is observed between the two curves, reflecting the small error on the EoS, illustrated in Fig. 1.15.

The dotted orange line shows the mass-radius relation for the SLy4 EoS without considering the clustering of matter at low density, that is assuming that the NS star interior consists of  $npe\mu$  matter at all densities. In that case, for a canonical NS, we find  $R_{1.4} = 11.1$  km, which is almost 1 km lower with respect to the result with a crust EoS.



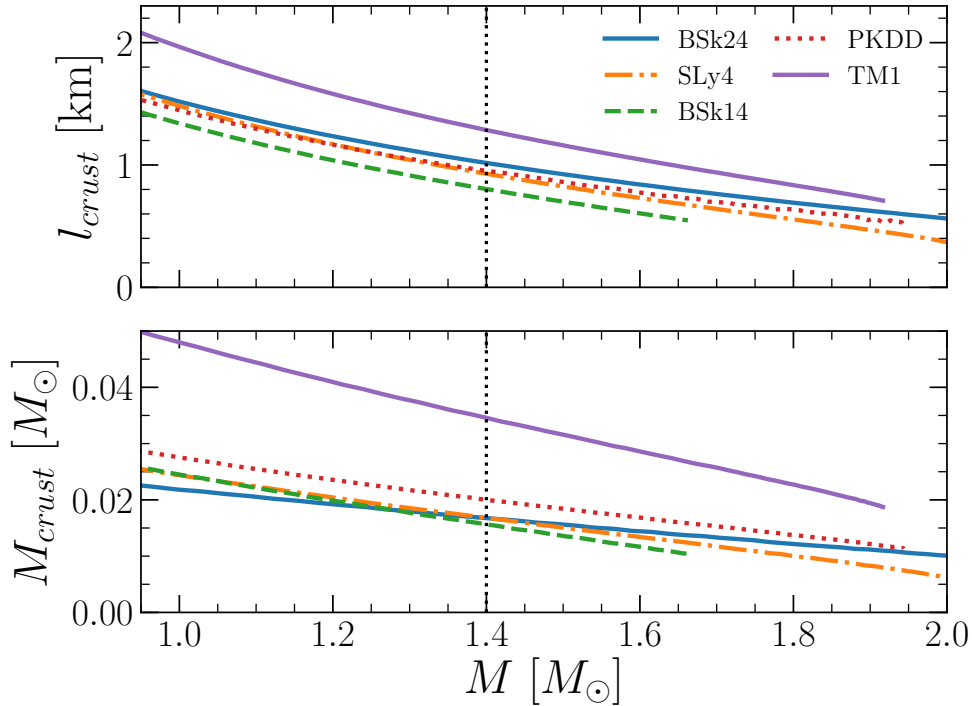


Figure 2.2: Crust thickness  $l_{crust}$  (upper panel) and mass  $M_{crust}$  (lower panel) as a function of the NS mass  $M$  for several popular EoS calculated within the metamodeling technique. The black dotted line marks the NS canonical mass, that is  $1.4M_{\odot}$ .

Moreover, this difference is found to be larger with decreasing mass. This shows that the crust EoS is essential to properly predict NS radii. However, we can see that the NS maximum mass is entirely determined by the stellar core EoS.

The black point marks the maximum mass for the SLy4 EoS,  $M_{max} = 2.05M_{\odot}$ . Let us notice that the branch left to the point is unstable, because from this point the mass decreases with central density increasing.

### 2.1.1.2 Crust thickness and mass

As explained in Chapter 1, a precise estimation of the CC transition point is required as far as crustal properties are concerned. In particular, the transition pressure  $P_t$  is essential for the determination of the thickness and mass of the crust, given respectively by  $l_{crust} = R - R_{core}$  and  $M_{crust} = M - M_{core}$ , with  $R_{core}$  ( $M_{core}$ ) the radius (mass) of the stellar core. Indeed, in order to calculate the core radius and mass, which are involved in the calculation of the crustal observables, one has to integrate the hydrostatic equilibrium equations, Eqs. (2.1) and (2.2), from  $r = 0$  to  $r = R_{core}$ , defined as  $P(r = R_{core}) = P_t$ .

Fig. 2.2 shows the variation with NS mass of crust thickness  $l_{crust}$  and mass  $M_{crust}$  for the same EoS as in Fig. 2.1. As previously, the high-order empirical parameters are

reevaluated to calculate the EoS entering into the TOV equation. However, the transition pressure is calculated using the values of  $Q_{sat(sym)}$  and  $Z_{sat(sym)}$  directly derived from the selected functionals. The reason is that, as showed in Fig. 1.12, the determination of CC transition point is sensitive to orders  $N > 2$ .

We find the interesting result that the crust is thicker for low-mass neutron stars, and in consequence that the  $M_{crust}$  drops continuously with increasing NS mass. At  $M = 1.4M_\odot$  (black dotted line), we observe that the crust is  $\approx 1$  km thick, and that the crust mass is approximately  $0.02 - 0.04M_\odot$ , which is  $(1.5 - 3)\%$  of the total mass. Once again, a model dependence is observed. Nonetheless, the correlation of the crust thickness with the slope of the symmetry energy is not clear. In Fig. 2.1, we have seen that the radius of the star is positively correlated with  $L_{sym}$ . In fact, the same is true for the core radius, explaining why the correlations cancel in the crust thickness  $l_{crust} = R - R_{core}$ .

### 2.1.2 Moment of inertia within the slow rotation approximation

We now turn to the calculation of the moment of inertia within the slow rotation approximation [Har67]. In fact, this approximation is not only applicable to slowly rotating pulsars but is also reliable for most of them. Indeed, while many observed pulsars show rapid rotation [SLL13], their structure is almost not altered by it since centrifugal forces are small in comparison to the gravity,  $R^3\Omega^2/(GM) \ll 1$ ,  $\Omega$  being the angular frequency. For instance, let us consider a rapidly rotating pulsar with  $\Omega = 1000 \text{ s}^{-1}$  and canonical values for its mass and radius,  $M = 1.4M_\odot$  and  $R = 10$  km. In that case, we find  $R^3\Omega^2/(GM) \approx 0.025$ .

In the following, we first calculate the total moment of inertia and the fraction of it contained in the NS crust for several popular EoS. Then, we explain the connection between the fraction of crust moment of inertia and the glitch behavior exhibited by some pulsars [Esp+11].

#### 2.1.2.1 Total moment of inertia and fraction contained in the crust

In a slowly rotating NS, the total moment of inertia is given by [Har67]

$$I = \frac{8\pi}{3} \int_0^R dr r^4 \left( \rho_B + \frac{P}{c^2} \right) \frac{\bar{\omega}}{\Omega} e^{-\lambda - \Phi}, \quad (2.5)$$

where we have introduced the rotational drag function  $\bar{\omega}(r)$ , which satisfies

$$\frac{d}{dr} \left( r^4 j \frac{d\bar{\omega}}{dr} \right) = -4r^3 \bar{\omega} \frac{dj}{dr}, \quad (2.6)$$

with  $j = e^{-\Phi - \lambda}$ . Eq. (2.6) satisfy the following boundary conditions at the surface and center of the NS, respectively:

$$\frac{1}{\Omega} \bar{\omega}(r = R) = 1 - \frac{2GI}{R^3 c^2} \quad \text{and} \quad \left. \frac{d\bar{\omega}}{dr} \right|_{r=0} = 0. \quad (2.7)$$

One can translate Eq. (2.6) into a first-order differential equation by introducing  $w = (1/\Omega)d \ln \bar{\omega} / d \ln r$ , yielding

$$\frac{dw}{dr} = \frac{4\pi G}{c^2} \frac{(P + \rho_B c^2)(4 + w)r^2}{rc^2 - 2Gm} - \frac{w}{r}(3 + w), \quad (2.8)$$



Figure 2.3: Moment of inertia  $I$  (upper panel) and fraction of crust moment of inertia  $I_{\text{crust}}/I$  (lower panel) as a function of the NS mass  $M$  for several popular EoS calculated within the metamodeling technique. In the upper panel, the black dotted line marks the NS canonical mass, that is  $1.4M_{\odot}$ . In the lower panel, the minimum values needed to justify Vela glitches with (Delsate *et al.* [Del+16] and Andersson *et al.* [And+12]) and without (Link *et al.* [LEL99]) crustal entrainment are represented.

with the boundary condition  $w(r = 0) = 0$ . This equation is solved together with Eqs. (2.1) and (2.2) in order to calculate the total moment of inertia, which can be rewritten as

$$I = \frac{c^2}{G} \frac{w(R)R^3}{6 + 2w(R)}. \quad (2.9)$$

The fraction of crust moment of inertia can also be calculated as [LHS19]

$$\begin{aligned} \frac{I_{\text{crust}}}{I} &= 1 - \frac{I_{\text{core}}}{I} \\ &= 1 - \left(\frac{R_{\text{core}}}{R}\right)^3 \frac{w(R_{\text{core}})}{w(R)} \exp \left[ - \int_{R_{\text{core}}}^R \frac{w(r)}{r} dr \right], \end{aligned} \quad (2.10)$$

where we have introduced the moment of inertia of the core  $I_{\text{core}}$ , defined by the integration of Eq. (2.5) up to the core radius  $R = R_{\text{core}}$ .

The total moment of inertia is plotted as a function of the NS mass in the upper panel of Fig. 2.3 for EoS based on Skyrme-type functionals BSk24, SLy4, and BSk14,

and relativistic models PKDD and TM1. For  $M = 1.4M_\odot$ , it is found that  $I$  ranges from  $\approx 1.2 \times 10^{45} \text{ g cm}^2$  for the soft BSk14 EoS to  $\approx 2 \times 10^{45} \text{ g cm}^2$  for the very stiff TM1 EoS. This shows the strong dependence on the EoS. More particularly, an apparent positive correlation with  $L_{sym}$  is observed. It is seen that the model dependence is even larger for high-mass NS.

The lower panel of Fig. 2.3 shows the variation with NS mass of  $I_{crust}/I$  for the same EoS. It is seen that the more massive a NS is, the less moment of inertia it stores in its crust, which follows the trends observed in Fig. 2.2. This was predictable from the sensitivity of crustal properties to the CC transition pressure. Given the selected EoS, the value of  $I_{crust}/I$  at  $M = 1.4M_\odot$  ranges from  $\approx 2\%$  for BSk14 to  $\approx 5\%$  for TM1, showing once again the dramatic dependence on the EoS.

### 2.1.2.2 Connection to pulsar glitches

The so-called pulsar glitches are sudden jumps in the rotational frequency of a compact star. They are thought to originate from an abrupt transfer of angular momentum from the liquid interior, acting as an angular momentum reservoir, to the solid crust of the star due to the unpinning of the superfluid vortices from the crystal lattice [AI75]. This occurs when the differential lag between the slower solid crust and faster superfluid vortices reach some threshold and can no longer be sustained, so as to recover a close equilibrium between the two. Eventually, stresses start to build up again, ultimately leading to another glitch event.

At the time of writing, 555 glitches have been observed in 190 pulsars through high-precision pulsar timing [Esp+11; Jod]. For a given glitcher, three quantities are usually measured: its spin frequency  $\Omega$ , its average spin-down rate  $\dot{\Omega}$ , and its glitch activity parameter  $\mathcal{A}$ , defined as

$$\mathcal{A} = \frac{1}{t} \sum_{i=1}^N \frac{\Delta\Omega_i}{\Omega}, \quad (2.11)$$

where  $\sum_i \Delta\Omega_i/\Omega$  represents the cumulative spin-up rate over the  $N$  glitches occurring during a time  $t$ .

It was demonstrated in [LEL99] that the ratio of the moment of inertia associated to the neutron superfluid which drives the glitches  $I_{sf}$  to the moment of inertia of the solid crust – plus any portion of the star strongly coupled to it –  $I_c$  must satisfy

$$\frac{I_{sf}}{I_c} \geq \frac{\Omega}{|\dot{\Omega}|} \mathcal{A} \equiv \mathcal{G}, \quad (2.12)$$

where we have introduced the coupling parameter  $\mathcal{G}$ . In view of relating the observational constraint  $\mathcal{G}$  to the fractional moment of inertia  $I_{crust}/I$  to ultimately set a constraint on the stellar EoS, we assume that the angular momentum reservoir is confined to the neutron superfluid coexisting with the crystal lattice of the inner crust. In that case, we have  $I_c = I - I_{res} \simeq I$  and  $I_{crust} \simeq I_{sf}$ , yielding [LEL99]

$$\frac{I_{sf}}{I_c} \simeq \frac{I_{crust}}{I} \geq \mathcal{G}. \quad (2.13)$$

In reality, the neutron superfluid is strongly coupled to the solid crust due to nondissipative entrainment effects [Cha13], limiting the amount of angular momentum that can be transferred during a glitch event. The importance of the entrainment coupling is related to the neutron effective mass  $m_n^*$  in the inner crust, which is proportional to the ratio of unbound neutrons to those that are not entrained,  $m_n^* \equiv m_n n_g / n_g^c$ . Taking into account the entrainment effects, the previous constraint, Eq. (2.13), then becomes [And+12]

$$\frac{I_{crust}}{I} \geq \mathcal{G} \frac{\langle m_n^* \rangle}{m_n}, \quad (2.14)$$

with  $\langle m_n^* \rangle$  the average effective neutron mass. It should be stressed that systematic calculations of  $m_n^*$  throughout the NS crust are very costly from the computational point of view and that, following [And+12; PFH14], we adopt the value  $\langle m_n^* \rangle / m_n = 4.3$ , which is inferred from the results of [Cha12]. Naturally, one recovers the expression without crustal entrainment in the case of  $\langle m_n^* \rangle = m_n$ . We should stress that the importance of the effect of crustal entrainment, that is the value of the average neutron effective mass throughout the solid crust, is currently under debate [MU16; WP17].

The Vela pulsar (B083345-45) is one of the most active glitchers known, with glitches occurring four times per decade on average. Since the beginning of its monitoring, almost than 50 years ago, it has exhibited 20 glitches in total [Jod], allowing for a precise estimation of the corresponding coupling parameter,  $\mathcal{G}_{Vela} = (1.62 \pm 0.03) \times 10^{-2}$  [Ho+15]. This can be translated into a constraint for the fraction of crust moment of inertia,

$$\left( \frac{I_{crust}}{I} \right)_{Vela} \geq (1.62 \pm 0.03) \times 10^{-2} \frac{\langle m_n^* \rangle}{m_n} \approx 0.07. \quad (2.15)$$

In other words, in the standard picture where the angular momentum reservoir is exclusively confined to the neutron superfluid permeating the inner crust, the crust must store more than 7% (with entrainment effects included) of the total moment of inertia in order to justify the large glitches occurring in the Vela pulsar.

The Vela constraints on  $I_{crust}/I$  are displayed in the lower panel of Fig. 2.3. Neglecting entrainment effects [LEL99], it is seen that the constraint is easily satisfied. For BSk14 EoS, which is the softest EoS here, it leads to  $\approx 1.6M_\odot$  for the maximum mass of Vela. Now including crustal entrainment with  $\langle m_n^* \rangle / m_n = 4.3$  [And+12], it is not clear, due to dependence on the EoS, whether the superfluid neutron in the crust can carry enough angular momentum to explain Vela glitches. For the three nonrelativistic Skyrme-based EoS that we consider, we see that the constraint is unlikely to be satisfied, because it would require the Vela mass to be very low,  $M_{Vela} \lesssim 0.9M_\odot$ . While theoretically plausible [HPY07], such low-mass NS have never been observed to this day. On the other hand, it is observed that very stiff EoS such as TM1 can be consistent with the constraint [PFH14]. Finally, given the constraint on the fractional moment of inertia calculated with the present largest estimation of crustal entrainment [Del+16] (black thick line), we find that the inferred mass of Vela is at most  $\approx 0.85M_\odot$  for the stiff TM1 EoS and  $\approx 0.65M_\odot$  for the soft EoS based on Skyrme functionals. Clearly the question “*Is the crust enough?*” deserves further investigations. In that sense, it will be addressed in 2.3.3 in the context of a complete statistical analysis [CGM19b].

### 2.1.3 Tidal deformability

The tidal deformability describes how much a body is deformed by tidal forces, which arise when two massive bodies orbit each other. The simplest example corresponds of course to the Moon causing the tides observed in Earth's oceans. The detection of gravitational radiation emitted by inspiraling binary NS is possible using ground-based gravitational wave (GW) detectors such as LIGO and Virgo. Shortly before merging, once the relative distance between the stars is small enough, the tidal distortion of the NS becomes so large that, in some cases (in strongest signals), it becomes possible to infer the tidal deformability from the GW signal. Theoretically, the tidal deformability parameter is the observable which makes the link between the EoS and the gravitational signal. In that sense, GW observations can ultimately provide new constraints on the EoS. In particular, the first detection of GW from the coalescence of two NS, the GW170817 event, have yielded a value for the tidal deformability of a  $1.4M_\odot$  NS [Abb+18].

Let us consider a static, spherically symmetric NS, placed in a static external quadrupolar tidal field  $\mathcal{E}_{tid}$ . The induced quadrupole moment of this star is given by

$$Q_{tid} = -\lambda \mathcal{E}_{tid}, \quad (2.16)$$

where  $\lambda$  is the tidal deformability, which is related to the so-called tidal Love number  $k_2$  by

$$\lambda = \frac{2}{3} k_2 \left( \frac{Rc^2}{G} \right)^5. \quad (2.17)$$

Let us also define the dimensionless tidal deformability,

$$\Lambda = \frac{\lambda}{M^5} = \frac{2}{3} k_2 \beta^{-5}, \quad (2.18)$$

where we have introduced the compactness of the star  $\beta = GM/(Rc^2)$ . Similarly to the moment of inertia in the slow rotation approximation,  $k_2$  can be obtained from the solution of the following first-order differential equation [Hin+10],

$$\begin{aligned} \frac{dy}{dr} = & -\frac{y^2}{r} - \frac{y-6}{r-2Gm/c^2} \\ & - \frac{4\pi G}{c^2} r^2 \frac{(5-y)\rho_B + (9+y)P/c^2 + (P + \rho_B c^2)/c_s^2}{r-2Gm/c^2} \\ & + \frac{1}{r} \left[ \frac{2G}{c^2} \frac{(m + 4\pi p r^3/c^2)}{r-2Gm/c^2} \right]^2, \end{aligned} \quad (2.19)$$

where we have introduced the speed of sound  $c_s = \sqrt{\partial P / \partial \rho_B}$ , and that must satisfy the boundary condition  $y(r=0) = 2$ . This equation together with the hydrostatic equilibrium equations, Eqs. (2.1) and (2.2), are solved simultaneously. The tidal Love number reads

$$k_2 = \frac{8}{5} \beta^5 (1-2\beta)^2 [2 - y(R) + 2\beta(y(R) - 1)]/a, \quad (2.20)$$

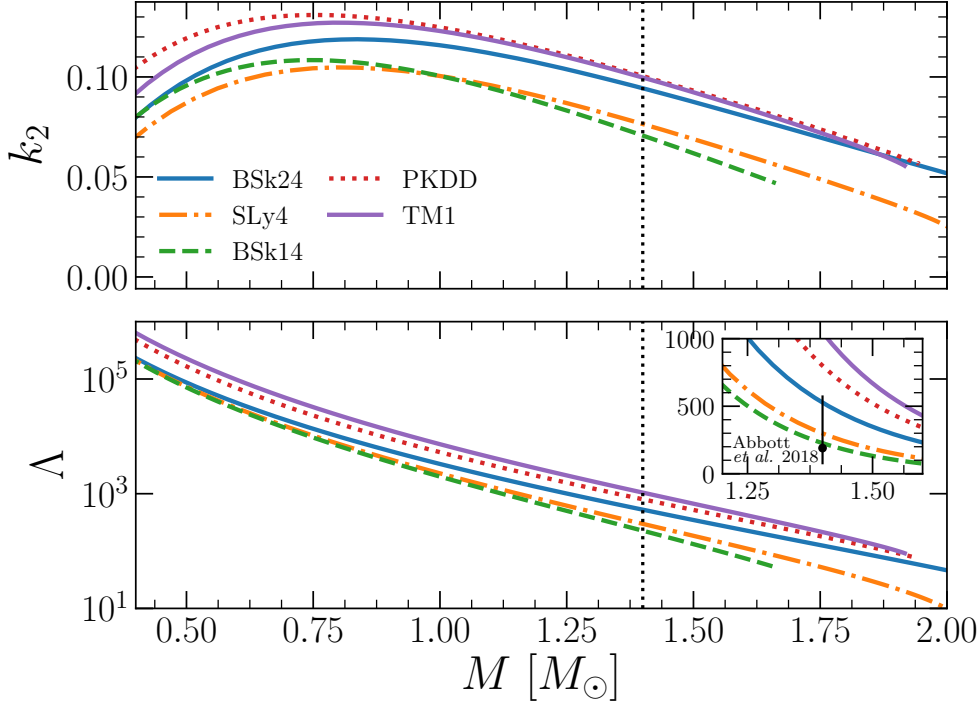


Figure 2.4: Tidal Love number  $k_2$  (upper panel) and dimensionless tidal deformability  $\Lambda$  (lower panel) as a function of the NS mass  $M$  for several popular EoS calculated within the metamodeling technique. The black dotted line marks the NS canonical mass, that is  $1.4M_\odot$ . In the inset of the lower panel, the value  $\Lambda_{1.4} = 190^{+390}_{-120}$  (at 90% level) inferred from GW170817 event is represented [Abb+18].

with

$$\begin{aligned}
 a = & 6\beta[2 - y(R) + \beta(5y(R) - 8)] \\
 & + 4\beta^3[13 - 11y(R) + \beta(3y(R) - 2) + 2\beta^2(1 + y(R))] \\
 & + 3(1 - 2\beta)^2[2 - y(R) + 2\beta(y(R) - 1)] \ln(1 - 2\beta).
 \end{aligned} \tag{2.21}$$

Fig. 2.4 shows the variation with NS mass of  $k_2$  for the several popular EoS that we consider in this chapter. The tidal Love number measures how easily the bulk of the matter in a NS is deformed. In the case of the mass being concentrated at the center of the star, the tidal deformation is small, and so is  $k_2$ . In agreement with [Hin+10], we find that  $k_2$  lies in the range  $\approx 0.05 - 0.15$ . The relative difference between the present EoS is of the order of  $\approx 35\%$  for  $M = 1.4M_\odot$ , showing once again the dramatic dependence on the EoS. Let us point out that the tidal Love number is very sensitive to the description of the crustal component of the EoS [PF19].

The dimensionless tidal deformability is plotted as a function of  $M$  in the lower panel of Fig. 2.4. In the mass range considered, it is seen that the more massive the NS is, the less it gets deformed when orbiting another massive compact object. In the inset,

we show the value of the dimensionless tidal deformability for a  $1.4M_\odot$  NS inferred from the GW170817 event,  $\Lambda_{1.4} = 190^{+390}_{-120}$  at the 90% level. We can observe that the EoS which predict the largest tidal effects, PKDD and TM1, are unlikely. One can notice the tug or war between the tendency of wanting soft EoS, typically based on nonrelativistic Skyrme-type functionals, which predict small tidal deformations, and wanting very stiff EoS, typically based on relativistic models, to explain the glitch phenomenon as a transfer of angular momentum from the neutron superfluid confined in the crystal lattice to the solid crust.

## 2.2 Bayesian determination of the equation of state parameters

### 2.2.1 Principle of Bayesian inference

The core of Bayesian inference is based on well-known Bayes' theorem, which provides an expression for the conditional probability (or posterior probability) of  $A$  given  $B$ ,

$$P(A|B) = \frac{P(B|A)P(A)}{P(B)}. \quad (2.22)$$

In this equation,  $P(B|A)$  is the conditional probability of  $B$  given  $A$ ,  $P(A)$  is the prior probability of  $A$ , and  $P(B)$  that of  $B$ . In the model form of Bayes' theorem, one replaces  $B$  by data,  $A$  with a parameter set  $\mathbf{X}$ , and  $P$  with probability density functions (PDF)  $p$ , resulting in [SS06]

$$p(\mathbf{X}|\text{data}) = \frac{p(\text{data}|\mathbf{X})p(\mathbf{X})}{p(\text{data})}. \quad (2.23)$$

The denominator is not a distribution but a normalization constant which makes sure that the posterior distribution  $p(\mathbf{X}|\text{data})$  is a true probability distribution, by ensuring that the sum of the distribution is equal to 1. It is given by

$$p(\text{data}) = \int p(\text{data}|\mathbf{X})p(\mathbf{X})d\mathbf{X}. \quad (2.24)$$

The model-based formulation of Bayes' theorem, Eq. (2.23), is therefore often written as

$$p(\mathbf{X}|\text{data}) \propto p(\text{data}|\mathbf{X})p(\mathbf{X}). \quad (2.25)$$

The marginal one- and two-parameter posterior distributions can be extracted from the posterior distribution. Then are defined respectively as

$$p(X_j|\text{data}) = \left\{ \prod_{i \neq j} \int dX_i \right\} p(\mathbf{X}|\text{data}), \quad (2.26)$$

$$p(X_j, X_k|\text{data}) = \left\{ \prod_{i \neq j, k} \int dX_i \right\} p(\mathbf{X}|\text{data}). \quad (2.27)$$

Let us describe the different components of Bayesian inference. The prior distribution  $p(\mathbf{X})$  represents our knowledge or bias on the model parameters. It can thus be more



or less informative.  $p(\text{data}|\mathbf{X})$  is the likelihood of observing the data given the model parameters  $\mathbf{X}$ . It is determined from the data comparison between the model and the measurement via an error estimator. In that sense, the link between the data and the model parameters is encoded in the likelihood distribution. Multiplying the prior by the likelihood function, one obtains the unnormalized joint posterior PDF, which corresponds to the conditional distribution of model parameters  $\mathbf{X}$  given the data. Let us notice that the prior distribution affects the posterior distribution.

Bayesian inference is an appealing statistical method in the sense that it allows to calculate the joint posterior distribution of model parameters using prior beliefs updated with the likelihood. In the context of NS physics, and from the nuclear physicist point of view, it allows for instance to update our knowledge on the nuclear EoS by taking into consideration the data arising from various astrophysical observations, that are for example the NS mass inferred from Shapiro delay or the tidal deformability from GW measurements. Conversely, one can account for nuclear physics constraints in the prediction of macroscopic observables related to NS. In our case, we first generate a large number of EoS, each of them being defined by a set of parameters  $\mathbf{X}$ , to numerically sample the prior parameter distribution. Then, we filter those EoS through constraints, which arise from physical requirements, astrophysical observations, and ab initio calculations of SNM and PNM. This gives us the joint posterior distribution of model parameters, from which we can finally calculate the posterior EoS and make general predictions for NS observables. Those predictions are presented in Section 2.3.

Let us describe the parameter set  $\mathbf{X}$ . Our calculation of the stellar EoS is based on the metamodeling of nuclear matter energy [MHG18a] introduced in 1.2.1.1. The empirical parameters  $E_{\text{sat}}$ ,  $n_{\text{sat}}$ ,  $K_{\text{sat}}$ ,  $Q_{\text{sat}}$ ,  $Z_{\text{sat}}$ ,  $E_{\text{sym}}$ ,  $L_{\text{sym}}$ ,  $K_{\text{sym}}$ ,  $Q_{\text{sym}}$ , and  $Z_{\text{sym}}$  – which we recall are given by the successive derivatives of the NM energy at saturation density in isoscalar and isovector sectors – therefore enter into the parameter set. The parameter  $b$  is also added in order to explore different functional behaviors close to the  $n \rightarrow 0$  limit. The parameter  $b$ , together with the isoscalar effective mass  $m_{\text{sat}}^*/m$  and isospin splitting  $\Delta m_{\text{sat}}^*/m$ , completes the parameter space, associated to the description of bulk matter. In the solid crust, the metamodeling technique is supplemented by a surface plus curvature term in order to describe the surface properties of finite nuclei. Thus, one could consider adding the surface and curvature parameters, respectively  $\sigma_0$ ,  $b_s$ ,  $p$ , and  $\sigma_{0c}$ ,  $\beta$ , as extra parameters to  $\mathbf{X}$ . However, we have seen in Fig. 1.5 that the  $\chi^2$  function, which evaluates the goodness of the fit of those parameters to experimental nuclear masses given BSk24 empirical parameters and  $p = 3$ , is strongly peaked on a given set  $\mathbf{S} = \mathbf{S}_{\text{opt}}$ . We expect this feature to be recovered for every model and we therefore choose to not include the surface and curvature parameters in  $\mathbf{X}$  but to take the set  $\mathbf{S}_{\text{opt}}$  that minimizes  $\chi^2$ , as explained in 1.2.1.2. We consider two choices for the value of the parameter  $p$ , which determines the behavior of the surface tension for extreme isospin values. Either it is fixed to the educated value  $p = 3$  or  $p = \{2.5, 3, 3.5\}$  is considered for each set of empirical parameters.

Parameter	Unit	$N$	Min	Max
$E_{sat}$	MeV	0	-17	-15
$n_{sat}$	$\text{fm}^{-3}$	1	0.15	0.17
$K_{sat}$	MeV	2	190	270
$Q_{sat}$	MeV	3	-1000	1000
$Z_{sat}$	MeV	4	-3000	3000
$E_{sym}$	MeV	0	26	38
$L_{sym}$	MeV	1	10	80
$K_{sym}$	MeV	2	-400	200
$Q_{sym}$	MeV	3	-2000	2000
$Z_{sym}$	MeV	4	-5000	5000
$m_{sat}^*/m$			0.6	0.8
$\Delta m_{sat}^*/m$			0.0	0.2
$b$			1	10

Table 2.2: Minimum value and maximum value of each of the parameters of parameter set  $\mathbf{X}$ . The associated unit and derivative order  $N$  are also reported.

### 2.2.2 Prior distribution of empirical parameters

We now discuss the prior distribution of empirical parameters. We first give insights concerning the selected intervals for the flat prior. A sensitivity analysis of the CC transition point is performed afterwards.

#### 2.2.2.1 Flat prior compatible with empirical constraints

The prior distribution of  $\mathbf{X}$  is given by an uncorrelated ansatz and an uniform distribution of each parameter within the interval specified in Table 2.2,

$$p(\mathbf{X}) = \prod_{i=1}^{2(N+1)+3} \mathcal{U}(X_i^{min}, X_i^{max}; X_i), \quad (2.28)$$

where the parameter  $X_i$  is uniformly distributed from  $X_i^{min}$  to  $X_i^{max}$ . The use of a flat prior means that all possible values of  $\mathbf{X}$  inside the intervals are equally likely a priori.

The range of variation for the model parameters reflect the degree of uncertainty on the EoS parameters, as measured by their variation in the functionals which reproduce successfully low energy nuclear physics data [MHG18a]. We can distinguish three groups of empirical parameters, depending on how well they are experimentally constrained. The first group consists of the low-order isoscalar parameters  $E_{sat}$ ,  $n_{sat}$ ,  $K_{sat}$ , the symmetry energy at saturation density  $E_{sym}$ , as well as the effective mass  $m_{sat}^*$  and isospin splitting  $\Delta m_{sat}^*/m$ . These parameters are rather well determined by nuclear experiments, with associated relative uncertainties lower than 15%. The second group consists of the parameters that are still poorly determined by nuclear experiments. The uncertainties on the isoscalar skewness  $Q_{sat}$ , the slope of the symmetry energy  $L_{sym}$ ,

and the isovector incompressibility  $K_{sym}$  are very large, yet one can expect a better accuracy to be reached in the near future. In particular, the determination of  $L_{sym}$  is an highly topical issue nowadays [Li+14]. Finally, the last group concerns the parameters which are inaccessible by nuclear experiments to the present day, that is the isovector skewness  $Q_{sym}$  as well as the parameters of order  $N = 4$ ,  $Z_{sat}$  and  $Z_{sym}$ . In consequence, we explore very large ranges for these parameters.

The choice of the interval for the low-density parameter,  $b = [1, 10]$ , is discussed thoroughly in [Ant+19].

### 2.2.2.2 Sensitivity analysis of the crust-core transition point\*

Let us recall that one of the advantages of the metamodeling approach is that all the EoS model parameters, such as the empirical parameters, are a priori uncorrelated. It is therefore possible to vary each one of them independently of the others, which is not feasible using specific functional behaviors such as Skyrme, Gogny, or the different versions of RMF, as discussed in detail in [MHG18a; MG19]. Indeed, common functionals describing infinite NM have a smaller number of parameters compared to the metamodel, which creates a priori correlations that are not fully controlled. If we take the example of Skyrme forces, the density dependence of SNM is controlled by only three parameters, which are fixed to reproduce the coordinates of the saturation point:  $n_{sat}$ ,  $E_{sat}$ , and  $K_{sat}$ . Thus, the isoscalar skewness  $Q_{sat}$  is correlated with the three other parameters, and it turns out that this correlation is not the same, for instance, in Skyrme, Gogny, or in RMF functionals [MG19]. A similar argument applies to the symmetry energy: when the density dependence of the symmetry energy is governed by a single parameter as assumed in heavy ion transport models, an isosoft behavior at suprasaturation density is necessarily associated to an isostiff behavior at subsaturation density, while more complicated behaviors are possible [MG19]. The independent variation of the different empirical parameters allows a simple determination of the most influential parameters impacting a given observable under study.

Such a sensitivity analysis is presented in Fig. 2.5 for the CC transition density  $n_t$  and transition pressure  $P_t$ . The reference metamodel used in the sensitivity analysis can have an influence. For this reason, the sensitivity analysis varying one by one the EoS parameters is performed around two different reference parameter sets  $\mathbf{X}_{ref}$ , namely the parameter set corresponding to the BSk24 functional (solid blue lines), and the one corresponding to the BSk22 functional (dashed orange lines). For both of them, we fix  $p = 3$  as well as  $b = 10 \ln(2)$ .

The range of variation for each of the EoS parameters correspond to the intervals for the prior distribution  $p(\mathbf{X})$  specified in Table 2.2. Let us however notice that some sets produce nonviable EoS, for which a solution for the CC transition point cannot be found. This occurs in particular when varying the value of the isovector incompressibility parameter  $K_{sym}$  with BSk24 as reference parameter set. Indeed, we find an extremely soft behavior for the symmetry energy for the minimum value  $K_{sym}^{min} = -400$  MeV. In that very situation, we increase the minimum value of the parameter to  $K_{sym}^{min} \approx -300$

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\*Partial results of the presented work have been published in Eur. Phys. J. A (2019) **55**: 188. With kind permission of The European Physical Journal (EPJ).

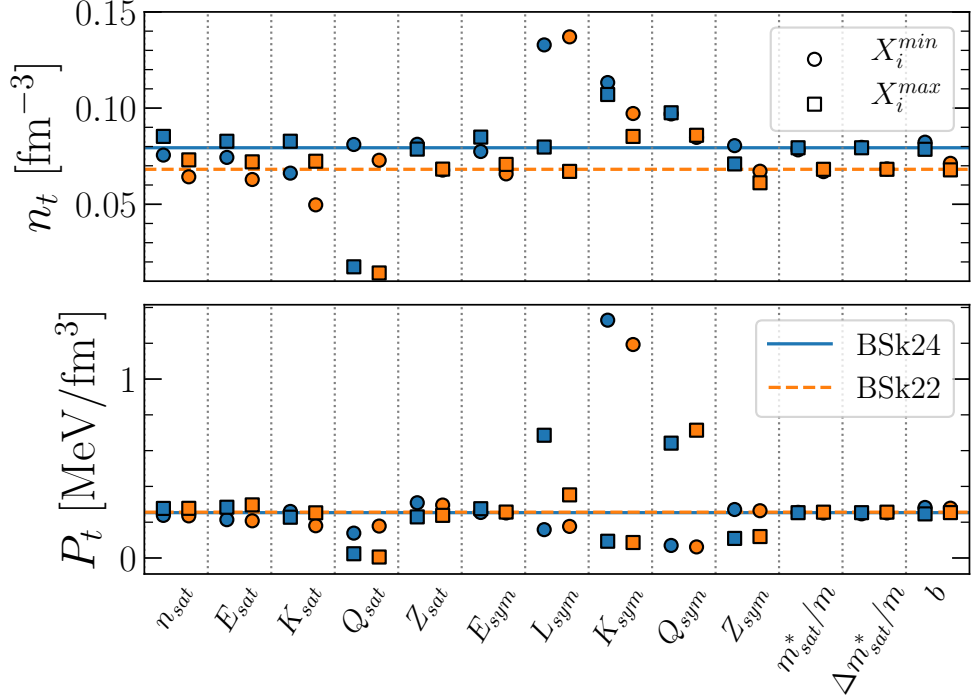


Figure 2.5: Sensitivity analysis of the transition density  $n_t$  (upper panel) and transition pressure  $P_t$  (lower panel) with respect to EoS parameters.  $p = 3$  is fixed and two different reference point are chosen: BSk24 parameters (solid blue lines) and BSk22 parameters (orange dashed lines). For each parameter, the symbols on left (blue) correspond to BSk24 and those on right (orange) to BSk22. Circles and squares are respectively associated to the minimum value  $X_i^{min}$  and maximum value  $X_i^{max}$  of the parameter, which are listed in Table 2.2.

MeV so as to find a solution for  $n_t$  and  $P_t$ .

The symbols in Fig. 2.5 give the transition density and pressure domain obtained when the EoS parameters are one by one varied around the reference model, within the interval of Table 2.2. More specifically, circles and squares are respectively associated to the minimum value and maximum value of each parameter. Since the uncertainty on the different parameters is not the same, the relative distance between circles and squares is a qualitative measurement of the propagation of the uncertainty on the transition point brought by each parameter. We can see that the sensitivity of each parameter depends on the value of the other parameters, that is on the chosen reference set  $\mathbf{X}_{ref}$ . Still, universal trends clearly emerge. We can see that the CC phase transition, at variance with the standard liquid-gas of SNM, is virtually insensitive to isoscalar parameters. Even if extremely large variations of  $Q_{sat}$  and  $Z_{sat}$  are considered, see Table 2.2, the influence of the isoscalar parameters in the transition pressure is almost negligible, and it is a bit larger for  $n_t$  but still smaller than the isovector parameters. This underlines

the importance of the energetics of the neutron gas on the transition point. Concerning the isovector sector, we can see that  $L_{sym}$  is the most important parameter as far as the transition density is concerned. This result is in agreement with previous findings by many authors [Duc+11] and was already anticipated in 1.2.6. The symmetry energy at saturation  $E_{sym}$  and the effective mass splitting  $\Delta m_{sat}^*/m$  do not play any role on the transition, which can be partially explained by the fact that these parameters are already relatively well constrained. The transition density and pressure show also a great sensitivity to the isovector incompressibility parameter  $K_{sym}$ . This can explain why the transition pressure exhibits an irregular behavior when plotted as a function of  $L_{sym}$  (see Fig. 1.13): the different functionals considered in the literature have very different values of  $K_{sym}$ , which blurs the correlation with  $L_{sym}$ . This effect is also amplified by the fact that, depending on the reference point, the dependence of  $P_t$  with  $L_{sym}$  is not monotonic. Finally, we can notice that the influence of the fully unknown high order derivatives  $Q_{sym}$  and  $Z_{sym}$ , though less important than the one of  $L_{sym}$ , is not negligible and comparable to the one of the isovector surface energy parameter  $p$  that can be inferred from Fig. 1.13. We have checked that similar conclusions can be drawn if the sensitivity analysis is performed using the definition of the transition point from the dynamical spinodal [Ant+19].

### 2.2.3 Determination of the likelihood function

The likelihood function corresponds to the probability of observing the data given the model parameters  $\mathbf{X}$ . In the context of our analysis, it is defined as

$$p(\text{data}|\mathbf{X}) = w_{LD}(\mathbf{X}) \times w_{HD}(\mathbf{X}) \times p_{\text{AME2016}}(\mathbf{X}). \quad (2.29)$$

In this expression,  $w_{LD}$  and  $w_{HD}$  are strict filters that concentrate on constraining the low density EoS and high density EoS, respectively. Both filters are given by sharp  $\delta$  functions that output 1 if the constraint is satisfied, and 0 otherwise. Constraints on nuclear physics observables, yielding the filter  $w_{LD}$  and the probability  $p_{\text{AME2016}}$  are presented in 2.2.3.1. In 2.2.3.2, we present the constraints corresponding to the filter  $w_{HD}$ , which are associated to general physical requirements and measurements of NS observables.

#### 2.2.3.1 Constraints on nuclear physics observables

Let us concentrate first on the strict filter  $w_{LD}$ , hereafter referred to as low density (LD) filter. It is introduced to verify whether the metamodel associated to a set of parameter  $\mathbf{X}$  is compatible with the recent chiral effective field theory (EFT) calculations for SNM and PNM by Drischler *et al.* [DHS16]. The authors have calculated the energy per particle for several values of isospin at low density within the many-body perturbation theory, based on a set of seven different Hamiltonians with chiral NN and three-nucleon interactions.

The resulting predicted bands for energy per nucleon as well as pressure of SNM and PNM are shown in Fig. 2.6 (pink bands), together with the predictions of the popular BSk24 (solid blue lines) and DD-ME $\delta$  (green dashed lines) functionals. The pressure band is calculated by taking the density derivative of the energy per nucleon for the

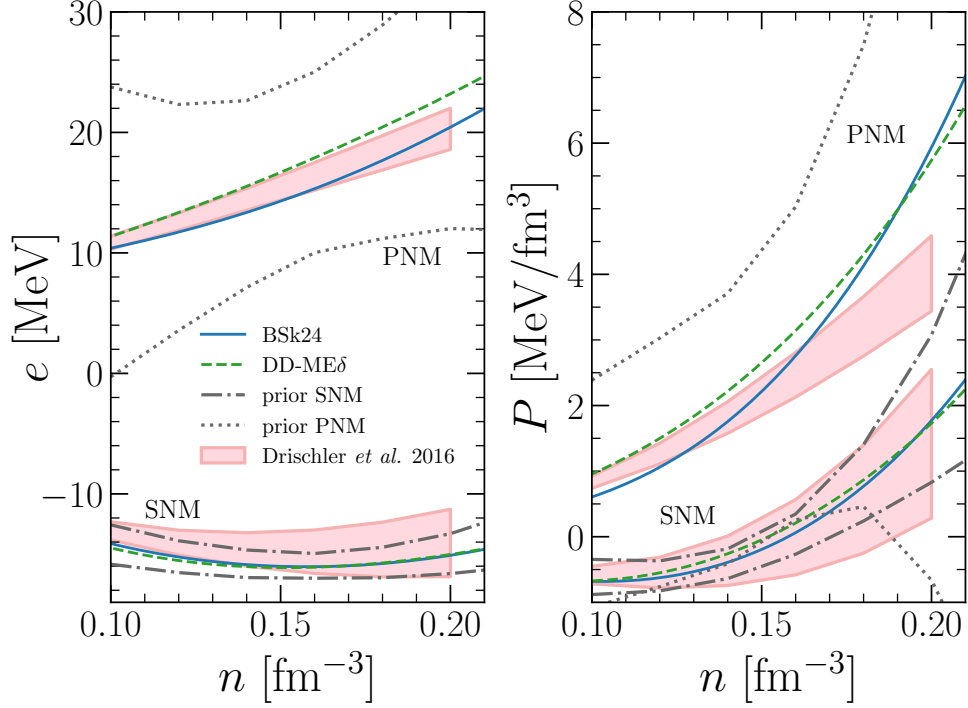


Figure 2.6: Energy per nucleon (left panel) and pressure (right panel) as a function of density for both SNM and PNM from chiral ETF calculations of [DHS16] (pink bands). The predictions of BSk24 and DD-ME $\delta$  functionals are represented in solid blue lines and dashed green lines, respectively. The gray lines give the minimum and maximum values of the prior distribution.

seven Hamiltonians, thus only few density dependencies of the pressure are explored. This can explain why even realistic models such as BSk24 are not compatible with the theoretical band for the pressure of PNM above  $\approx n_{sat}$ . It is therefore debatable whether we should add the chiral EFT bands for pressure of NM in our list of constraints defining our pass-band filters  $w_{LD}$ . In the following, we will not consider this particular constraint by default, yet we will still investigate how it affects the posterior distribution EoS in 2.3. In the context of the pass-band filter  $w_{LD}$ , the chiral EFT predicted bands are widened by 5% in order to account for other existing ab initio calculations in an effective manner. In addition, we will not apply the filter in the low density region  $n < 0.10 \text{ fm}^{-3}$  because the width of the band becomes extremely small, causing numerical issues.

The minimum values and maximum values of the prior distribution are also represented in Fig. 2.6 (gray lines). Let us notice that one can infer the posterior band graphically by the comparison with the chiral EFT bands. Interestingly, the graphically inferred posterior band of the prior energy per nucleon and pressure of SNM around saturation density is narrower than the one corresponding to the ab initio calculation. The reason is that low energy nuclear physics experiments provide strong empirical con-

straints for SNM [MHG18a], which have been embedded effectively in the intervals for the empirical parameters of the prior distribution in Table 2.2. The same is not true for the PNM case. Indeed, the present day knowledge on the empirical parameters of the isovector sector is insufficient. For this reason, we can see that the corresponding ab initio bands are very thin in comparison to the prior distribution.

We now turn the calculation of the probability  $p_{\text{AME2016}}$ , which describe the ability of the CLDM defined from a parameter set  $\mathbf{X}$ , to fit the measured nuclear masses of the AME2016 [Hua+17]. The goodness of fit is evaluated by the reduced  $\chi^2$ , Eq. 1.62. We can define the associated likelihood probability as

$$p_{\text{AME2016}}(\mathbf{X}) = \exp \left[ -\frac{1}{2} \chi^2(\mathbf{X}) \right]. \quad (2.30)$$

### 2.2.3.2 Physical requirements and constraints on neutron star observables

The filter  $w_{HD}$ , hereafter referred to as high density (HD) filter, is applied subsequently to the LD filter  $w_{LD}$ . It imposes general physical constraints to the global density behavior of the functional, as follows:

- the positiveness of the symmetry energy at all densities,
- the stability of the EoS,
- the causality condition,  $0 < c_s/c < 1$ ,
- the maximum observed NS mass,  $M_{\text{max}}(\mathbf{X}) \geq M_{\text{max}}^{\text{obs}}$ ,

where  $M_{\text{max}}(\mathbf{X})$  is the maximum mass supported by the EoS calculated for a parameter set  $\mathbf{X}$ , and  $M_{\text{max}}^{\text{obs}}$  is the maximum observed NS mass. By default we choose  $M_{\text{max}}^{\text{obs}} = 1.97 M_\odot$ , which corresponds to a  $1\sigma$  conservative estimate, based on the observation of PSR J0348+0432 [Ant+13]. Let us however notice that it does not correspond to the mass of the heaviest NS observed to the present day [Cro+20]. The reason is that the observation of PSR J0740+6620 with  $M = 2.14_{-0.09}^{+0.10} M_\odot$  is very recent and it was not yet published at the time of this study.

As we have already discussed, for a given functional, the redefinition of high-order parameters  $Q_{\text{sat}}$ ,  $Z_{\text{sat}}$ ,  $Q_{\text{sym}}$ , and  $Z_{\text{sym}}$  is necessary in order to achieve faster convergence at high density and thus to recover the right value for the maximum mass. The LD filter is applied in the density range  $0.10 \text{ fm}^{-3} < n < 0.20 \text{ fm}^{-3}$ , that is in the vicinity of the saturation density  $n_{\text{sat}}$ . In this density range, the effect of parameters of orders  $N \geq 2$  on the energy per nucleon of NM is very small, and consequently we can legitimately expect that  $Q_{\text{sat}}$ ,  $Z_{\text{sat}}$ ,  $Q_{\text{sym}}$ , and  $Z_{\text{sym}}$  will not be strongly constrained by the LD filter. In that sense, for each parameter set passing through the pass-band filter  $w_{LD}$ , we sample 10 new parameter sets by randomly drawing new values for the high-order parameters with the aim of obtaining more sets passing through both LD and HD filters. Nevertheless, we have seen in Fig. 1.12 that the inclusion of orders  $N \geq 2$  is important to precisely estimate the CC transition point. In that sense,  $n_t$  and  $P_t$  will be calculated using the high-order parameters that originate from the passage through the LD filter and which are kept by the HD filter.

The possibility of a negative symmetry energy at high density was sometimes evoked in the literature [Li17; WFF88]. However, extremely soft functionals are generally incompatible with the maximum mass constraint, thus relaxing the condition of the positiveness of the symmetry energy will not alter our results.

Finally, let us notice that we do not include the constraint on the tidal deformability inferred from the GW170817 event [Abb+17] in the  $w_{HD}$  filter, the reason being that we are interested in confronting our results with the observations of the LIGO-Virgo collaboration.

#### 2.2.4 Posterior distribution of empirical parameters

### 2.3 General predictions for neutron star observables

#### 2.3.1 Confrontation with GW170817 event

##### 2.3.1.1 Equation of state

##### 2.3.1.2 Mass-radius relation

##### 2.3.1.3 Tidal deformability

#### 2.3.2 Bayesian analysis of the crust-core transition

#### 2.3.3 Crustal moment of inertia and pulsar glitches

### 2.4 Conclusion



## Appendix A

# Energy density of a relativistic electron gas

We give here the derivation of the energy density of a relativistic electron gas of density  $n_e$  at zero temperature.

From  $\approx 10^{14}$  g/cm<sup>3</sup>, electrons are essentially free. In this case, the energy density is given by

$$\varepsilon_e = \frac{c}{\pi^2} \int_{k=0}^{k_e} dk k^2 \sqrt{\hbar^2 k^2 + m_e^2 c^2}, \quad (\text{A.1})$$

$c$  being the speed of light,  $\hbar$  the reduced Planck constant, and  $m_e$  the electron mass. The electron Fermi wave number  $k_e$  is related to the electron density via  $k_e = (3\pi^2 n_e)^{1/3}$ . Making a simple variable,  $x = \hbar k / (m_e c)$ , we can rewrite the energy density as

$$\varepsilon_e = \frac{m_e^4 c^5}{\pi^2 \hbar^3} \xi(x_r) \quad (\text{A.2})$$

with  $x_r = \hbar k_e / (m_e c)$  and

$$\xi = \int_{x=0}^{x_r} dx x^2 \sqrt{x^2 + 1}. \quad (\text{A.3})$$

Integrating by parts, it follows

$$\begin{aligned} \xi &= \left[ \frac{x^3}{3} \sqrt{x^2 + 1} \right]_0^{x_r} - \int_0^{x_r} dx \frac{x^4}{3\sqrt{x^2 + 1}} \\ 3\xi &= x_r^3 \sqrt{x_r^2 + 1} - \int_0^{x_r} dx x^2 \left( \frac{x^2 + 1}{\sqrt{x^2 + 1}} - \frac{1}{\sqrt{x^2 + 1}} \right) \\ 4\xi &= x_r^3 \sqrt{x_r^2 + 1} + \int_0^{x_r} dx \frac{x^2}{\sqrt{x^2 + 1}}. \end{aligned} \quad (\text{A.4})$$

We have

$$\frac{d}{dx} \left( x \sqrt{x^2 + 1} \right) = \sqrt{x^2 + 1} + \frac{x^2}{\sqrt{x^2 + 1}}, \quad (\text{A.5})$$

and

$$\frac{d}{dx} \left[ \ln(x + \sqrt{x^2 + 1}) \right] = \sqrt{x^2 + 1} - \frac{x^2}{\sqrt{x^2 + 1}}. \quad (\text{A.6})$$

thus,

$$\frac{1}{2} \frac{d}{dx} \left[ x \sqrt{x^2 + 1} - \ln \left( x + \sqrt{x^2 + 1} \right) \right] = \frac{x^2}{\sqrt{x^2 + 1}}. \quad (\text{A.7})$$

Hence,

$$4\xi = x_r^3 \sqrt{x_r^2 + 1} + \frac{1}{2} \left[ x \sqrt{x^2 + 1} - \ln \left( x + \sqrt{x^2 + 1} \right) \right]_0^{x_r}, \quad (\text{A.8})$$

which finally gives

$$\varepsilon_e(n_e) = \frac{m_e^4 c^5}{8\pi^2 \hbar^3} \left[ x_r (2x_r^2 + 1) \gamma_r - \ln(x_r + \gamma_r) \right], \quad (\text{A.9})$$

where we have introduced the parameter  $\gamma_r = \sqrt{x_r^2 + 1}$ .

## Appendix B

# Neutron and proton chemical potentials in the metamodel

From the metamodeling of the nuclear matter energy, Eq. 1.37, we give here the complete expressions of the neutron and proton chemical potentials for a homogeneous system characterized by a neutron density  $n_n$  and proton density  $n_p$  at zero temperature. Let us introduce the variables  $x = (n - n_{sat})/(3n_{sat})$  and  $\delta = (n_n - n_p)/n$ ,  $n = n_n + n_p$  being the total baryon density.

The neutron chemical potential, with the rest mass energy  $m_n c^2$ , is given by

$$\mu_{HM,n} = e_{HM}(n_n, n_p) + n \left( \frac{\partial e_{HM}}{\partial n_n} \right)_{n_p} + m_n c^2 \quad (\text{B.1})$$

$$\mu_{HM,n} = e_{HM}(n, \delta) + \frac{1+3x}{3} \left( \frac{\partial e_{HM}}{\partial x} \right)_{\delta} + (1-\delta) \left( \frac{\partial e_{HM}}{\partial \delta} \right)_x + m_n c^2. \quad (\text{B.2})$$

Equivalently, the proton chemical potential is given by

$$\mu_{HM,p} = e_{HM}(n, \delta) + \frac{1+3x}{3} \left( \frac{\partial e_{HM}}{\partial x} \right)_{\delta} - (1+\delta) \left( \frac{\partial e_{HM}}{\partial \delta} \right)_x + m_p c^2. \quad (\text{B.3})$$

We first go through the derivative with respect to  $x$ , yielding

$$\begin{aligned} \left( \frac{\partial e_{HM}}{\partial x} \right)_{\delta} &= \frac{5}{1+3x} t_{HM}^{FG}(n, \delta) \\ &\quad - \frac{3}{1+3x} \frac{1}{2} t_{sat}^{FG} (1+3x)^{2/3} \left[ (1+\delta)^{5/3} + (1-\delta)^{5/3} \right] \\ &\quad + \sum_{\alpha \geq 0} (v_{\alpha}^{is} + v_{\alpha}^{iv} \delta^2) \frac{1}{\alpha!} \left( \alpha x^{\alpha-1} u_{\alpha}(x) + x^{\alpha} \frac{du_{\alpha}}{dx} \right). \end{aligned} \quad (\text{B.4})$$

Let us notice that it gives the expression for the total pressure,

$$P_{HM}(n, \delta) = n^2 \left( \frac{\partial e_{HM}}{\partial n} \right)_{\delta} \quad (\text{B.5})$$

$$= \frac{1}{3} n_{sat} (1+3x)^2 \left( \frac{\partial e_{HM}}{\partial x} \right)_{\delta}. \quad (\text{B.6})$$

We now turn to the derivative with respect to the asymmetry  $\delta$ ,

$$\begin{aligned}
\left(\frac{\partial e_{HM}}{\partial \delta}\right)_x &= \frac{1}{2}t_{sat}^{FG}(1+3x)^{5/3} \left[(1+\delta)^{5/3} - (1-\delta)^{5/3}\right] \\
&\quad + \frac{5}{6}t_{sat}^{FG}(1+3x)^{2/3} \left[(1+\delta)^{2/3} \frac{m}{m_n^*} - (1-\delta)^{2/3} \frac{m}{m_p^*}\right] \\
&\quad + 2\delta \sum_{\alpha \geq 0} v_\alpha^{iv} \frac{x^\alpha}{\alpha!} u_\alpha(x).
\end{aligned} \tag{B.7}$$

## Appendix C

### Total pressure in the inner crust



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