## Shell-model fits for Sn isotopes

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[NUCLEAR STRUCTURE, shell model]

The Sn isotopes provide us with an ideal test for conventional shell-model calculations on the basis of the microscopic effective interaction theory and also the knowledge of nuclear forces. Because of the  $Z{=}50$  semi-magic structure, the deformation is not significant, and the the shell-model description within a limited one-major shell is expected be reasonable. In fact, shell-model calculations using an effective interaction derived microscopically from a realistic nucleon-nucleon potential have been successfully applied for the analysis of new experimental data and also the prediction for unobserved properties  $^{1)}$ .

However, such a microscopic approach is not necessarily successful in the middle of the neutron shell as shown in Ref.  $^2$  for example, where the excitation energy of the  $2_1^+$  state is predicted to be too high by 0.4MeV for  $^{118}$ Sn. On the other hand, it has been shown that the microscopic effective interaction can be improved for a practical use by fitting to the experimental data. Therefore, it is interesting to investigate how accurate the structure of Sn isotopes can be described by using such a fitted effective interaction. We report the results of such an attempt.

The adopted model space consists of five singleparticle orbits  $1d_{5/2}$ ,  $0g_{7/2}$ ,  $0h_{11/2}$ ,  $2s_{1/2}$  and  $1d_{3/2}$ in which the Sn isotopes are described by valence neutrons, and the T=1 part of the two-body interactions is relevant. We start with an effective interaction obtained in a microscopic way as described in Ref.<sup>3)</sup> using the N<sup>3</sup>LO interaction<sup>4)</sup>. A series of fitting calculations was carried out by taking 313 experimental energy data including 29 binding energies. We utilized the Linear-Combination (LC) method as in our previous approach<sup>5)</sup>. We found that only a small modification was necessary to improve the microscopic interaction. In fact a reasonable convergence was obtained by varying only 22 LC's of 165 Hamiltonian parameters (5 single-particle energies and 160 two-body matrix elements). In the final fit, the rms deviation of 147 keV was obtained.

As an example of the results, the excitation energies of the yrast states of even-N isotopes are shown in Fig.1. One can find a reasonable agreement between the experimental data and the shell-model results throughout the isotope chain. The quality of the

fit is similar for odd-N isotopes. The predicted ground-state spin of  $^{101}\mathrm{Sn}$  is  $7/2^+$  rather than  $5/2^+$ , in good agreement with recent experimental observation. The fit of the binding energy is basically successful for heavier isotopes, but the shell-model result begins to show underbinding as approaching  $^{100}\mathrm{Sn}$ . This result suggests the development of collectivity which can not be described by the present model space. Such a picture is consistent with the large  $B(\mathrm{E2})$  values below N=64 observed in recent extensive measurements.

This fitted interaction can be a good starting point for extending the model space so as to describe the <sup>100</sup>Sn core excitations and also for the study of more neutron-rich isotopes.

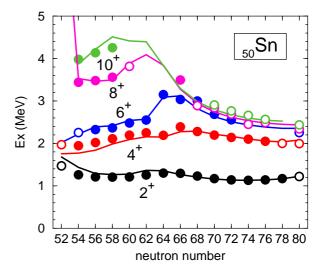


Fig. 1. Comparison of excitation energies of even-spin yrast states between the experimental data (symbols) and the shell-model results (lines). Experimental data are taken from Ref.<sup>8)</sup> Open symbols indicate that the spin assignment is uncertain. The shell-model results are obtained by using the efficient shell-model code MSHELL<sup>9)</sup>.

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