Mean-field approach to nuclear structure with semi-realistic nucleon-nucleon interactions

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Semi-realistic nucleon-nucleon interactions applicable to the self-consistent mean-field (both Hartree-Fock and Hartree-Fock-Bogolyubov) calculations are developed, by modifying the M3Y interaction. The modification is made so as to reproduce binding energies and rms matter radii of doubly magic nuclei, single-particle levels in $^{208}$Pb, and even-odd mass differences of the Sn isotopes. We find parameter-sets with and without the tensor force. The new interactions are further checked by the saturation properties of the uniform nuclear matter including the Landau-Migdal parameters. By the mean-field calculations, interaction-dependence of the neutron drip line is investigated for the O, Ca and Ni isotopes, and of the single-particle energies for the $N = 16, 32, 50, 82$ and $Z = 50$ nuclei. Results of the semi-realistic interactions including the tensor force are in fair agreement with available experimental data for all of these properties.

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I. INTRODUCTION

Mean-field (MF) theories provide us with a good first approximation to the nuclear structure problems. They are able to describe the saturation and the shell structure simultaneously, both of which are basic nuclear properties, based on effective nucleon-nucleon ($NN$) interactions. As far as we constrain to the non-relativistic approaches, most of the MF calculations have been performed with the Skyrme interaction $^{[1]}$. Finite-range interactions have rarely been applied, except the Gogny interaction $^{[2]}$ which has the Gaussian form for the central channels. Most popular parameter-sets of the Skyrme and the Gogny interactions have been adjusted mainly to the nuclear properties around the $\beta$-stability. However, it is a question whether such phenomenological effective interactions work well for nuclei far off the $\beta$-stability. For instance, whereas role of the tensor force in nucleus-dependence of the MF has attracted interest $^{[3, 4]}$, the tensor force is usually ignored in those parameter-sets.

Although direct application of the bare $NN$ (and $NNN$) interaction to the nuclear structure problems $^{[5, 6, 7]}$ is yet limited to light nuclei or made to medium-mass nuclei but with limited accuracy, guide from microscopic theories will be valuable even in heavy-mass nuclei. The Michigan 3-range Yukawa (M3Y) interaction $^{[8]}$, which was derived by fitting the Yukawa functions to Brueckner's $G$-matrix, has been used in nuclear structure as well as in low-energy nuclear reaction studies. There have been a few attempts applying the M3Y-type interaction to MF calculations $^{[9, 10]}$. In Ref. $^{[10]}$, the author has developed an M3Y-type interaction which is applicable to the Hartree-Fock (HF) calculations. The original M3Y interaction is incapable of reproducing the saturation and the spin-orbit splitting within the MF regime. To cure this problem a density-dependent contact term has been added and some of the strength parameters have been modified. Such interactions, which have originally been derived from microscopic theories but are slightly modified from phenomenological standpoints, may be called semi-realistic interactions. It has been shown $^{[10, 11]}$ that semi-realistic $NN$ interactions could give different shell structure from the widely used Skyrme and Gogny interactions. However, the pairing properties have not been taken into account in the parameter-set M3Y-P2 that was proposed in Ref. $^{[10]}$. This implies that the singlet-even channel in M3Y-P2 is not quite appropriate as long as the pairing interaction is taken to be consistent with the HF interaction, while this problem seems to be masked in the HF approximation. In order to apply semi-realistic interactions to the MF studies extensively, we explore new parameter-sets of the M3Y-type interaction, taking the pairing properties into account. Special attention is paid also to role of the tensor force. Using the recently developed algorithm $^{[12, 13, 14]}$, we apply the new semi-realistic interactions to the Hartree-Fock-Bogolyubov (HFB) as well as to the HF calculations of medium- to heavy-mass spherical nuclei.

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II. M3Y-TYPE INTERACTION

We express a non-relativistic nuclear effective Hamiltonian by

\begin{equation}
H_N = K + V_N; \quad K = \sum_i \frac{p_i^2}{2m}, \quad V_N = \sum_{i<j} v_{ij},
\end{equation}

with \( i \) and \( j \) representing the indices of individual nucleons. For the effective \( NN \) interaction \( v_{ij} \), we consider the following form,

\begin{align}
v_{ij} &= v_{ij}^{(C)} + v_{ij}^{(LS)} + v_{ij}^{(TN)} + v_{ij}^{(DD)}; \\
v_{ij}^{(C)} &= \sum_{n=1}^2 (t_n^{(SE)} P_{SE} + t_n^{(TE)} P_{TE} + t_n^{(SO)} P_{SO} + t_n^{(TO)} P_{TO}) f_n^{(C)} (r_{ij}), \\
v_{ij}^{(LS)} &= \sum_{n=1}^2 (t_n^{(LS)} P_{TE} + t_n^{(LSO)} P_{TO}) f_n^{(LS)} (r_{ij}) \mathbf{L}_{ij} \cdot (\mathbf{s}_i + \mathbf{s}_j), \\
v_{ij}^{(TN)} &= \sum_{n=1}^2 (t_n^{(TNE)} P_{TE} + t_n^{(TNO)} P_{TO}) f_n^{(TN)} (r_{ij}) r_{ij}^2 S_{ij}, \\
v_{ij}^{(DD)} &= (t_p^{(SE)} P_{SE} \cdot [\rho(r_i)]^{(SE)} + t_p^{(TE)} P_{TE} \cdot [\rho(r_i)]^{(TE)} ) \delta(r_{ij}).
\end{align}

where \( r_{ij} = r_i - r_j, r_{ij} = |r_{ij}|, p_{ij} = (p_i - p_j)/2, L_{ij} = r_{ij} \times p_{ij}, S_{ij} = 4 [3(s_i \cdot r_{ij})(s_j \cdot r_{ij}) - s_i \cdot s_j] \) with \( \hat{r}_{ij} = r_{ij}/r_{ij} \), and \( \rho(r) \) denotes the nucleon density. The Yukawa function \( f_n(r) = e^{-\mu_n r}/\mu_n r \) is assumed for all channels except \( v^{(DD)} \) in the M3Y-type interactions. The projection operators on the singlet-even (SE), triplet-even (TE), singlet-odd (SO) and triplet-odd (TO) two-particle states are defined as

\begin{align}
P_{SE} &= \frac{1 - P_\sigma}{2} \frac{1 + P_\tau}{2}, & P_{TE} &= \frac{1 + P_\sigma}{2} \frac{1 - P_\tau}{2}, \\
P_{SO} &= \frac{1 - P_\sigma}{2} \frac{1 - P_\tau}{2}, & P_{TO} &= \frac{1 + P_\sigma}{2} \frac{1 + P_\tau}{2},
\end{align}

where \( P_\sigma \) (\( P_\tau \)) expresses the spin (isospin) exchange operator.

We shall start from the M3Y-Paris interaction \[13\], which will be denoted by M3Y-P0 in this article as in Ref. \[10\]. We change none of the range parameters \( \mu_n \) of M3Y-P0 in \( v^{(C)}, v^{(LS)} \) and \( v^{(TN)} \). In M3Y-P0, the longest range part in \( v^{(C)} \) is kept identical to the central channels of the one-pion exchange potential (OPEP), \( v^{(C)}_{OPEP} \). We also maintain this reasonable assumption. As is well known, the spin-orbit (\( \ell s \)) splitting plays a significant role in the nuclear shell structure. Even though higher-order effects may account for the observed \( \ell s \) splitting \[16\], it is desired to enhance \( v^{(LS)} \) in order to describe the shell structure within the MF regime. We here use an overall enhancement factor to \( v^{(LS)} \), which is determined from the single-particle (s.p.) spectrum of \(^{208}\text{Pb} \), as will be shown in Sec. \[IV\].

Influence of the tensor force on s.p. energies is a current topic, which could be relevant to the new magic numbers in \(^{208}\text{Pb} \). We develop two parameter-sets having \( v^{(TN)} \) without any modification from M3Y-P0, as well as a parameter-set in which we impose \( v^{(TN)} = 0 \).

The saturation properties are important to describe many nuclei in a wide mass range. Since it is still hard to describe accurately the saturation properties by the bare \( NN \) (and \( NNN \)) interaction despite certain progress \[1\], it will be appropriate to modify realistic effective interaction so as to reproduce the saturation properties. Density-dependence in the effective interaction has been known to be essential in obtaining the saturation. We therefore add a density-dependent contact force \( v^{(DD)} \) \[10\]. The parameter \( \alpha^{(TE)} \) in \( v^{(DD)} \), power to \( \rho \), is taken to be 1/3, by which the incompressibility \( K \) becomes close to a reasonable value as shown later. On the other hand, \( \alpha^{(SE)} \) is not quite sensitive to \( K \), because the major source of the saturation lies in the TE channel, not in the SE channel \[13\]. Although we simply assumed \( \alpha^{(SE)} = \alpha^{(TE)} = 1/3 \) in M3Y-P2 \[10\], this assumption makes it difficult to reproduce pairing properties and to avoid instability of the neutron matter \[10,12\] simultaneously. To overcome this problem, we adopt \( \alpha^{(SE)} = 1 \) in the new parameter-sets. The difference between \( \alpha^{(SE)} \) and \( \alpha^{(TE)} \) may be attributed to the difference in origin of the \( \rho \)-dependence; the short-range repulsion in the bare \( NN \) interaction in the SE channel while primarily the tensor force in the TE channel. Validity of the choice \( \alpha^{(SE)} = 1 \) is further discussed in Sec. \[III\].

The remaining parameters are \( t_n \) in \( v^{(C)} \) (except those of \( v^{(OPEP)} \)) and \( t_p \) in \( v^{(DD)} \). We fit them to the measured binding energies of \(^{16}\text{O} \) and \(^{208}\text{Pb} \), in the HF approximation (see Sec. \[IX\]). The proton and neutron Fermi energies of \(^{208}\text{Pb} \), which are primarily relevant to the symmetry energy, are checked additionally. To determine \( t_n^{(SE)} \) and \( t_p^{(SE)} \), we also use the even-odd mass differences of the Sn isotopes, by comparing results of the Hartree-Fock-Bogolyubov
In the set M3Y-P5, we keep both $v^{(\text{TN})}$ and the odd-channel (SO and TO) strengths in $v^{(\text{C})}$ of M3Y-P0. The set M3Y-P4 is obtained by assuming $v^{(\text{TN})} = 0$, while changing $t_n^{(\text{SO})}$ and $t_n^{(\text{TO})}$ ($n = 1, 2$) substantially. In the set M3Y-P5, we somewhat modify $t_2^{(\text{SO})}$ and $t_2^{(\text{TO})}$ while keeping $v^{(\text{TN})}$, so as to reproduce the binding energies of several doubly magic nuclei better than M3Y-P3, as will be shown in Sec. IV. Thus the number of adjusted parameters are 7, 11 and 9 (including the overall enhancement factor to $v^{(\text{LS})}$) for M3Y-P3, P4 and P5, respectively. It will be useful to compare results of these parameter-sets for pinning down which part of the interaction is important to individual physical quantities. In particular, role of the tensor force will be of interest. It is remarked that, while schematic tensor forces have been introduced into some of the recent MF studies [4, 18], the present $v^{(\text{TN})}$ in M3Y-P3 and P5 is much more realistic.

| parameters | M3Y-P0 | M3Y-P2 | M3Y-P3 | M3Y-P4 | M3Y-P5 |
|------------|--------|--------|--------|--------|--------|
| $1/\mu_1^{(\text{C})}$ (fm) | 0.25   | 0.25   | 0.25   | 0.25   | 0.25   |
| $t_1^{(\text{SE})}$ (MeV) | 11466. | 8027.  | 8027.  | 8027.  | 8027.  |
| $t_1^{(\text{TE})}$ (MeV) | 13967. | 6080.  | 7130.  | 5503.  | 5576.  |
| $t_1^{(\text{SO})}$ (MeV) | -1418. | -11900.| -1418. | -12000.| -1418. |
| $t_1^{(\text{TO})}$ (MeV) | 11345. | 3800.  | 11345. | 3700.  | 11345. |
| $1/\mu_1^{(\text{C})}$ (fm) | 0.40   | 0.40   | 0.40   | 0.40   | 0.40   |
| $t_2^{(\text{SE})}$ (MeV) | -3556. | -2880. | -2637. | -2637. | -2650. |
| $t_2^{(\text{TE})}$ (MeV) | -4594. | -4266. | -4594. | -4183. | -4170. |
| $t_2^{(\text{SO})}$ (MeV) | 950.   | 2730.  | 950.   | 4500.  | 2880.  |
| $t_2^{(\text{TO})}$ (MeV) | -1900. | -780.  | -1900. | -1000. | -1780. |
| $1/\mu_1^{(\text{LS})}$ (fm) | 0.25   | 0.25   | 0.25   | 0.25   | 0.25   |
| $t_1^{(\text{LS})}$ (MeV) | -5101. | -9181.8| -10712.1| -8671.7| -11222.2|
| $t_2^{(\text{LS})}$ (fm) | 0.40   | 0.40   | 0.40   | 0.40   | 0.40   |
| $t_1^{(\text{LO})}$ (MeV) | -1897. | -3414.6| -3983.7| -3224.9| -4173.4|
| $t_2^{(\text{LO})}$ (MeV) | -337.  | -606.6 | -707.7 | -572.9 | -741.4 |
| $t_1^{(\text{L})}$ (fm) | 0.40   | 0.40   | 0.40   | 0.40   | 0.40   |
| $t_2^{(\text{L})}$ (fm) | 0.40   | 0.40   | 0.40   | 0.40   | 0.40   |
| $t_1^{(\text{LNO})}$ (MeV) | 244.   | 29.28  | 244.   | 0.     | 244.   |
| $t_2^{(\text{LNO})}$ (MeV) | 15.6   | 15.6   | 15.6   | 0.     | 15.6   |
| $\alpha_1^{(\text{SE})}$ (fm) | -707.7 | -30.9  | -3.708 | 1/3    | 1      |
| $\alpha_2^{(\text{SE})}$ (MeV fm$^3$) | 0.     | 181.  a) | 220.   | 248.   | 126.   |
| $\alpha_1^{(\text{TE})}$ (fm) | 0.     | 1/3    | 1/3    | 1/3    | 1/3    |
| $\alpha_2^{(\text{TE})}$ (MeV fm) | 0.     | 1139.  | 1198.  | 1142.  | 1147.  |

a) (MeV·fm)
III. PROPERTIES OF NUCLEAR MATTER AT AND AROUND SATURATION POINT

We first view properties of the infinite nuclear matter that are predicted by the semi-realistic $NN$ interactions. In the HF approximation, energy of the nuclear matter can be expressed by the following variables:

\[
\begin{align*}
\rho &= \sum_{\sigma} \rho_{\tau\sigma}, \\
\eta_s &= \frac{\sum_{\sigma} \sigma \rho_{\tau\sigma}}{\rho} = \frac{\rho_p - \rho_p + \rho_n - \rho_n}{\rho}, \\
\eta_t &= \frac{\sum_{\sigma} \tau \rho_{\tau\sigma}}{\rho} = \frac{\rho_p + \rho_p - \rho_n - \rho_n}{\rho}, \\
\eta_{st} &= \frac{\sum_{\sigma} \sigma \tau \rho_{\tau\sigma}}{\rho} = \frac{\rho_p - \rho_p - \rho_n + \rho_n}{\rho}.
\end{align*}
\]

(4)

\(\rho_{\tau\sigma}\) (\(\tau = p, n\) and \(\sigma = \uparrow, \downarrow\), which are sometimes substituted by \(\pm 1\) without confusion) stands for densities depending on the spin and the isospin, and is related to the Fermi momentum \(k_{F\tau\sigma}\) by

\[
\rho_{\tau\sigma} = \frac{1}{6\pi^2} k_{F\tau\sigma}^3.
\]

(5)

In the interaction of Eq. (2), only \(v^{(C)} + v^{(DD)}\) contributes to the energy of the uniform nuclear matter. Formulas to calculate the nuclear matter energy and its derivatives for given \(k_{F\tau\sigma}\) have been derived in Ref. [10]. Note that, even when superfluidity makes the nuclear matter energy somewhat lower, it is not much different from the energy in the HF approximation.

The spin-saturated symmetric matter is characterized by \(\eta_s = \eta_t = \eta_{st} = 0\), for which we denote \(k_{F\tau\sigma}\) simply by \(k_F\). The minimum of the energy per nucleon \(E = E/A\), given by

\[
\left. \frac{\partial E}{\partial \rho} \right|_0 = \left. \frac{\partial E}{\partial k_F} \right|_0 = 0,
\]

(6)
defines the saturation density \(\rho_0\) (equivalently, \(k_{F0}\)) and energy \(E_0\). The expression \(|_0\) indicates evaluation at the saturation point. As well as \(\rho_0\) and \(E_0\), second derivatives of \(E\) carry basic information of the effective \(NN\) interaction. Two of the curvatures of \(E\) at the saturation point are called incompressibility and volume symmetry energy,

\[
K = k_F^2 \left. \frac{\partial^2 E}{\partial k_F^2} \right|_0 = 9 \rho^2 \left. \frac{\partial^2 E}{\partial \rho^2} \right|_0, \quad a_t = \frac{1}{2} \left. \frac{\partial^2 E}{\partial \eta_t^2} \right|_0,
\]

(7)

and are related to the Landau-Migdal (LM) parameters \(f_1\) and \(f_0\) as

\[
K = \frac{3k_{F0}^2}{M_0^*} (1 + f_0), \quad a_t = \frac{k_{F0}^2}{6M_0^*} (1 + f_0'),
\]

(8)

where \(M_0^*\) represents the effective mass (\(k\)-mass) at the saturation point. See Ref. [10] for definition of the LM parameters. The other curvatures of \(E\) with respect to \(\eta_s\) and \(\eta_{st}\), denoted by \(a_s\) and \(a_{st}\), are defined analogously to \(a_t\) and are expressed in terms of the LM parameters \(g_0\) and \(g_0'\) [10]. The \(k\)-mass is defined by a derivative of the s.p. energy \(\varepsilon(k\sigma\tau)\):

\[
\left. \frac{\partial \varepsilon(k\sigma\tau)}{\partial k} \right|_0 = \frac{k_{F0}}{M_0^*},
\]

(9)

and is connected to the LM parameter \(f_1\) by

\[
\frac{M_0^*}{M} = 1 + \frac{1}{3} f_1.
\]

(10)

In addition, density-dependence of the symmetry energy, which is represented by a third derivative of \(E\) as

\[
\mathcal{L}_t = \frac{1}{2} k_F \left. \frac{\partial^3 E}{\partial k_F \partial \eta_t^2} \right|_0 = \frac{3}{2} \rho \left. \frac{\partial^3 E}{\partial \rho \partial \eta_t^2} \right|_0,
\]

(11)
is under interest in relevance to structure of the neutron star crust \[19\]. These quantities calculated from the new semi-realistic interactions are tabulated in Tables \[\text{III}\] We here set \(M = (M_p + M_n)/2\), where \(M_p\) (\(M_n\)) is the measured mass of a proton (a neutron) \[20\]. For comparison, the values obtained by the D1S parameter-set \[21\] of the Gogny interaction and from M3Y-P2 are also displayed.

Related to the global systematics of the binding energies and the radii, \(k_{F0} \approx 1.33 - 1.34 \text{ fm}^{-1}\) and \(E_0 \approx -16\) MeV have been established empirically. Although the M3Y-P3 interaction yields deeper \(E_0\) than the other interactions, it is still within the range of ambiguity in extracting the volume energy from the experimental data \[22\]. In practice, M3Y-P3 does not yield overbinding for any of the doubly magic nuclei presented in Sec. \[\text{IV}\].

Global characters of the spin and isospin responses are customarily discussed in terms of the LM parameters. By using the formulas given in Ref. \[10\], we evaluate the LM parameters for the new semi-realistic interactions, as shown in Table \[\text{III}\].

Table II: Nuclear matter properties at the saturation point.

|                  | D1S | M3Y-P2 | M3Y-P3 | M3Y-P4 | M3Y-P5 |
|------------------|-----|--------|--------|--------|--------|
| \(k_{F0}\) (fm) | 1.342 | 1.340 | 1.340 | 1.340 | 1.340 |
| \(E_0\) (MeV)   | -16.01 | -16.14 | -16.51 | -16.13 | -16.12 |
| \(K\) (MeV)     | 202.9 | 220.4 | 245.8 | 235.3 | 235.6 |
| \(M_0^*/M\)     | 0.697 | 0.652 | 0.658 | 0.665 | 0.629 |
| \(a_t\) (MeV)   | 31.12 | 30.61 | 29.75 | 28.71 | 29.59 |
| \(a_s\) (MeV)   | 26.18 | 21.19 | 20.17 | 15.61 | 19.56 |
| \(a_{st}\) (MeV) | 29.13 | 38.19 | 36.45 | 39.89 | 41.01 |
| \(\mathcal{L}_t\) (MeV) | 22.44 | 27.98 | 25.30 | 17.87 | 24.63 |

In Fig. 2, contributions of the SE, TE, SO and TO channels in \[v^{(C)} + v^{(DD)}\] to \(E\) of the symmetric matter are shown as a function of \(k_F\). The contribution of the TE and the SO channels in M3Y-P4 is hard to be distinguished from that in M3Y-P5. So is the contribution of the TE channel in M3Y-P2. The TE channel takes a minimum at \(k_F \approx 1.5\) fm\(^{-1}\), primarily responsible for the saturation at \(k_F = k_{F0} \approx 1.3\) fm\(^{-1}\). Both the SO and the TO channels do not contribute to \(E\) significantly at \(\rho \lesssim \rho_0\) (i.e. \(k_F \lesssim k_{F0}\)). While the SO channel becomes attractive in the D1S interaction, it is repulsive in the M3Y-type interactions at \(\rho > \rho_0\). The TO channel is repulsive in M3Y-P3 and P5, while attractive in M3Y-P4, although the attraction in M3Y-P4 is not so strong as to cause spin polarization in the pure neutron matter up to \(\rho = 1.3\) fm\(^{-1}\)(\(\approx 8\rho_0\)). Remember that the odd channels in M3Y-P3 are unchanged from M3Y-P0.

Energy per nucleon in the spin-saturated neutron matter (i.e. \(\eta_s = -1\)) is presented in Fig. 4. The result from a microscopic calculation in Ref. \[27\] is also shown as a reference. The unphysical behavior at high \(\rho\) in the D1S result, which comes from the absence of density-dependence in the SE channel, was pointed out in Refs. \[10, 12\]. The present M3Y-P3 to P5 interactions have relatively strong \(\rho\)-dependence at high \(\rho\) for the neutron matter, if compared to M3Y-P2 and D1S. This originates in \(\alpha^{(SE)}(=1)\), and tends to make \(E(\rho)\) closer to the result of Ref. \[27\].

IV. PROPERTIES OF DOUBLY MAGIC NUCLEI

We next turn to doubly magic nuclei, for which the spherical HF approach is expected to be a good approximation.
with \( \nu \) expansion method (GEM) \[12, 13\]. In this method we employ the s.p. bases of apart from the isospin index. Here \( Y \) parameter \( \nu \) parameter of a nuclide, we adopt the following basis parameters \[14\]:

\[
\text{M3Y-P3 yields underbinding by about 9 MeV (5 MeV) accuracy. Though the accuracy is slightly worse, the new interactions also reproduce experimental data. Influence of the c.m. motion on the matter radii is subtracted in a similar manner to the c.m.}
\]

\[0.743 \quad 0.607 \quad 0.578 \quad 0.538 \quad 0.502\]

\[0.470 \quad 0.635 \quad 0.670 \quad 0.797 \quad 0.692\]

\[0.342 \quad 0.245 \quad 0.271 \quad 0.286 \quad 0.267\]

\[0.100 \quad 0.096 \quad 0.104 \quad 0.106 \quad 0.100\]

\[0.466 \quad 0.113 \quad 0.070 \quad -0.164 \quad -0.007\]

\[-0.184 \quad 0.273 \quad 0.214 \quad 0.374 \quad 0.299\]

\[0.245 \quad 0.162 \quad 0.160 \quad 0.190 \quad 0.178\]

\[0.091 \quad 0.078 \quad 0.079 \quad 0.085 \quad 0.081\]

\[0.631 \quad 1.006 \quad 0.933 \quad 1.136 \quad 1.081\]

\[0.610 \quad 0.202 \quad 0.213 \quad 0.109 \quad 0.087\]

\[-0.038 \quad 0.040 \quad 0.063 \quad 0.016 \quad 0.029\]

\[-0.036 \quad -0.002 \quad 0.005 \quad -0.008 \quad -0.002\]

To all the following calculations of finite nuclei, we apply the recently developed algorithm based on the Gaussian expansion method (GEM) \[12, 13\]. In this method we employ the s.p. bases of

\[
\varphi_{\nu\ell j m}(r) = R_{\nu\ell j}(r)|Y^{(\ell)}(\hat{r})\chi_\nu(r)|^m; \quad R_{\nu\ell j}(r) = N_{\nu\ell j} r^\ell \exp(-\nu r^2),
\]

(12)

apart from the isospin index. Here \( Y^{(\ell)}(\hat{r}) \) expresses the spherical harmonics and \( \chi_\nu \) the spin wave function. The parameter \( \nu = \nu_r + i\nu_i \) indicates a complex number corresponding to the range of the Gaussian. Irrespective to nuclide, we adopt the following basis parameters \[14\]:

\[
\nu_r = \nu_0 b^{-2n}, \quad \nu_i = \begin{cases} 
0 & (n = 0, 1, \ldots, 5) \\
\pm \frac{\pi}{2} & (n = 0, 1, 2) 
\end{cases}
\]

(13)

with \( \nu_0 = (2.40 \text{fm})^{-2} \) and \( b = 1.25 \) for each \( (\ell, j) \). It is notable that, without parameters specific to mass number or nuclide, a single set of the GEM bases is applicable to wide range of the nuclear mass table \[14\]. The Hamiltonian is \( H = H_N + V_C - H_{c.m.} \), where \( V_C \) and \( H_{c.m.} \) represent the Coulomb interaction and the center-of-mass (c.m.) Hamiltonian, while \( H_N \) has been given in Eq. (1). The exchange term of \( V_C \) is treated exactly, in the same manner as the nuclear force \( V_N \). Both the one- and the two-body terms of \( H_{c.m.} \) are subtracted before iteration.

The calculated binding energies and rms matter radii of several doubly magic nuclei are displayed in Table IV. The results of the new semi-realistic interactions are compared with those of D1S and M3Y-P2 as well as with the experimental data. Influence of the c.m. motion on the matter radii is subtracted in a similar manner to the c.m. energies \[10\]. The binding energies of these nuclei obtained from D1S (M3Y-P2) are in agreement with the measured values within the 3 MeV (5 MeV) accuracy. Though the accuracy is slightly worse, the new interactions also reproduce the binding energies moderately well. M3Y-P3 yields underbinding by about \( 9 - 17 \text{ MeV} \) except for \( ^{208}\text{Pb} \). For M3Y-P4 and P5, maximum deviation in the binding energies shown in Table IV is \( \sim 7 \text{ MeV} \). Since correlations due to the residual interaction could influence, we do not take this deviation seriously at the present stage. The rms matter radii of these nuclei calculated from the semi-realistic interactions are comparable to those from the D1S interaction, in fair agreement with the data.

The s.p. levels in \( ^{208}\text{Pb} \) are depicted in Fig. 3. The levels obtained from M3Y-P5 are compared with those from D1S and the experimental levels. M3Y-P3 and P4 give s.p. levels similar to, though slightly different from, those of M3Y-P5. The experimental s.p. energies are obtained from the levels of the neighboring nuclei: \( ^{207,209}\text{Pb}, ^{207,210}\text{Tl} \) and \( ^{209}\text{Bi} \). In the HF results, the overall level spacing is relevant to \( M_0^\nu \) shown in Table II. In the usual HF calculations we have larger level spacing than in the data, and it is not (should not be) remedied until correlations due to the residual interaction (or the \( \omega \)-mass) are taken into account \[24\]. We thus confirm that the present interactions yield as plausible s.p. levels as D1S does.
FIG. 1: Energy per nucleon $E = E/A$ in the symmetric nuclear matter for several effective interactions. Purple dashed, green dot-dashed and red solid lines represent the results with the M3Y-P3, P4 and P5 interactions, respectively. Those with the M3Y-P2 and D1S interactions are also displayed for comparison by orange and blue solid lines.

TABLE IV: Binding energies and rms matter radii of several doubly magic nuclei. Experimental data are taken from Refs. [28, 29, 30, 31].

|       | Exp. | D1S  | M3Y-P2 | M3Y-P3 | M3Y-P4 | M3Y-P5 |
|-------|------|------|--------|--------|--------|--------|
| $^{16}$O | $-E$ (MeV) | 127.6 | 129.5 | 127.2 | 118.6 | 126.3 | 126.1 |
|       | $\sqrt{\langle r^2 \rangle}$ (fm) | 2.61 | 2.61 | 2.61 | 2.65 | 2.60 | 2.59 |
| $^{24}$O | $-E$ (MeV) | 168.5 | 168.6 | 165.7 | 158.2 | 164.0 | 166.7 |
|       | $\sqrt{\langle r^2 \rangle}$ (fm) | 3.19 | 3.01 | 3.06 | 3.08 | 3.04 | 3.03 |
| $^{40}$Ca | $-E$ (MeV) | 342.1 | 344.6 | 338.8 | 325.2 | 337.0 | 335.1 |
|       | $\sqrt{\langle r^2 \rangle}$ (fm) | 3.47 | 3.37 | 3.38 | 3.42 | 3.37 | 3.37 |
| $^{48}$Ca | $-E$ (MeV) | 416.0 | 416.8 | 411.9 | 401.0 | 409.4 | 414.1 |
|       | $\sqrt{\langle r^2 \rangle}$ (fm) | 3.57 | 3.51 | 3.53 | 3.56 | 3.52 | 3.50 |
| $^{90}$Zr | $-E$ (MeV) | 783.9 | 785.9 | 779.4 | 767.9 | 775.1 | 779.8 |
|       | $\sqrt{\langle r^2 \rangle}$ (fm) | 4.32 | 4.24 | 4.25 | 4.27 | 4.24 | 4.23 |
| $^{132}$Sn | $-E$ (MeV) | 1102.9 | 1104.1 | 1099.0 | 1089.3 | 1095.7 | 1098.4 |
|       | $\sqrt{\langle r^2 \rangle}$ (fm) | 4.77 | 4.79 | 4.81 | 4.77 | 4.77 | 4.76 |
| $^{208}$Pb | $-E$ (MeV) | 1636.4 | 1639.0 | 1636.5 | 1635.2 | 1632.1 | 1633.2 |
|       | $\sqrt{\langle r^2 \rangle}$ (fm) | 5.49 | 5.51 | 5.54 | 5.55 | 5.51 | 5.51 |

In Ref. [4], it has been shown that the $Z = N = 20$ shell gaps are narrowed by the tensor force. It is also true in the M3Y-type interactions. The s.p. energy difference $\varepsilon_\tau(0f_{7/2}) - \varepsilon_\tau(0d_{3/2})$ ($\tau = p, n$) in $^{40}$Ca obtained by the D1S interaction is in good agreement with the experimental values, both for protons and neutrons, as viewed in Fig. 5. While M3Y-P2 and P4 give almost the same size of the shell gaps as D1S does, we have narrower gaps in the HF calculations with M3Y-P3 and P5. However, the shell gaps do not collapse by $\nu^{(TN)}$ of M3Y, in contrast to the zero-range tensor force of Ref. [4]. We still have 5.2 MeV (7.7 MeV) gap for the proton (neutron) orbits with M3Y-P5. These gaps are close to those obtained from the tensor-free Skyrme interaction ‘T22’ in Ref. [4]. It is also worth
FIG. 2: Contribution of the SE, TE, SO and TO channels to $\mathcal{E}$. See Fig. 1 for conventions.

FIG. 3: Energy per nucleon $\mathcal{E} = E/A$ in the neutron matter for several effective interactions. Circles are the results of Ref. [27]. See Fig. 1 for the other conventions.
FIG. 4: Single-particle energies for $^{208}\text{Pb}$. Experimental values are extracted from Refs. [31, 32].

commenting that, for M3Y-P5, the octupole correlations significantly influence the ground state of $^{40}\text{Ca}$, as will be discussed elsewhere. This can make the shell gap look wider, having possibility to account for the observed gap.

V. PAIRING PROPERTIES

The M3Y-P2 interaction seems to have reasonable characters in the HF regime, as exemplified in Table IV. However, M3Y-P2 has too strong pair correlations, indicating too strong attraction in the SE channel at low densities though almost invisible in Fig. 2. This character is inherited from the original M3Y interaction. We have developed the M3Y-P3 to P5 parameter-sets by taking the pairing properties into consideration. In this section we shall show characters of the new interactions with respect to the pairing. We restrict ourselves to the pairing among like-nucleons, as usual.

We implement the spherical HFB calculations for finite nuclei, using the GEM bases of Eqs. (12, 13) together with the $\ell \leq 7$ truncation. The blocked HFB calculations are applied to the odd-mass nuclei, by assuming that a quasiparticle (q.p.) occupies a specified spherical orbital. When several q.p. levels lie closely in energy, we compare the total energies by filling each q.p. level and adopt the lowest-energy solution.

A. Even-odd mass difference in Sn isotopes

The $t_n^{(\text{SE})}$ ($n = 1, 2$) and $t_p^{(\text{SE})}$ parameters of M3Y-P3 to P5 are adjusted to the even-odd mass differences of the Sn isotopes with $66 < N < 80$. For the mass difference we use the three-point formula $\Delta_{\text{mass}}^Z(N) = E(Z, N) - \frac{1}{2}[E(Z, N + 1) + E(Z, N - 1)]$, with $Z = 50$ and $N = \text{odd}$. The mass differences calculated with M3Y-P4 and P5 are displayed in Fig. 6 in comparison with the experimental data and with those of D1S. Though not shown to keep the figure viewable, M3Y-P3 gives similar $\Delta_{\text{mass}}^Z(N)$ to M3Y-P4. The calculations are not fully convergent for the $\ell$ truncation. Moreover, the restoration of the particle-number conservation [33] and the non-spherical mean fields [34] could influence $\Delta_{\text{mass}}^Z(N)$. Each of them could vary $\Delta_{\text{mass}}^Z(N)$ by up to a few hundred keV. Not attempting fine
tuning of the parameters, we just point out that some of these effects tend to compensate one another in the mass differences, and that the new interactions give $\Delta Z_{\text{mass}}$ to comparable accuracy to the D1S interaction in the same model space.

At $N \sim 50, 64$ and 90, we find that the calculated mass differences depend on the interactions. This is ascribed to interaction-dependence of the shell structure. Irregularity and discrepancy at $N = 63$ and 65 should be relevant to the $N = 64$ subshell. In M3Y-P5 the subshell effect seems stronger than in the other interactions. At $N \sim 90$ all the M3Y-type interactions yield larger mass differences than D1S. This takes place because $n1f_{7/2}$ and $n2p_{3/2}$ well mix due to the pairing, in the M3Y-type interactions. At $N \sim 50$ M3Y-P5 yields larger mass difference than the other interactions. This is traced back to appreciable excitation from $n1d_{5/2}$ to $n0g_{7/2}$, which takes place since these two orbits are close in energy. Possibly carrying information of the shell structure, data on the masses in $N \sim 50$ and $N \sim 90$ will be of interest.

B. Pairing gap in nuclear matter

We next view the pairing property in the nuclear matter obtained from the new semi-realistic interactions. In phenomenological studies using the Skyrme energy density functionals, it has been argued whether and how much the pair correlations are dominated by the nuclear surface region. Results of the semi-realistic interactions for the nuclear matter may provide certain information on this point. However, we find that, in calculating the pairing properties, the Yukawa function gives quite slow convergence for the maximum momentum of the s.p. states. In practice, even if we cut off the momentum at $k = 50 \text{fm}^{-1}$, which corresponds to $\varepsilon \approx 50 \text{GeV}$, the pairing gap is not yet fully convergent. It is impractical to include such high energy states in calculations of finite nuclei. As shown in
the preceding subsection, we have fixed the SE channel parameters from the HFB calculations of the Sn nuclei using the basis-parameters of Eq. (13). High momentum components are automatically excluded in the basis set. It will be natural to introduce a certain cut-off in arguing the pairing in the nuclear matter, and the cut-off should desirably be consistent with the basis set of Eq. (13).

The basis set is composed of radial Gaussians, whose Fourier transforms are again Gaussians in the momentum space. We here consider a cut-off factor for the s.p. momentum space of

\[ g(k) = \theta(k_c - k) + \theta(k - k_c) \exp \left[ -\frac{(k - k_c)^2}{k_d} \right]. \]  

(14)

The measure in the \( k \) integration is multiplied by \( g(k) \). Among the bases of Eq. (13), the highest \( k \) component is given by the \( \nu = \nu_0(1 \pm \frac{2}{\pi}i) \) basis. Since the Fourier transform of this basis is proportional to \( \exp \left[ -k^2(1 \pm \frac{2}{\pi}i)/4\nu_0(1 + \frac{2}{\pi}^3) \right] \), \( k \lesssim k_0 = 2\sqrt[3]{\nu_0(1 + \frac{2}{\pi}^3)}(\approx 1.55 \text{ fm}^{-1}) \) components are well included in the set. To be consistent with the basis set, it will be reasonable to take \( k_c \approx k_0 \) for the nuclear matter calculation. We here consider three cases, \((k_c, k_d) = (k_0, k_0)\), \((2k_0, k_0)\) and \((4 \text{ fm}^{-1}, 0)\). The last choice of the \( k_d \to 0 \) limit indicates a sharp cut-off, and \( k_c = 4 \text{ fm}^{-1} \) approximately corresponds to the maximum q.p. energy in the HFB calculations of the Sn nuclei.

In Fig. 7, the pairing gap at the Fermi energy, which is obtained from the Bardeen-Cooper-Schrieffer (BCS) calculation in the symmetric nuclear matter using the method of Ref. [36], is plotted as a function of \( k_F = (3\pi^2\rho/2)^{1/3} \). The pairing among like-nucleons arises from the SE channel in the effective interaction. The result of the M3Y-P5 interaction is compared with that of the Gogny D1S interaction. The cut-off is not needed for the Gogny interaction, and the present cut-off does not influence the D1S gap. The pairing gaps of M3Y-P3 and P4 are similar to that of M3Y-P5.

The new semi-realistic interactions are not drastically different from D1S, in respect to the pairing properties in the nuclear matter. The gap has a peak at \( k_F \approx 0.8 \text{ fm}^{-1} \approx 0.6k_F0 \), namely at \( \rho \approx 0.2\rho_0 \), for all cases. However, the peak height and the behavior at \( \rho > 0.2\rho_0 \) are different between the M3Y-type interactions and D1S. With rapid decrease at \( \rho > 0.2\rho_0 \), the M3Y-type interactions have more surface-dominant pairing than D1S. In D1S, the pair correlation may have sizable contribution from the bulk, even though it is stronger at the nuclear surface. Note that the cut-off parameter does not influence the nuclear matter pairing qualitatively, as long as it is more or less harmonious with the basis-set adopted in the calculations of finite nuclei.
Prediction of the neutron drip line depends on effective interactions to a certain degree. In this section we compare location of the neutron drip line predicted by the spherical HFB calculations with the present semi-realistic interactions and with the Gogny D1S interaction for the O, Ca and Ni isotopes. Although complete description of the drip line may require fine tuning of the parameters as well as taking account of correlation effects, it will be interesting to see what is relevant to location of the drip line.

A. \( Z = 8 \) nuclei

We present the two-neutron separation energies \( S_{2n} \) for the O isotopes in Fig. 8. The calculated values are compared with the experimental data. We do not show \( S_{2n} \) if the neutron chemical potential is positive. Though not displayed in Fig. 8, \( S_{2n} \) obtained from M3Y-P3 is close either to that from D1S or M3Y-P4. Whereas \( 25^{\text{–}} - 28^{\text{O}} \) have experimentally been established to be unbound \cite{37}, most MF calculations so far have failed to reproduce this nature. It would be noteworthy that one of the present semi-realistic interactions, M3Y-P5, correctly describes the location of the neutron drip line for oxygen within the spherical HFB approximation: \( 24^{\text{O}} \) is the heaviest bound oxygen isotope. The \( 25^{\text{O}} \) nucleus has higher energy than \( 24^{\text{O}} \), and in \( 26^{\text{–}} - 28^{\text{O}} \) the chemical potential becomes positive. In contrast, \( 26^{\text{O}} \) is bound in the HFB calculations with M3Y-P3 and P4, as in the calculation with the D1S interaction.

The reason why \( 26^{\text{O}} \) is not bound with M3Y-P5 can be traced back to \( \varepsilon_n(0d_{3/2}) \), the s.p. energy of \( n0d_{3/2} \). In Fig. 8 the neutron s.p. energies in the HF calculations are depicted. We show the energies obtained from the bases of Eq. (13) even when the s.p. energy is positive, not treating the boundary condition carefully. We view that M3Y-P5 gives higher \( \varepsilon_n(0d_{3/2}) \) than the other interactions, which originates from the slightly stronger \( v^{(LS)} \) as well as from the relatively small \( M^*_0 \). Note that we have fixed the enhancement factor for \( v^{(LS)} \) in M3Y-P5 so as to reproduce the s.p. spectrum around \( 208^{\text{Pb}} \), not adjusting accurately to, e.g., the \( ls \) splitting around \( 16^{\text{O}} \). It is commented that \( v^{(\text{TN})} \) has small but attractive contribution to \( \varepsilon_n(0d_{3/2}) \), and therefore is irrelevant to the higher \( \varepsilon_n(0d_{3/2}) \) in the M3Y-P5 result.
FIG. 8: $S_{2n}$ of the O isotopes ($N = \text{even}$). The calculations are performed in the spherical HFB approximation. Experimental data are taken from Ref. [31]. See Fig. 6 for conventions.

FIG. 9: HF single-particle energies in the O isotopes. Blue, green and red lines represent the results with the D1S, M3Y-P4 and M3Y-P5 interactions, respectively. For each interaction, dot-dashed line is for $\varepsilon_n(0d_{5/2})$, solid line for $\varepsilon_n(1s_{1/2})$ and dashed line for $\varepsilon_n(0d_{3/2})$. 
TABLE V: Neutron numbers of the heaviest bound Ca and Ni nuclei predicted by the spherical HFB calculations with several interactions.

| Isotope | D1S | M3Y-P3 | M3Y-P4 | M3Y-P5 |
|---------|-----|--------|--------|--------|
| Ca      | 44  | 50     | 48     | 50     |
| Ni      | 58  | 64     | 62     | 60     |

FIG. 10: Difference between the HF and HFB energies for the Ca isotopes ($N = \text{even}$), obtained from D1S, M3Y-P4 and P5. See Fig. 6 for conventions. Dotted lines are drawn to guide eyes.

B. $Z = 20$ and 28 nuclei

Location of the neutron drip line for the Ca and Ni nuclei could be investigated by the currently constructed or designed experimental facilities [38]. We tabulate location of the neutron drip line predicted by the spherical HFB calculations with the M3Y-type and the D1S interactions, in Table V.

If we use the D1S interaction, the heaviest bound Ca nucleus is $^{64}\text{Ca}$, because the neutron chemical potential is positive in $N \geq 45$. The M3Y-P3 and P5 interactions predict that $^{70}\text{Ca}$ is bound, while $^{68}\text{Ca}$ is the heaviest bound Ca isotope in the M3Y-P4 result. We depict difference between the HF and the HFB energies for the Ca nuclei in Fig. 10, which represents the pair correlation. Though not shown, M3Y-P3 gives similar results to M3Y-P5. While the pairing effects are in good agreement among all of these interactions in $N \leq 32$, the M3Y-type interactions give stronger pairing than D1S in $N \geq 34$.

Both in the predicted position of the neutron drip line and in the pair correlation in $N \geq 34$, the s.p. energy of $n 0g_{9/2}$ plays an important role. The neutron drip line can extend up to $^{70}\text{Ca}$ if $\varepsilon_n(0g_{9/2})$ is sufficiently low. At $^{60}\text{Ca}$ we have $\varepsilon_n(0g_{9/2}) = +0.73, -0.63, +0.23$ and $-0.65$ MeV in the HF calculations with D1S, M3Y-P3, P4 and P5, respectively, well correlated to the location of the drip line. The lower $\varepsilon_n(0g_{9/2})$ leads to the smaller shell gap at $N = 40$, $\varepsilon_n(0g_{9/2}) - \varepsilon_n(0f_{5/2})$, which makes the pair excitation across $N = 40$ easier. The gap is 3.8, 2.0, 3.7 and 1.7 MeV in D1S, M3Y-P3, P4 and P5. It is noted that this quenching of the shell gap in M3Y-P3 and P5 comes from the relatively strong $v^{(\text{LS})}$ to some degree, but not from $v^{(\text{TN})}$, because $v^{(\text{TN})}$ hardly contributes to the s.p. energies in an $\ell s$-closed shell.

The vanishing difference between the HF and HFB energies is often connected to the shell (or sub-shell) closure. Figure 10 indicates that, while $^{60}\text{Ca}$ is stiff against the pair excitation with D1S and M3Y-P4, significant pair excitation occurs by M3Y-P5 (and by M3Y-P3), because of the small shell gap at $N = 40$. On the contrary, $N = 50$ is stiff against the pair excitation in the present M3Y-P3 and P5 results, in which $^{70}\text{Ca}$ is bound. We have $\varepsilon_n(0g_{9/2}) \approx -2$ MeV.
FIG. 11: Difference between the HF and HFB energies for the Ni isotopes ($N = \text{even}$). See Fig. 6 for conventions.

at $^{70}\text{Ca}$, and the gap between $n0g_{9/2}$ and the continuum seems large enough for $N = 50$ to keep the magic nature against the pairing, if we use the present semi-realistic pairing interaction. Figure 10 also suggests shell closure at $N = 32$. We shall return to this point in Sec. VII.

In Fig. 11, difference between the HF and HFB energies is shown for the Ni isotopes. For all the interactions, the energy difference becomes vanishingly small at $N = 20, 28, 40, 50$ and 58. We view significant interaction-dependence in $28 < N < 40$. In particular, the pair correlation is suppressed at $N = 32$ with D1S, while no such effect is found with M3Y-P5. It is noted that, with M3Y-P5, $^{68}\text{Ni}$ seems almost doubly magic as is consistent with experiments [39], although $^{60}\text{Ca}$ is not, as has been seen in Fig. 10. For the neutron-rich Ni region, energy sequence of the s.p. orbitals above $N = 50$ is $1d_{5/2}, 2s_{1/2}, 1d_{3/2}$ and $0g_{7/2}$, from the lower orbit to the higher. The hindrance of the pair excitation at $^{86}\text{Ni}$ suggests magic or submagic nature of $N = 58$ due to the gap between $n2s_{1/2}$ and $n1d_{3/2}$. Unlike the M3Y-type interactions, by D1S the pair excitation is hindered also at $N = 56$. The predicted neutron drip line appreciably depends on $\varepsilon_n(1d_{3/2})$ and $\varepsilon_n(0g_{7/2})$. The D1S interaction yields higher $\varepsilon_n(1d_{3/2})$ than the M3Y-type interactions, which causes the drip line at $N = 58$. The low $\varepsilon_n(0g_{7/2})$ in M3Y-P3 induces pair excitation to $0g_{7/2}$, leading to the binding up to $^{92}\text{Ni}$.

In the highly neutron-rich region, the diffuseness of the nuclear surface becomes larger than in the $\beta$-stable region. Then the pair correlation could be relatively strong if it has the surface-dominant nature. However, it is not easy to argue precisely the extent of surface dominance in the pairing from location of the drip line, since it is obscured by influence of the shell structure. In practice, when we use M3Y-P5 for the HF Hamiltonian and D1S for the pair potential, predicted location of the drip line for Ca is the same as the result of the pure M3Y-P5 prediction. For the Ni case, although $^{88}\text{Ni}$ becomes unbound when D1S is used for the pair potential, the chemical potential is only $-0.07\text{MeV}$ in the pure M3Y-P5 result. It will be fair to say that the difference in the pairing channel between the M3Y-type interactions and D1S is not quite significant to location of the neutron drip line. Dependence of the rms matter radii on the pairing interaction is not apparent either, as long as we work with the present M3Y-type or the D1S interactions, whereas the radii near the drip line are sensitive to the separation energies.

VII. NUCLEUS-DEPENDENCE OF SINGLE PARTICLE ENERGIES

It has been pointed out that the shell structure, particularly its nucleus-dependence (sometimes called shell evolution), may be connected to characters of effective interactions [3, 10, 40]. In recent studies role of the tensor force in the shell structure has attracted great interest [3, 4]. It is known that observed energies of one-particle states on top of a certain core are appreciably disturbed by correlations beyond the MF regime. However, it is presumable that those correlations do not vary in a certain region of nuclei. In the Sb and the $N = 83$ nuclei, s.p. energies of a few orbitals are extracted from several fragmented states [41], by averaging their energies weighted by the spectroscopic factors. As a result, the averaged s.p. energies are shifted from the lowest states with specific spin-parity nearly by a constant, from nucleus to nucleus. While data on the averaged energies are not available in many cases, we proceed to investigate nucleus-dependence of the shell structure by using the measured energies of the lowest states.
In this section we shall investigate nucleus-dependence of s.p. energies, using the spherical HF or HFB calculations. Results of several interactions are compared. In the HF calculations for open-shell nuclei, the HF Hamiltonian is obtained by folding the interaction by the occupation numbers on each spherical orbital up to the Fermi level.

A. Neutron orbits and shell gap in $N=16$ nuclei

Nucleus-dependence of s.p. energies could be relevant to the new magic numbers in unstable nuclei [3, 40]. We investigated the s.p. energies in the $N=16$ and 32 nuclei in Refs. [10, 11], and disclosed role of $v^{(C)}_{\text{OPEP}}$ for the $N=16$ isotones, using the M3Y-P2 interaction. We shall reinvestigate nucleus-dependence of the s.p. energies in these nuclei, drawing attention also to the tensor force.

$Z$-dependence of the neutron s.p. energy $\varepsilon_n(0d_{3/2})$ relative to $\varepsilon_n(1s_{1/2})$ is appreciably affected by effective interactions [10, 42]. Figure 12 depicts $\Delta\varepsilon_n = \varepsilon_n(0d_{3/2}) - \varepsilon_n(1s_{1/2})$ for varying $Z$ obtained from the HF calculations in the $N=16$ isotones. Though not shown, $\Delta\varepsilon_n$ of M3Y-P3 resembles that of M3Y-P5. To clarify role of $v^{(TN)}$ and $v^{(C)}_{\text{OPEP}}$, we also plot their contributions to $\Delta\varepsilon_n$ in the M3Y-P5 result, which are calculated as $\sum_{j'} \langle N_{j'} | (2J+1) (jj'J|v|jj'J)/(2j+1)(2j'+1) and shifted by the values at $N=14$.

The present semi-realistic interactions yield increasing $\Delta\varepsilon_n$ as $Z$ goes from 14 to 8, in contrast to the D1S interaction. We view in Fig. 12 that $v^{(C)}_{\text{OPEP}}$ [42] and $v^{(TN)}$ produce this feature cooperatively. This $Z$-dependence of the s.p. energies could be relevant to the new magic number $N=16$ in the neutron-rich region [43]. It has been confirmed that several popular Skyrme interactions show similar behavior to D1S [42].

FIG. 12: $\Delta\varepsilon_n = \varepsilon_n(0d_{3/2}) - \varepsilon_n(1s_{1/2})$ for the $N=16$ isotones. Blue, green and red lines correspond to the results with the D1S, M3Y-P4 and P5 interactions, respectively. Thin red solid and dashed lines represent relative contributions of $v^{(TN)}$ and $v^{(C)}_{\text{OPEP}}$ in the M3Y-P5 result.
B. Neutron orbits and shell gap in $N = 32$ nuclei

In Fig. 13 $Z$-dependence of the neutron s.p. energies relative to $\varepsilon_n(1p_{3/2})$, $\Delta\varepsilon_n(j) = \varepsilon_n(j) - \varepsilon_n(1p_{3/2})$, is shown for the $N = 32$ nuclei, by taking $j = 0f_{5/2}$ and $1p_{1/2}$. As in the preceding subsection, contributions of $v^{(\text{TN})}$ and $v^{(\text{C})}_{\text{OPEP}}$ to $\Delta\varepsilon_n(0f_{5/2})$ in the M3Y-P5 result are also presented, after shifting by the values at $Z = 28$. With the M3Y-P5 interaction we obtain strong $Z$-dependence in $\Delta\varepsilon_n(0f_{5/2})$, which could be relevant to the magicity of $N = 32$ in the neutron-rich region \cite{ref44}. Once again this $Z$-dependence originates in $v^{(\text{C})}_{\text{OPEP}}$ \cite{ref11} and $v^{(\text{TN})}$.

This behavior of $\Delta\varepsilon_n(0f_{5/2})$ is reflected in the pair correlations shown in Figs. 10 and 11. The pairing effects are small at $N = 32$ for all the interactions in the calcium case, because $\Delta\varepsilon_n(0f_{5/2})$ as well as $\Delta\varepsilon_n(1p_{3/2})$ are greater than 2 MeV. Recall that this s.p. energy difference competes with the pairing gap, whose typical value is estimated to be $\Delta \approx 12A^{-1/2} \approx 1.7$ MeV. On the other hand, in the nickel case the narrow $\Delta\varepsilon_n(0f_{5/2})$ leads to substantial pair excitation at $N = 32$ for the M3Y-type interactions, while such excitation is kept suppressed in the D1S result.

In the neutron-rich region of $Z \sim 20$, there was a prediction that $N = 34$ should be a magic number, based on a shell model calculation \cite{ref47}. The present MF calculations with the semi-realistic interactions do not support this prediction. While the $N = 32$ shell gap is 2.7 MeV for $^{52}\text{Ca}$ in the HF calculation with M3Y-P5, the $N = 34$ gap is only 1.2 MeV for $^{54}\text{Ca}$. The pair excitation across $N = 34$ is sizable, as has been viewed in Fig. 10.

C. Proton orbits in $Z = 50$ nuclei

In recent studies, nucleus-dependence of s.p. energies in the Sn isotopes and in the $N = 82$ isotones has been disclosed from experiments \cite{ref41}. It has been pointed out that the tensor force seems to play a crucial role in the
The semi-realistic M3Y-P5 interaction reproduces $\delta \Delta \varepsilon_p(j)$ remarkably well. In the $64 \leq N \leq 82$ region, the $N$-dependence of $\Delta \varepsilon_p(j)$ takes place due primarily to the occupation of $1d_{5/2}$, to which contribution of $v^{(C)}$ is not, because the tensor force is absent. The same holds for M3Y-P4. Although M3Y-P4 reproduces the tendency of the observed $N$-dependence of $\varepsilon_p(0h_{11/2}) - \varepsilon_p(0g_{7/2})$, it gives wrong behavior for $\Delta \varepsilon_p(0h_{11/2})$; i.e., $N$-dependence of $\varepsilon_p(0h_{11/2})$ relative to $\varepsilon_p(1d_{5/2})$. It is also remarked that M3Y-P5 gives quite different behavior of $\delta \Delta \varepsilon_p(j)$ from D1S and M3Y-P4 in $N < 64$, and that the currently available data favor the result of M3Y-P5. We have confirmed that $v^{(C)}_{\text{OEP}}$ does not have important effects on the $N$-dependence of $\Delta \varepsilon_p(j)$.

It is emphasized that the M3Y-P5 interaction can reproduce the variation of the s.p. energy difference in the Sn isotopes without destroying the shell structure of the doubly magic nuclei shown in Sec. IV. This could be an advantage of the present realistic tensor force. In Ref. [4], no such parameters were found within the Skyrme density functional including the zero-range tensor force.
FIG. 15: \( \delta \Delta \varepsilon_n(j) \) in the \( N = 50 \) isotones (\( Z = \text{even} \)), for \( j = 0 \ell \) and \( 0 h_{11/2} \). Conventions are the same as in Fig. 14 except that \( j \) represents the neutron orbits and \( Z = 40 \) is taken to be a reference. Experimental values are taken from the lowest states of the \( N = 51 \) nuclei [32].

**D. Neutron orbits in \( N = 50 \) and \( N = 82 \) nuclei**

In the preceding subsection, we have seen that the tensor force affects \( \varepsilon_p(0\ell_{7/2}) \) and \( \varepsilon_p(0h_{11/2}) \) via the occupation of \( n0h_{11/2} \). This is accounted for by the attractive (repulsive) nature of the tensor force between a neutron occupying a \( j_n = \ell + 1/2 \) orbit and a proton occupying \( j_p' = \ell' - 1/2 \) (\( j_n' = \ell' + 1/2 \)) [18]. The same mechanism is expected for \( \varepsilon_n(0\ell_{7/2}) \) and \( \varepsilon_n(0h_{11/2}) \) in the \( N = 50 \) nuclei, as \( p0h_{9/2} \) is occupied. We define \( \Delta \varepsilon_n(j) = \varepsilon_n(j) - \varepsilon_n(1d_{5/2}) \), and take its value at \( Z = 40 \) to be \( \Delta \varepsilon_n^0(j) \). In Fig. 15 \( \Delta \varepsilon_n(j) = \Delta \varepsilon_n(j) - \Delta \varepsilon_n^0(j) \) is displayed for \( j = 0\ell_{7/2} \) and \( 0h_{11/2} \). Although \( \varepsilon_n(0\ell_{7/2}) \) varies almost in parallel to \( \varepsilon_n(0h_{11/2}) \) in \( Z \geq 40 \) for D1S, notable \( Z \)-dependence arises for M3Y-P5, in qualitative agreement with the observed s.p. levels. The tensor force has significant contribution to this behavior as in the \( Z = 50 \) case. More realistic than D1S in the central channels but not having the tensor force, M3Y-P4 yields \( \delta \Delta \varepsilon_n(j) \) in-between which is qualitatively good but quantitatively insufficient.

For the \( N = 82 \) nuclei, we consider \( \Delta \varepsilon_n(j) = \varepsilon_n(j) - \varepsilon_n(1f_{7/2}) \), from which \( \delta \Delta \varepsilon_n(j) = \Delta \varepsilon_n(j) - \Delta \varepsilon_n^0(j) \) is obtained by assuming the value at \( Z = 64 \) as \( \Delta \varepsilon_n^0(j) \). Figure 16 shows \( \delta \Delta \varepsilon_n(j) \) for \( j = 0h_{9/2} \) and \( 0i_{13/2} \). The M3Y-P5 interaction well describes \( \delta \Delta \varepsilon_n(j) \) in \( Z \geq 64 \), which is affected mainly by the occupation of \( p0h_{11/2} \). However, we cannot fully reproduce the tendency in \( Z < 64 \), which was argued in Ref. [41], although the M3Y-P5 results of \( \delta \Delta \varepsilon_n(j) \) are substantially better than those of D1S and M3Y-P4. Further investigation will be necessary.

**VIII. SUMMARY AND OUTLOOK**

We have developed semi-realistic effective interactions to describe low energy phenomena of nuclei. Starting from the M3Y interaction, we add a density-dependent contact force and modify several strength parameters in a phenomenological manner, whereas maintaining the OPEP part in the central force. We have obtained three new parameter-sets; two of them (M3Y-P3 and P5) keep the tensor force of the M3Y-Paris interaction, and the other (M3Y-P4) has no tensor force. Basic characters of the interactions are checked by the Hartree-Fock calculations for the infinite nuclear matter, and for the doubly magic nuclei. The singlet-even channels of the interactions, which are relevant to the pairing properties, are fixed from the even-odd mass differences in the Sn isotopes, by using the Hartree-Fock-Bogolyubov calculations.

We further implement the Hartree-Fock and the Hartree-Fock-Bogolyubov calculations for spherical nuclei, applying the new interactions. Predicted shell structure depends on the effective interactions to certain extent. This may significantly affect location of the drip lines. The new semi-realistic interaction M3Y-P5 correctly describes the
experimental consequence that the heaviest bound oxygen is $^{24}$O. We have argued location of the neutron drip line for the Ca and the Ni nuclei, and its relevance to the shell structure. Variation of the single-particle (s.p.) energies, particularly contribution of the tensor force to it, is a current topic. It is suggested that the tensor force as well as the OPEP part of the central force play a significant role in the magic numbers $N = 16$ and $32$ in the neutron-rich region. It has been shown that the new semi-realistic interactions including the tensor force, M3Y-P5 in particular, describe the variation of the s.p. levels fairly well in $Z = 50$, $N = 50$ and $N = 82$ nuclei. It is remarked that this interaction can reproduce the variation of the s.p. energy difference in the Sn isotopes without destroying the shell structure of the doubly magic nuclei.

It will be of interest to apply the semi-realistic interactions to deformed nuclei, and to excited states via the random-phase approximation (RPA). Both projects are in progress (for the latter, see Ref. [46]).

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