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Microscopic analysis of T = 1 and T = 0proton–neutron correlations in N = Z nuclei

M. Hasegawa^{a,*}, K. Kaneko^b

^a Laboratory of Physics, Fukuoka Dental College, Fukuoka 814-0193, Japan ^b Department of Physics, Kyushu Sangyo University, Fukuoka 813-8503, Japan

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Abstract

The competition between the isovector (T = 1) and isoscalar (T = 0) proton-neutron (p-n) correlations in N = Z nuclei is investigated by calculating their correlation energies with a realistic effective interaction which reproduces observed nuclear properties very well, in a strict shell model treatment. It is shown in the realistic shell model that the double-differences of binding energies (B(A + pn : T) + B(A) - B(A + p) - B(A + n)) (B(A) being the binding energy) are good indicators of the T = 1 and T = 0 p-n correlations. Each of them, however, originates in plural kinds of correlations with T = 1 or T = 0.

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1. Introduction

The competition between isovector (T = 1) and isoscalar (T = 0) pairing correlations has been a matter of renewed concern in nuclear structure studies of $N \approx Z$ nuclei [1–3]. The competition appears in the near degeneracy of the lowest T = 1 and T = 0 states in odd–odd N = Z nuclei. The T = 0 proton–neutron (p-n) pairing correlations in $N \approx Z$

Corresponding author.

E-mail address: hasegawa@college.fdcnet.ac.jp (M. Hasegawa).

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nuclei have been studied by the two approaches [2,3] with different treatments of the symmetry energy. The two conclusions about the importance of the T = 0 *p*–*n* pairing correlations are in opposition to one another. The T = 1 and T = 0 pairing correlations, which are considered to be induced by T = 1 and T = 0 nuclear interactions, should be treated consistently on the same footing [3]. The structure of N = Z nuclei has been well described by the shell model which does treat T = 1 and T = 0 pairing consistently. Large-scale shell model calculations were applied to the investigation of the isovector and isoscalar pairing correlations in Refs. [4–6], where the contributions of T = 1, J = 0 and T = 0, J = 1 interactions are compared [4] and the contributions of the quadrupole–quadrupole (QQ) force are also considered [5]. The authors have recently shown that the competition between the T = 1 and T = 0 pairing correlations are approximately explained with the T = 1, J = 0 pairing force (P_0) and a T = 0 p–*n* force ($V_{mp}^{T=0}$ below) [7]. In order to understand the competition in detail, it is important to evaluate the two types of correlations induced by realistic effective interactions.

In this paper, we investigate the competition between T = 1 and T = 0 *p*–*n* correlations in the lowest states of N = Z nuclei using a realistic effective interaction in a strict shell model treatment. The spherical shell model, which gives an excellent description of various properties of $N \approx Z$ nuclei (not only the binding energies but also other properties), has the advantage that the correlation energies of respective interactions are properly calculated. A shell model Hamiltonian is composed of the T = 0 and T = 1 interactions,

$$H = H_{\rm sp} + V^{T=0} + V^{T=1},\tag{1}$$

$$V^{T} = \sum_{a \leq b} \sum_{c \leq d} \sum_{JM} \sum_{TK} G_{JT}(ab:cd) A^{\dagger}_{JMTK}(ab) A_{JMTK}(cd) \quad (T = 0, 1),$$
(2)

where H_{sp} stands for the single-particle energies, $A_{JMTK}^{\dagger}(ab)$ is the creation operator of a nucleon pair with the spin JM and the isospin TK in the single-particle orbits (a, b), and $G_{JT}(ab:cd)$ denotes the interaction matrix elements. The so-called realistic effective interactions contain multipole $(J \ge 0)$ pairing forces of T = 0 and T = 1 in the expression (2). In this sense, the shell model with a realistic effective interaction is to deal with all the *multipole pairing correlations*. We investigate the competition between the T = 0and T = 1 correlations induced by the T = 0 and T = 1 interactions.

The realistic effective interactions have the property that the centroid of T = 0 diagonal interaction matrix elements $\overline{G_{T=0}(ab)} = \sum_{J} (2J+1)G_{J0}(ab:ab) / \sum_{J} (2J+1)$ has a roughly constant value, being independent on the orbits (*ab*) [8]. By setting $-k^0 = \sum_{ab} \overline{G_{T=0}(ab)} / \sum_{ab}$, we obtain the average T = 0 *p*-*n* force

$$V_{\rm mp}^{T=0} = -k^0 \sum_{a \leqslant b} \sum_{JM} A_{JM00}^{\dagger}(ab) A_{JM00}(ab).$$
(3)

Let us write residual T = 0 interactions as $V_{\text{res}}^{T=0} = V^{T=0} - V_{\text{mp}}^{T=0}$ and rewrite the Hamiltonian as

$$H = H_{\rm sp} + V_{\rm mp}^{T=0} + V_{\rm res}^{T=0} + V^{T=1}.$$
 (4)

The separation of $V_{mp}^{T=0}$ in Eq. (4) follows the procedure of Dufour and Zuker in Ref. [9], where the Hamiltonian is divided into the monopole and multipole parts as $H = H_m + H_M$.

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