

原子核結合エネルギーに対する 低運動量相互作用の切断運動量依存性

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1. はじめに

- ・現実的核力から導かれる低運動量相互作用 ($V_{\text{low-}k}$)
- ・ $V_{\text{low-}k}$ を用いた最近の構造計算

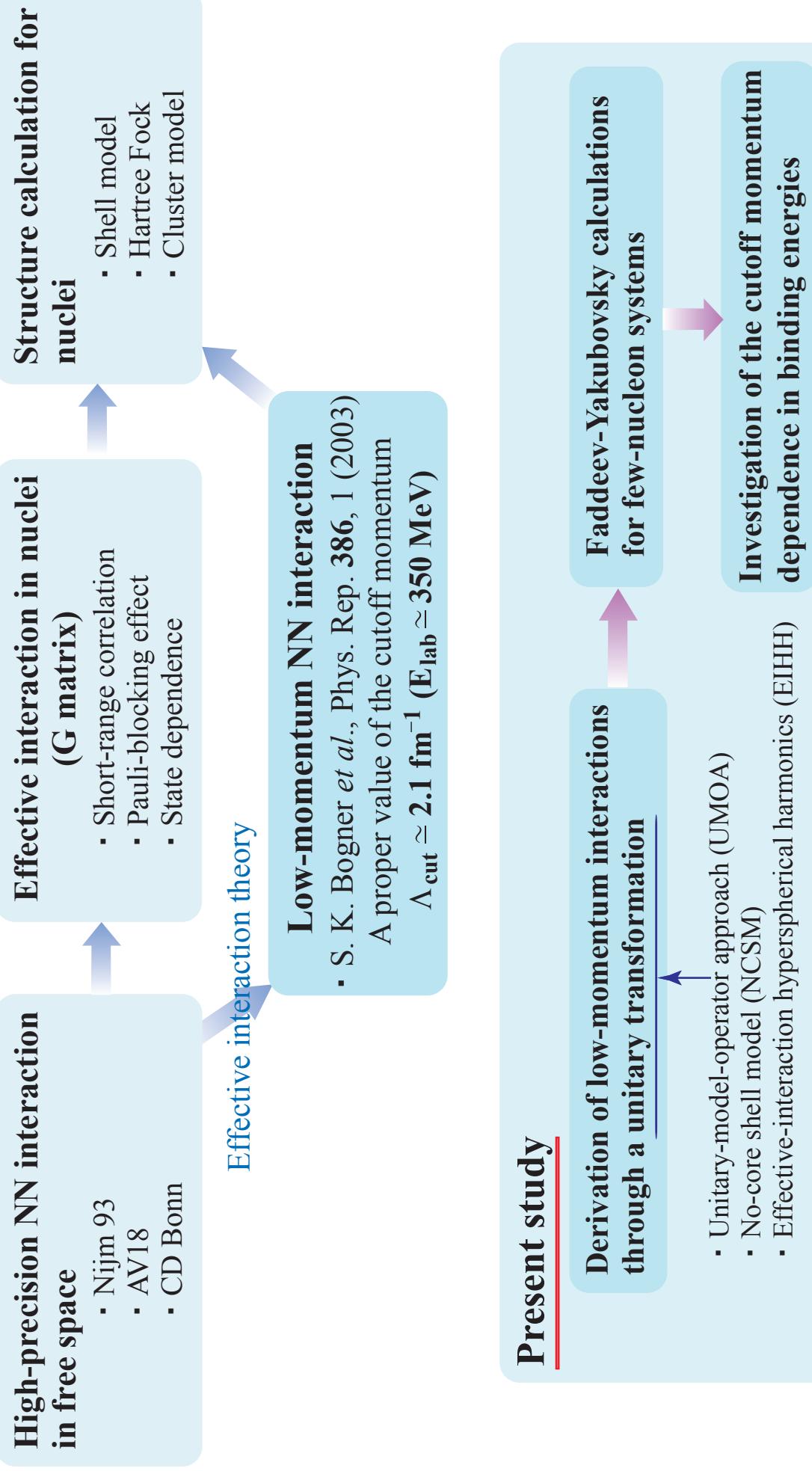
2. ユニタリー変換による $V_{\text{low-}k}$ の導出

3. $V_{\text{low-}k}$ を用いた結合エネルギーの計算結果

- ・ ${}^3\text{H}, {}^4\text{He} \cdots$ Faddeev-Yakubovsky
- ・ ${}^{16}\text{O} ({}^{15}\text{N}, {}^{15}\text{O}, {}^{17}\text{F}, {}^{17}\text{O}) \cdots$ UMOA

4. まとめ

Microscopic description of nuclear structure



Derivation of effective interaction (Hamiltonian) by means of unitary transformation

Hamiltonian

$$H = H_0 + V$$

Unitary transformation of H

$$\tilde{H} = U^{-1}HU$$

$$U = e^S, \quad (S : \text{anti-Hermitian}, S^\dagger = -S)$$

Decoupling equation

$$Q(e^{-S}He^S)P = 0$$

Solution

$$S = \operatorname{arctanh}(\omega - \omega^\dagger), \quad \omega = Q\omega P$$

(with the restrictive condition $PSP = QSQ = 0$)

K. Suzuki, Prog. Theor. Phys. **68** (1982), 246

Effective Hamiltonian

$$\underline{H_{\text{eff}} = P\tilde{H}P}$$

Effective interaction

$$\underline{V_{\text{eff}} = P\tilde{H}P - PH_0P}$$

Unitary transformation operator U in terms of ω

$$U = (1 + \omega - \omega^\dagger)(1 + \omega^\dagger\omega + \omega\omega^\dagger)^{-1/2}$$

$$= \begin{pmatrix} P(1 + \omega^\dagger\omega)^{-1/2}P & -P\omega^\dagger(1 + \omega\omega^\dagger)^{-1/2}Q \\ Q\omega(1 + \omega^\dagger\omega)^{-1/2}P & Q(1 + \omega\omega^\dagger)^{-1/2}Q \end{pmatrix}$$

S. Ōkubo, Prog. Theor. Phys. **12** (1954), 603

Application of effective interaction theory to the two-nucleon system in momentum space

Derivation of low-momentum interaction

$$\begin{array}{ccc}
 H = & \left(\begin{array}{cccccc} k_1 k_2 & \dots & \dots & \dots & \dots & k_{max} \\ \vdots & & & & & \vdots \\ k_1 & k_2 & & & & \\ \vdots & & & & & \\ \vdots & & & & & \\ k_1 & k_2 & & & & \\ \vdots & & & & & \\ \vdots & & & & & \\ k_1 k_2 \dots k_{cut} & & & & & \\ \vdots & & & & & \\ k_{max} & & & & & \end{array} \right) & \xrightarrow{\text{unitary transformation}} \\
 \widetilde{H} = U^{-1} H U & \xrightarrow{\quad\quad\quad} & \widetilde{H} = \left(\begin{array}{cc} P \widetilde{H} P & 0 \\ 0 & Q \widetilde{H} Q \end{array} \right)
 \end{array}$$

Unitary transformation operator

$$U = (1 + \omega - \omega^\dagger)(1 + \omega^\dagger \omega + \omega \omega^\dagger)^{-1/2}, \quad \omega = Q \omega P$$

Low-momentum interaction

$$\underline{V_{low-k} = P(\widetilde{H} - T)P}$$

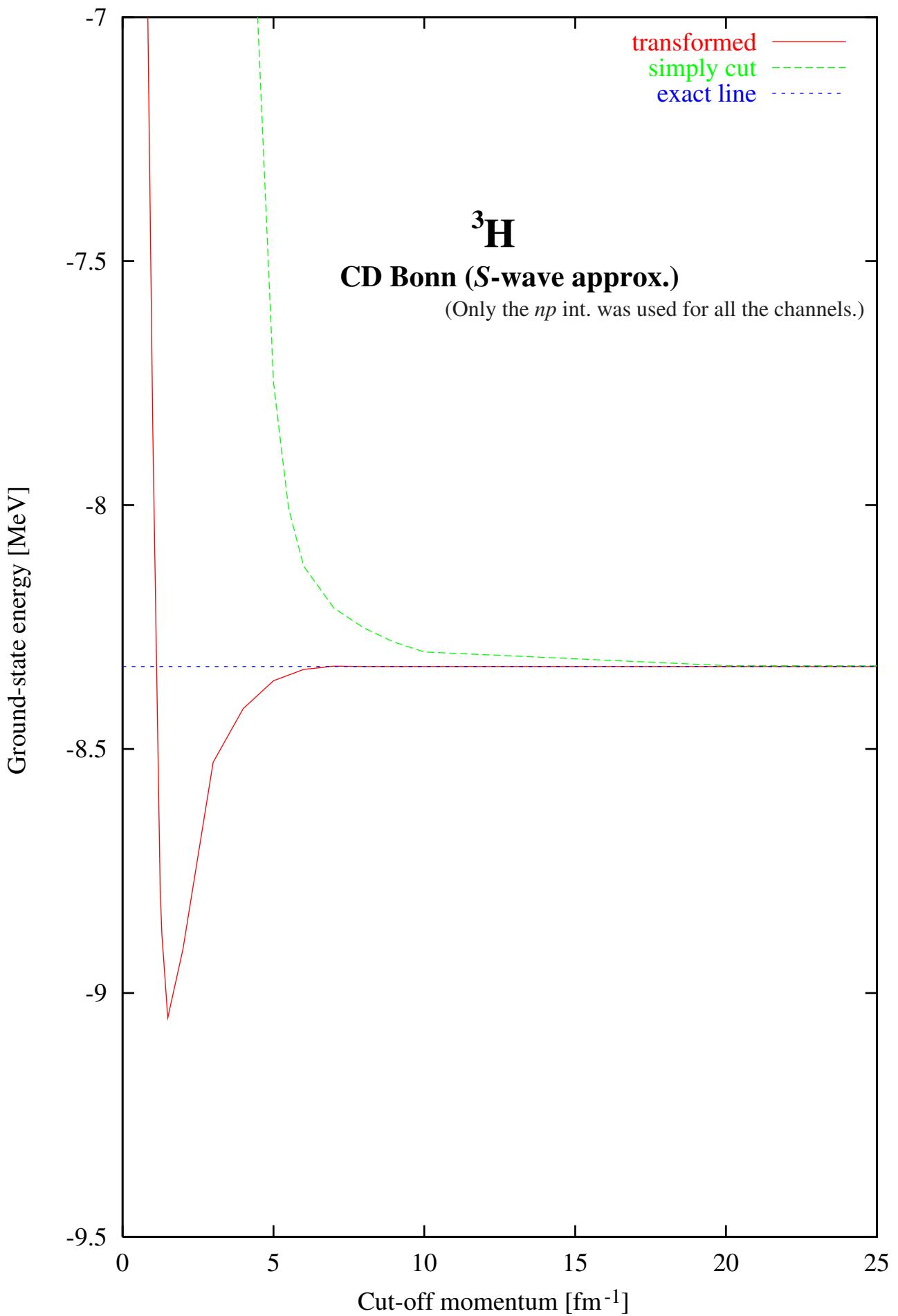
Properties of V_{low-k}

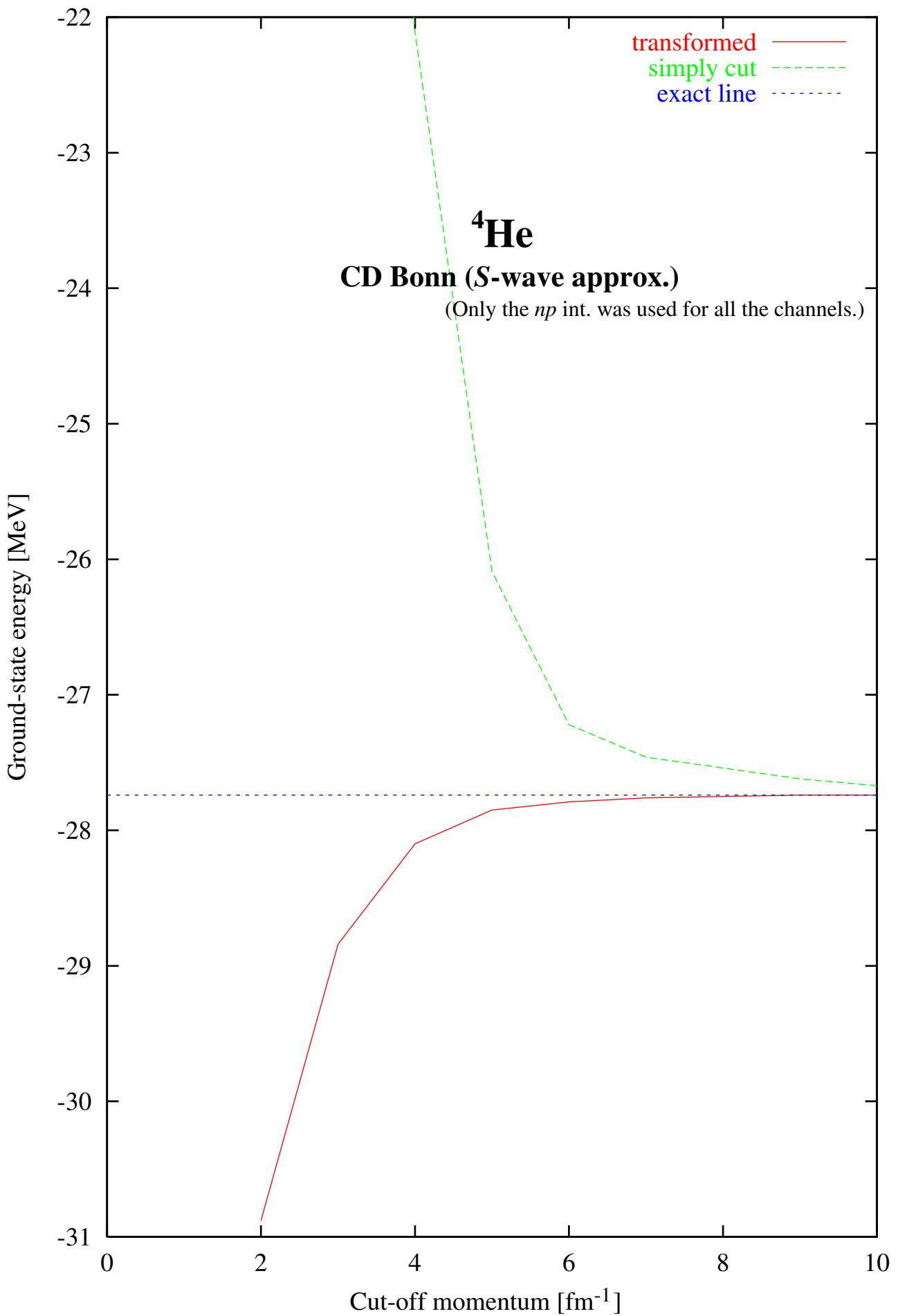
- $Q V_{low-k} P = P V_{low-k} Q = 0$
- Energy independent
- Hermitian

Accuracy of low-momentum interactions

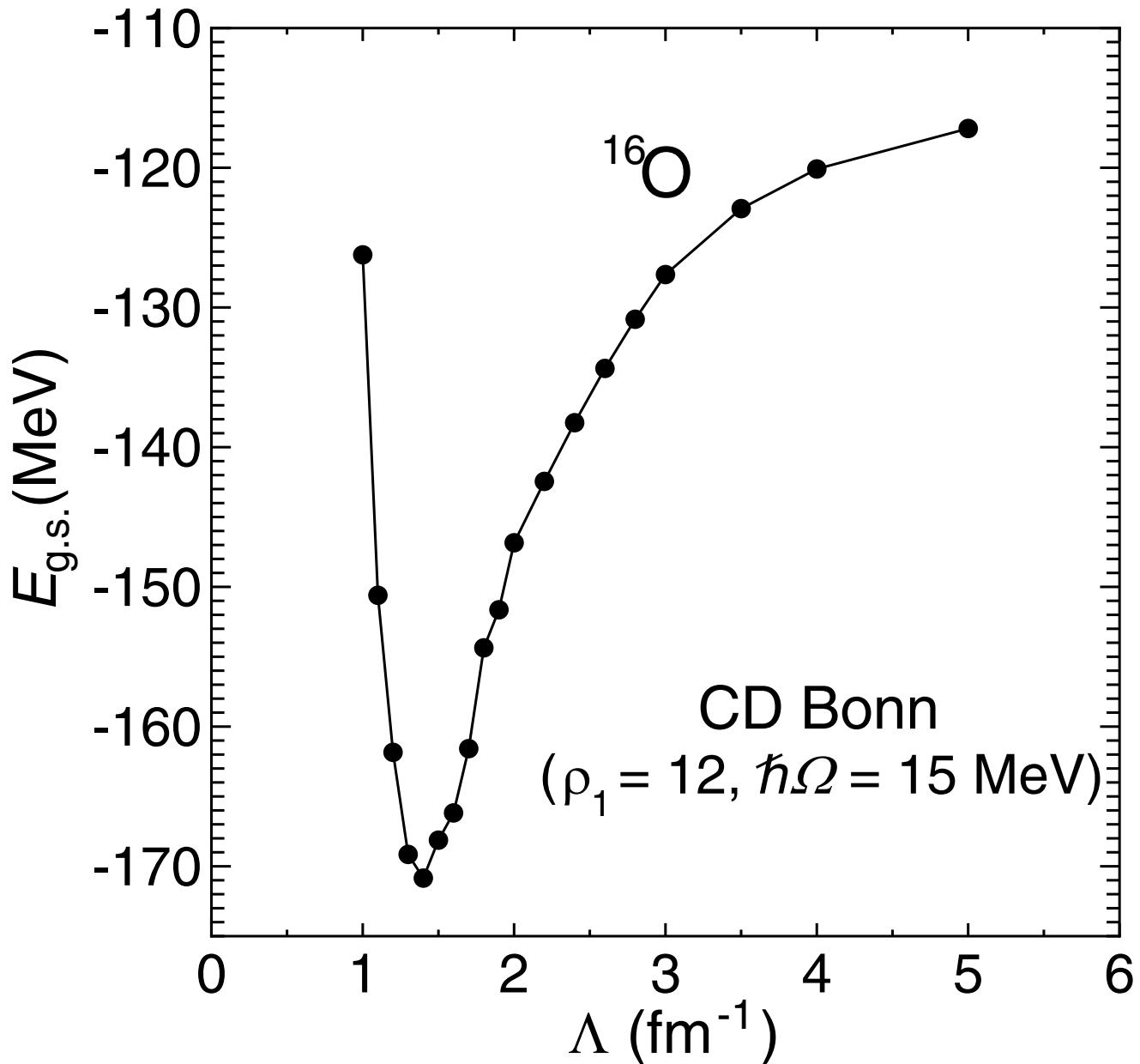
Deuteron properties

$\Lambda_{\text{cut}}(\text{fm}^{-1})$	CD Bonn	Nijm I
	BE(MeV)	P_D(%)
1.0	2.224576	1.212
2.0	2.224576	3.549
3.0	2.224576	4.546
4.0	2.224576	4.789
5.0	2.224576	4.830
6.0	2.224576	4.834
7.0	2.224576	4.833
quoted	2.224575	4.83

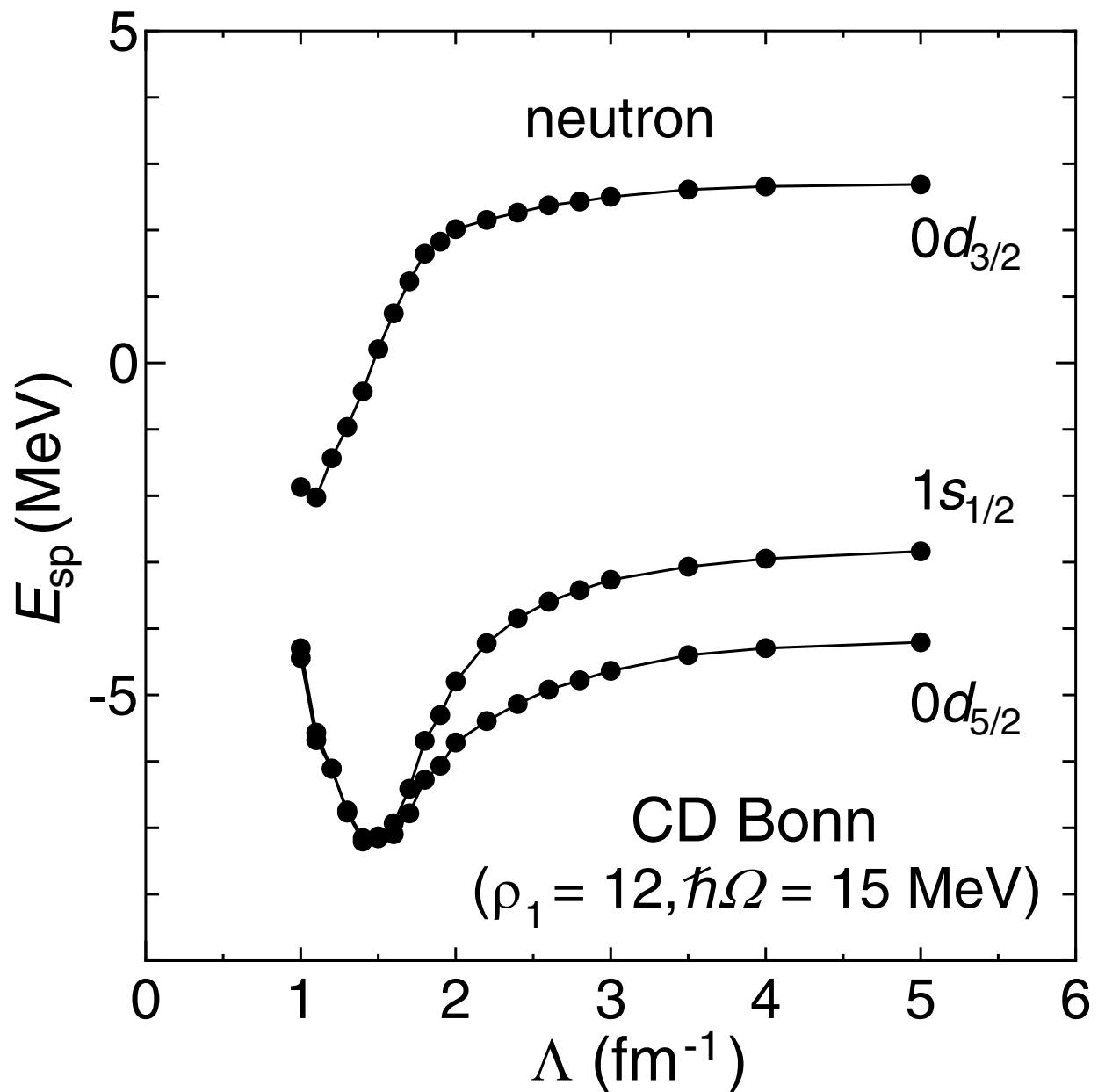




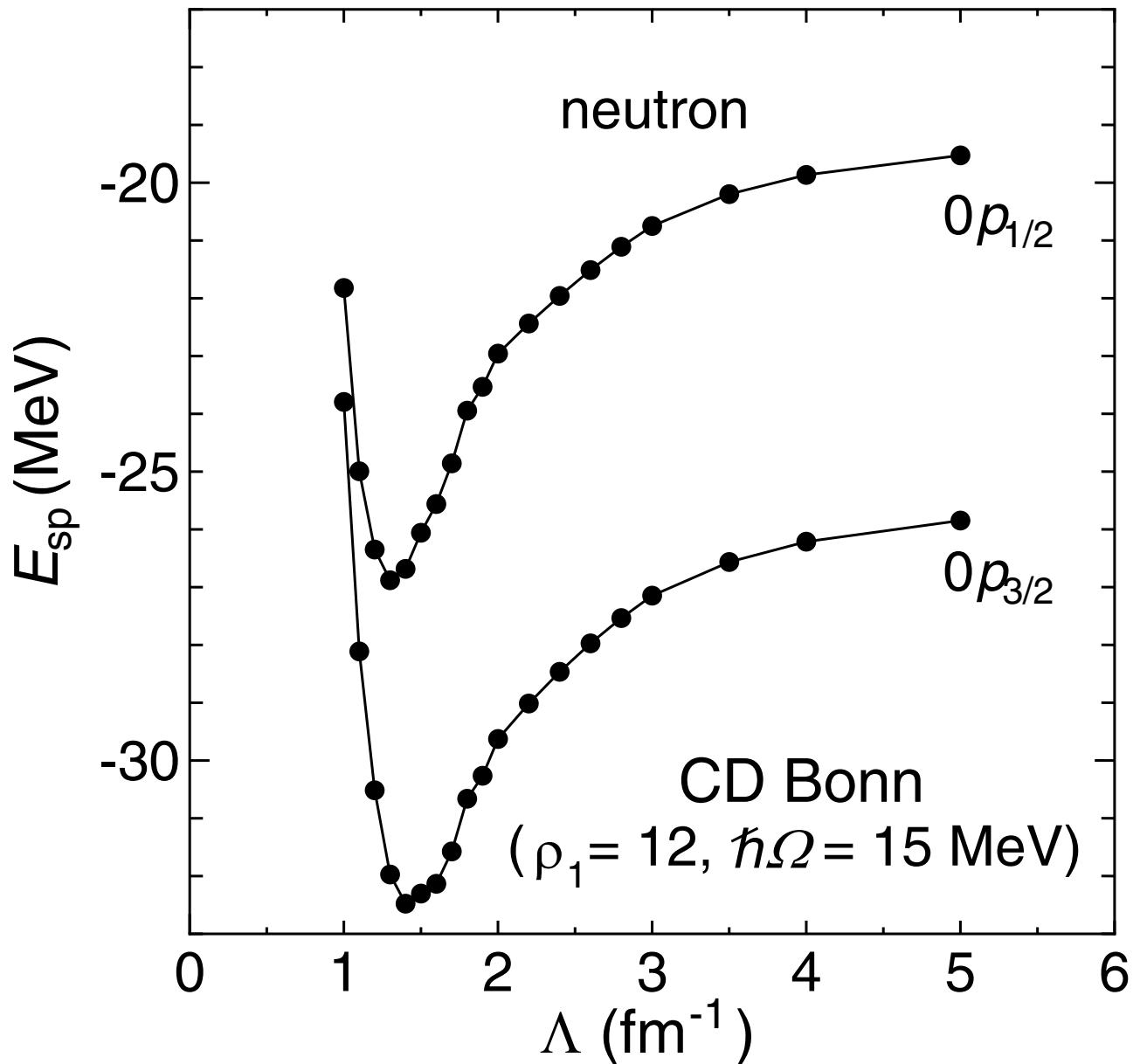
Ground-state energies of ^{16}O



Single-particle energies for particle states in ^{17}O



Single-particle energies for hole states in ^{16}O



Summary

- Low-momentum interactions were derived from realistic nucleon-nucleon interactions through a unitary transformation which has also been used in the unitary-model-operator approach (UMOA), the no-core shell model (NCSM), and the effective-interaction hyperspherical harmonics (EIHH).
- The low-momentum interactions obtained have high accuracy numerically, which was confirmed by the calculations of the deuteron binding energy.
- The low-momentum interactions were successfully applied to the Faddeev-Yakubovsky calculations for three- and four-nucleon systems.
- The calculated ground-state energies of the few-nucleon systems using the low-momentum interactions vary considerably at $\Lambda < 4 \sim 5 \text{ fm}^{-1}$, and there occurs the energy minimum at $\Lambda = 1 \sim 2 \text{ fm}^{-1}$.
- A similar tendency of the energy curve was obtained also in the calculations for ^{16}O . However, the magnitudes of relative spacings of single-particle levels are not so changed in the area $\Lambda > 2 \text{ fm}^{-1}$.
- The low-momentum interaction should be used with care especially in calculations of the total binding energy though the low-momentum interaction would be very useful in structure calculations as has been shown in shell-model calculations.