

Constraining the Nuclear Energy Density Functional by Low-Energy QCD \diamond

P. Finelli, N. Kaiser, D. Vretenar, and W. Weise

A relativistic nuclear energy density functional

$$E_0[\rho] = E_{\text{free}}[\rho] + E_{\text{H}}[\rho] + E_{\text{coul}}[\rho] + E_{\pi}[\rho] \quad (1)$$

has been developed, guided by two important features that establish connections with chiral dynamics and the symmetry breaking pattern of low-energy QCD:

1. large scalar and vector mean fields (with opposite sign) related to in-medium changes of QCD vacuum condensates [1], determining the Hartree energy functional $E_{\text{H}}[\rho]$;
2. long- and intermediate-range interactions generated by one- and two-pion exchange and derived from in-medium chiral perturbation theory [2], with explicit inclusion of $\Delta(1232)$ excitations, determining the exchange correlation energy functional $E_{\pi}[\rho]$.

We have constructed a point-coupling model with density dependent interaction terms, in which the single terms of (1) reads [3]

$$E_{\text{free}}[\rho] = \int d^3r \langle \phi_0 | \bar{\psi} [-i\boldsymbol{\gamma} \cdot \boldsymbol{\nabla} + M_N] \psi | \phi_0 \rangle, \quad (2)$$

$$E_{\text{H}}[\rho] = \frac{1}{2} \int d^3r [G_S^{(0)} \rho_S^2 + G_V^{(0)} \rho^2], \quad (3)$$

$$E_{\pi}[\rho] = \frac{1}{2} \left\{ \int d^3r [G_S^{(\pi)}(\rho) \rho_S^2 + G_V^{(\pi)}(\rho) \rho^2] + \int d^3r [G_{TS}^{(\pi)}(\rho) \rho_{S3}^2 + G_{TV}^{(\pi)}(\rho) \rho_3^2] - \int d^3r [D_S^{(\pi)}(\vec{\nabla} \rho_S)^2] \right\}, \quad (4)$$

$$E_{\text{coul}}[\rho] = \int d^3r A^0 e \frac{1 + \tau_3}{2} \rho. \quad (5)$$

The overall structure of $E_0[\rho]$ is reminiscent of what is commonly introduced for relativistic atomic systems in a DFT approach [4]. E_{π} plays the role of the exchange correlation term E_{exc} and E_{H} is the nuclear counterpart of the Hartree term.

This approach has been tested in the analysis of ground-state observables of spherical and deformed nuclei. In Fig. 1 we show recent results of an extensive study of ground-state properties of deformed isotope chains with $60 \leq Z \leq 80$, calculated in the Relativistic Hartree-Bogoliubov approximation [5].

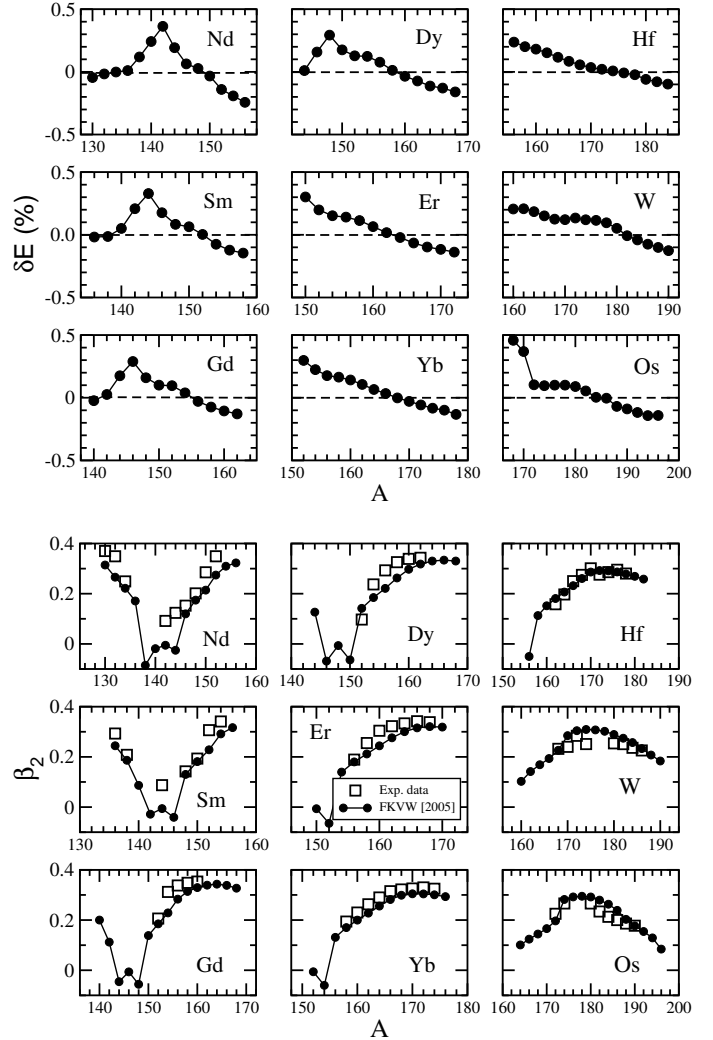


Fig. 1: The deviations (in percent) of the calculated binding energies and the predictions for the ground-state quadrupole deformations of the Nd, Sm, Gd, Dy, Er, Yb, Hf, W, and Os isotopes.

The agreement with experimental data is very good over a wide range of isotopic chains and it demonstrates the validity of our basic assumptions. Study of collective excitations is in progress.

References

- [1] T. D. Cohen, R. J. Furnstahl, D. K. Griegel and X.m. Jin, Prog. Part. Nucl. Phys. **35** (1995) 221
- [2] S. Fritsch, N. Kaiser and W. Weise, Nucl. Phys. **A750** (2005) 259
- [3] P. Finelli, N. Kaiser, D. Vretenar and W. Weise, Nucl. Phys. **A735** (2004) 449
- [4] R.M. Dreizler and E.K.U. Gross, Density Functional theory, Springer-Verlag, 1990; R.M. Dreizler, Lect. Notes Phys. **620** (2003) 123
- [5] P. Finelli, N. Kaiser, D. Vretenar and W. Weise, [arXiv:nucl-th/0509040], Nucl. Phys. A . (in print)

\diamond Work supported by BMBF, GSI, MURST and INFN