

Exotic molecular states in the $\alpha+{}^{6,8}\text{He}$
resonant scattering

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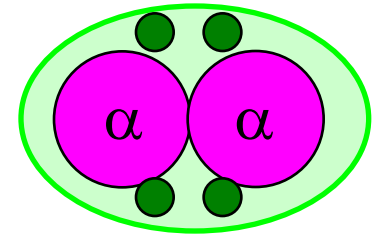
Nuclear clustering in N-rich nuclei

1. α -clustering in stable nuclear systems

$${}^8\text{Be} = 2\alpha, \quad {}^{12}\text{C} = 3\alpha, \quad {}^{16}\text{O} = \alpha + {}^{12}\text{C}, \quad {}^{20}\text{Ne} = \alpha + {}^{16}\text{O}$$

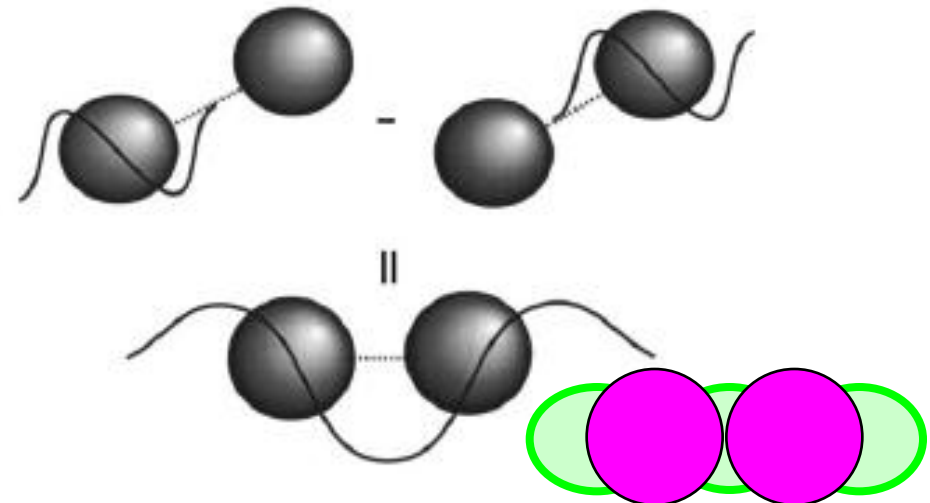
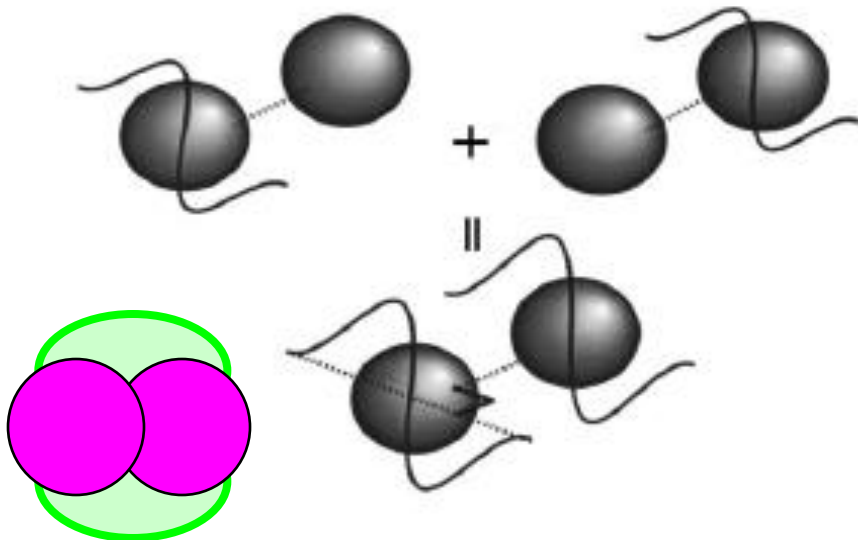
2. Clustering in N-rich systems : Clusters + XN

Example : Be isotopes = $2\alpha + \text{XN}$



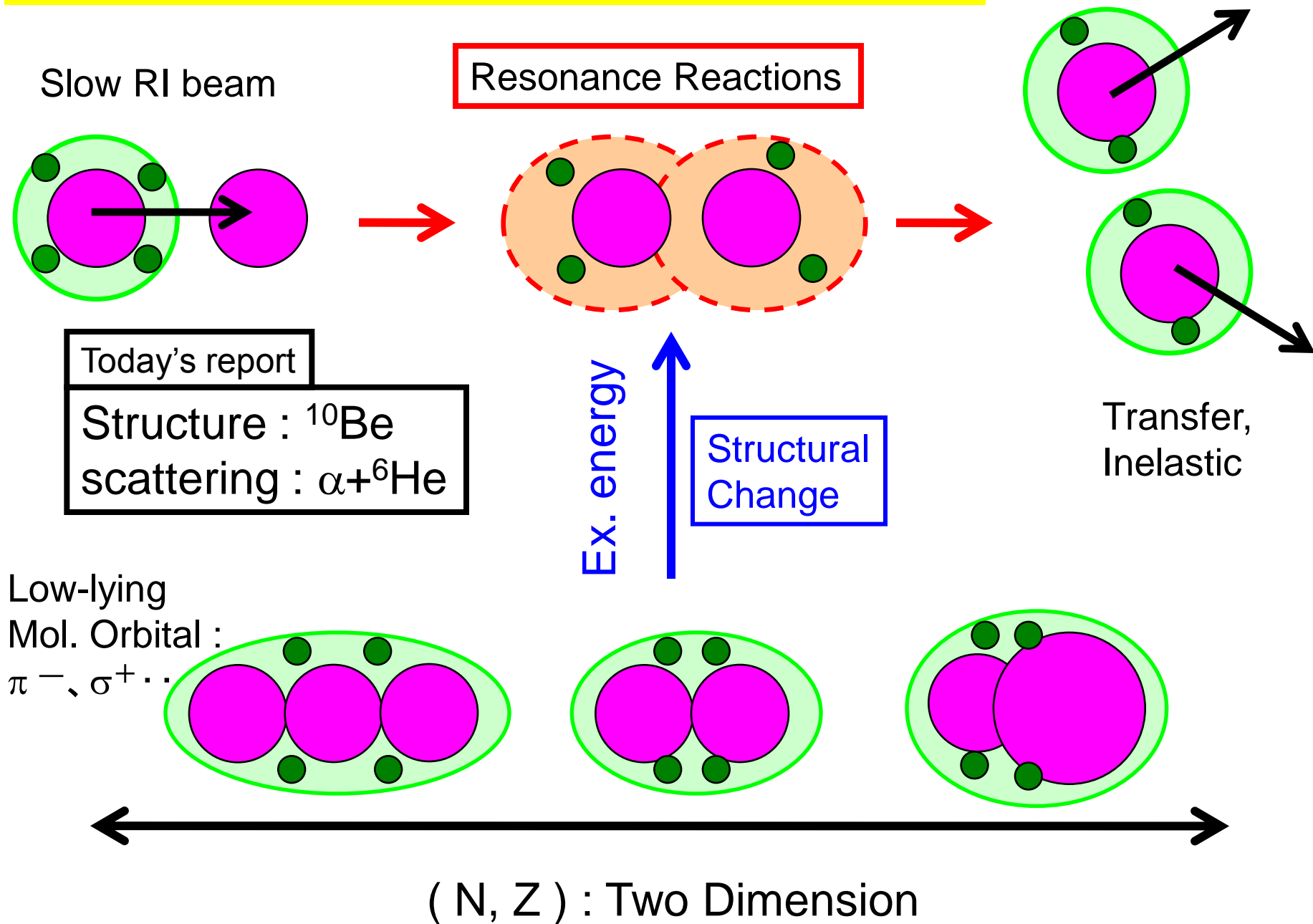
(a) π -orbit

(b) σ -orbit



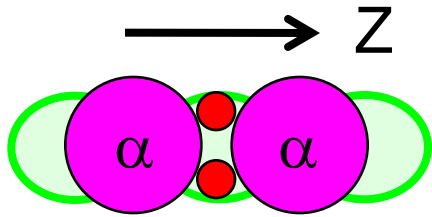
Molecular Orbital model : N. Itagaki et al., PRC61,62 (2000)

Studies on neutron-rich systems in 3-dimensional space



Formulation

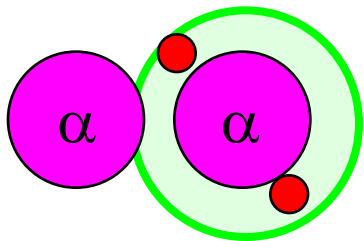
Linear Combination of Atomic Orbital (LCAO)



$$(\sigma^+)^2 = (P_z(L) - P_z(R))^2$$

$$= P_z(L) \cdot P_z(L) + P_z(R) \cdot P_z(R) - 2P_z(L) \cdot P_z(R)$$

$\begin{matrix} \text{6He} + \alpha & \alpha + \text{6He} & \text{5He} + \text{5He} \end{matrix}$



$$= P_x(R) \cdot P_x(R) + P_y(R) \cdot P_y(R) + P_z(R) \cdot P_z(R)$$

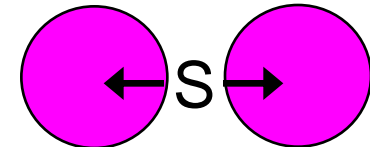
$\alpha + \text{6He}(0^+)$

Total wave function

$$\Psi = \sum_{\beta, S} \underline{C(\beta, S)} P_m(\mathbf{a}) \cdot P_n(\mathbf{b})$$

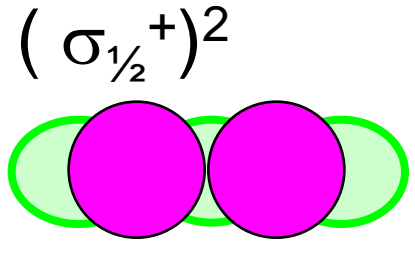
Variational PRM

$(m, n) = x, y, z$ $(\mathbf{a}, \mathbf{b}) = L, R$

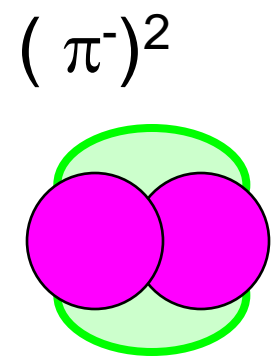


Adiabatic Energy surfaces : NN int. \Rightarrow Volkov No.2 + G3RS

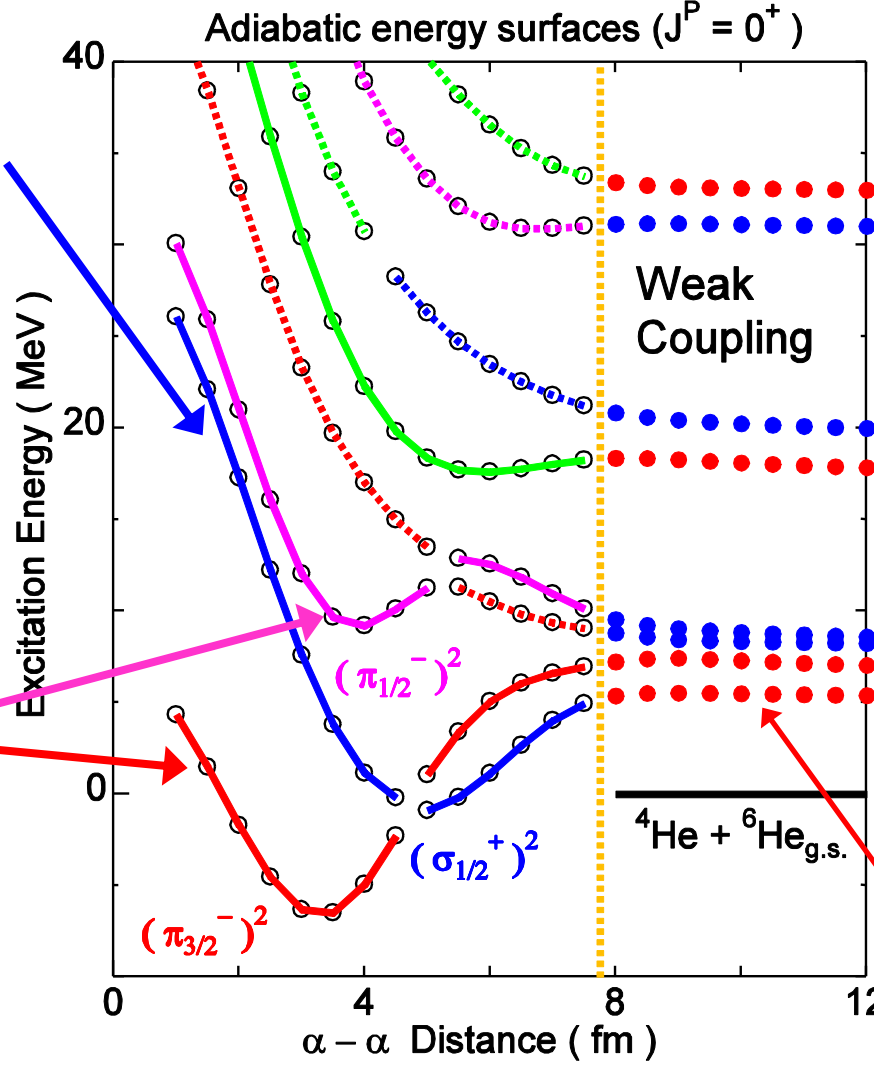
$$^{10}\text{Be} = \alpha + \alpha + n + n \quad (J^\pi = 0^+)$$



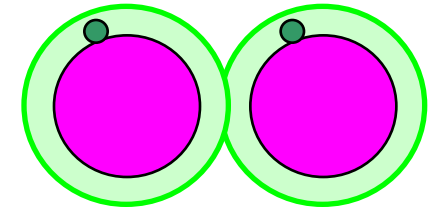
sd-orbital



0p-orbital

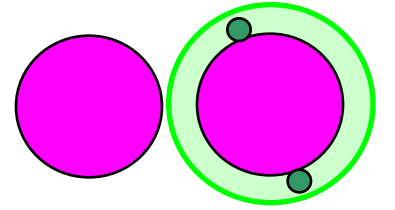


Blue Dots



$^5\text{He}(I_1) + ^5\text{He}(I_2) L$

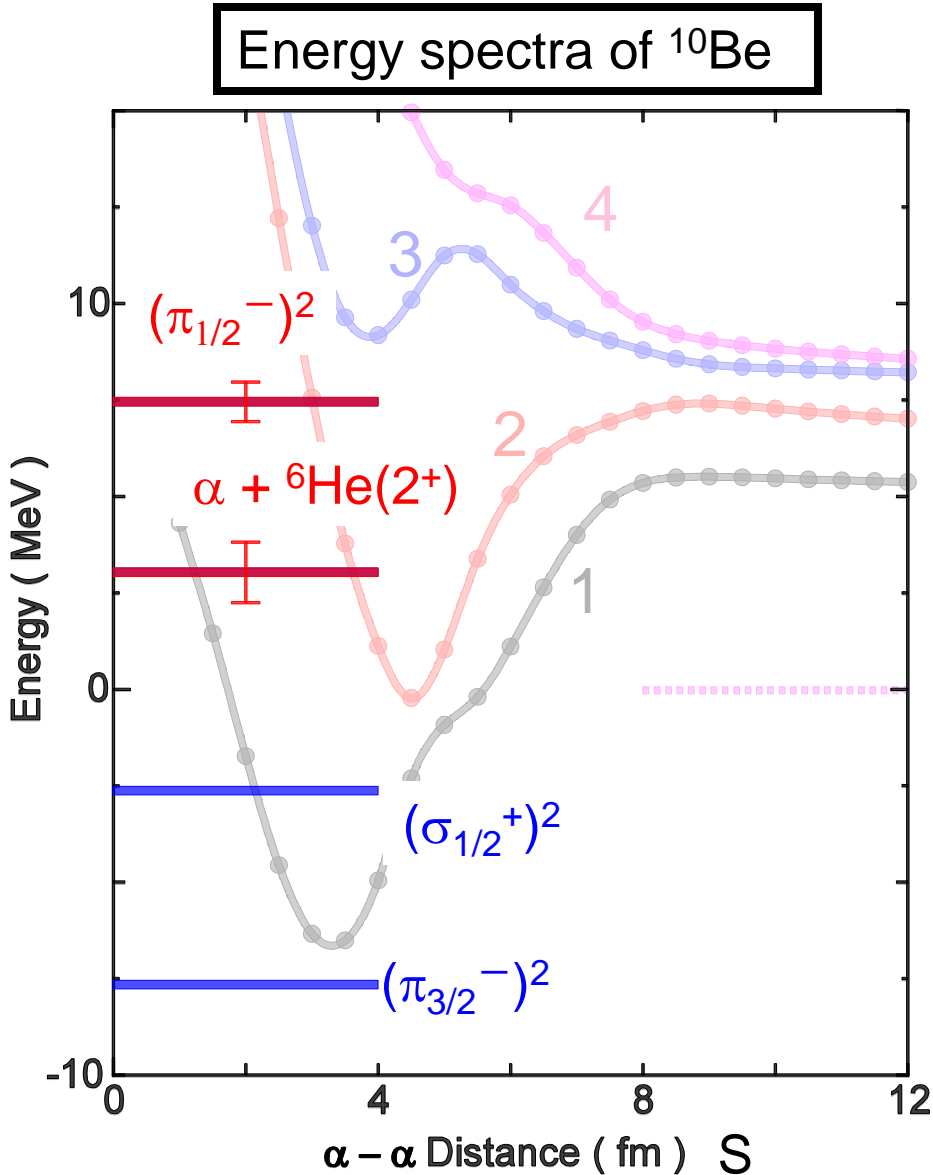
Red Dots



$\alpha + ^6\text{He}(I) L$

$\alpha + ^6\text{He}(0_1^+)$

Coupled channels with the adiabatic states



Adiabatic basis

$$\Psi = \sum_{k,S} f(k,S) \Phi^{AD}(k,S)$$

$k = 1, 2, 3, 4, \dots$

$$H = \sum_i t_i + \sum_{i>j} v_{i,j}$$

$$(H - E)\Psi = 0$$

Bound,
Resonant
B.C.

Scattering
B.C.

Energy spectra

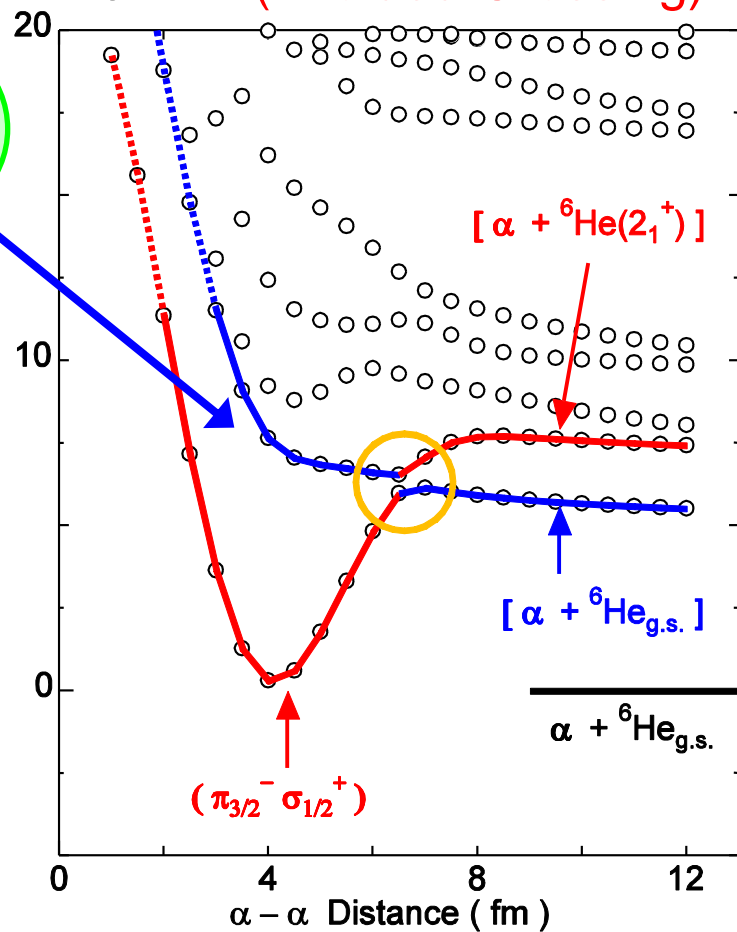
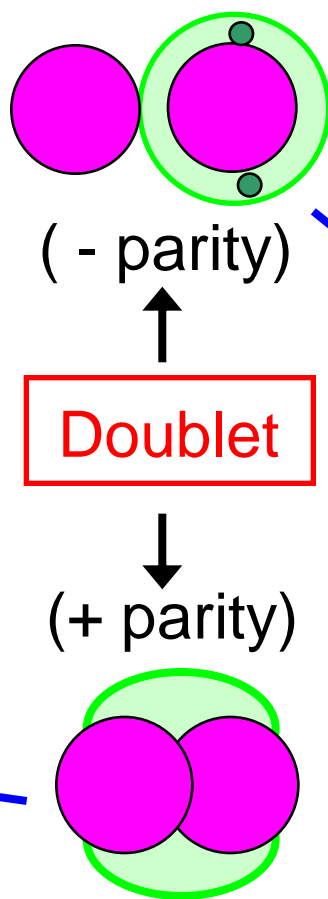
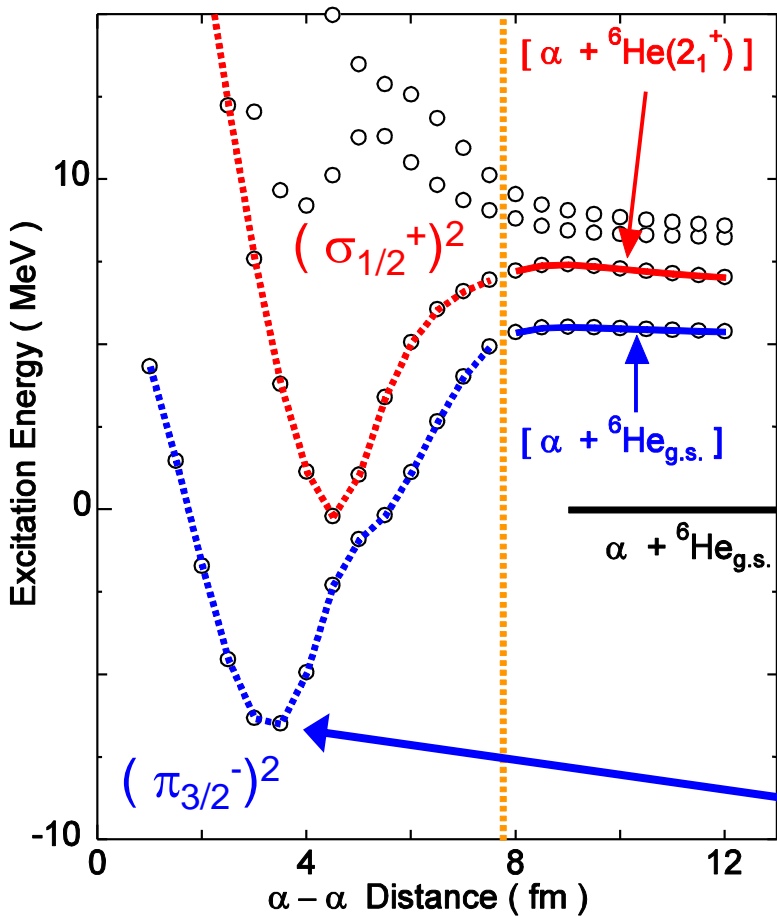
$\alpha + {}^6\text{He}$ scattering

Parity doublet formation in adiabatic energy surfaces

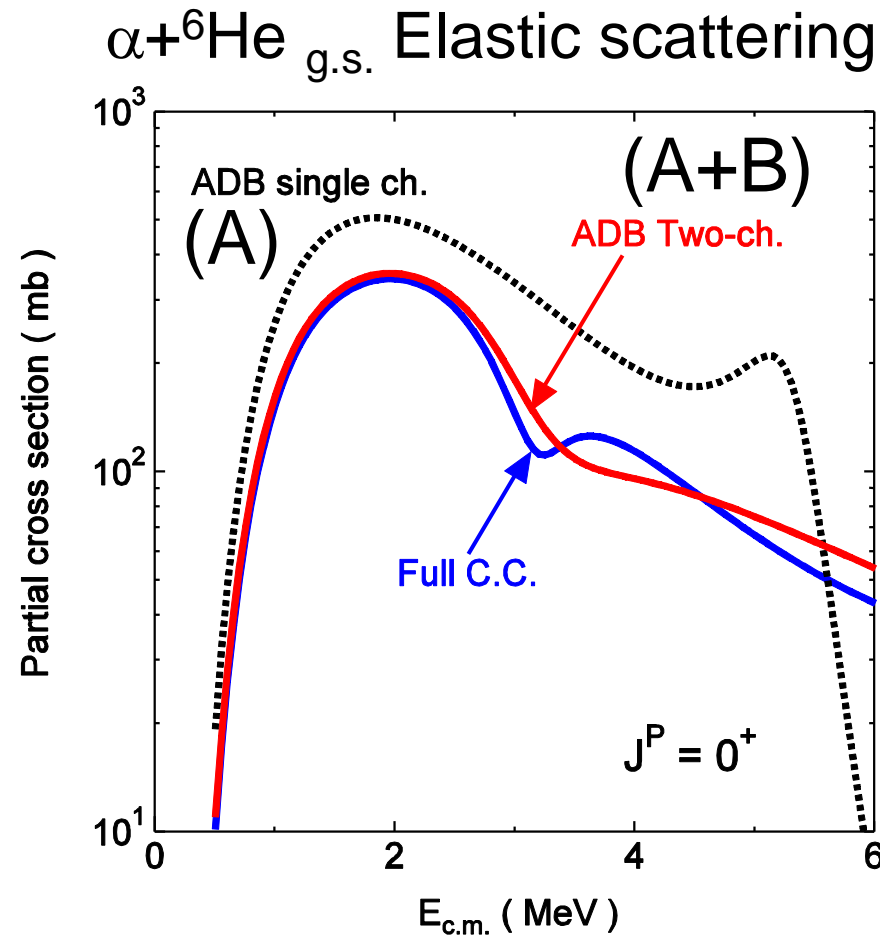
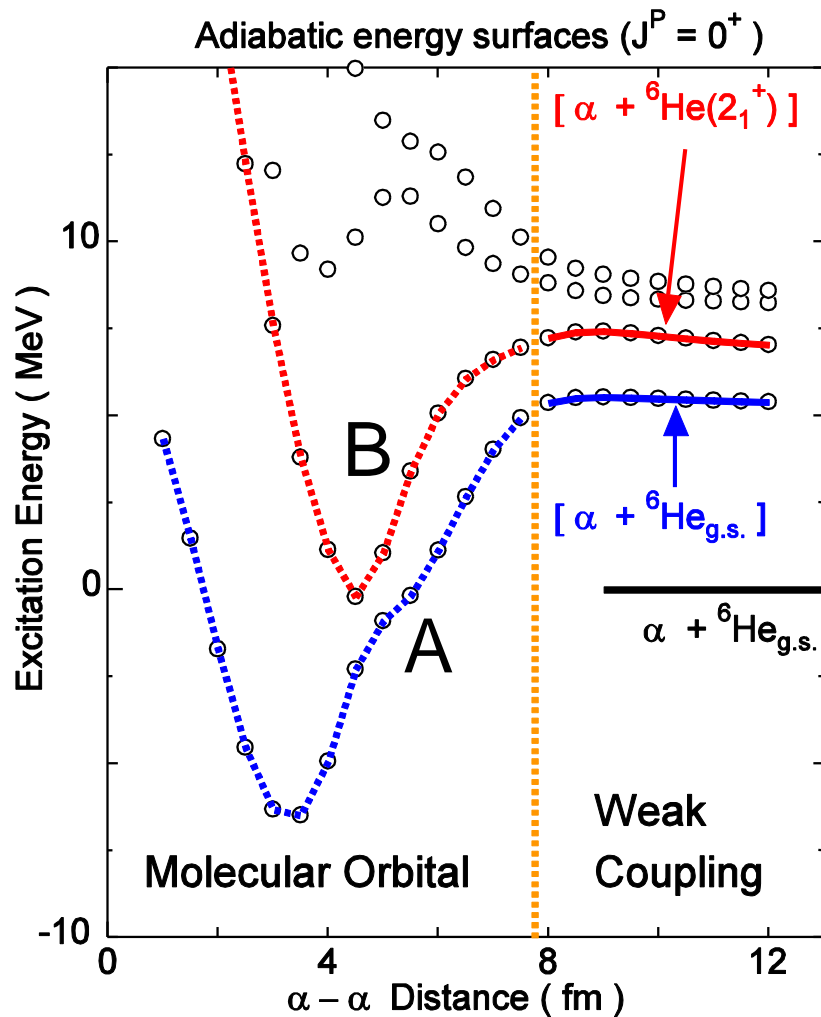
P-Doublet : $\Psi^{(\pm)} = \left\{ \begin{array}{c} \text{[Diagram 1]} \\ \pm \\ \text{[Diagram 2]} \end{array} \right\}$

$J^\pi = 0^+$ (Long range coupling)

$J^\pi = 1^-$ (Avoided Crossing)

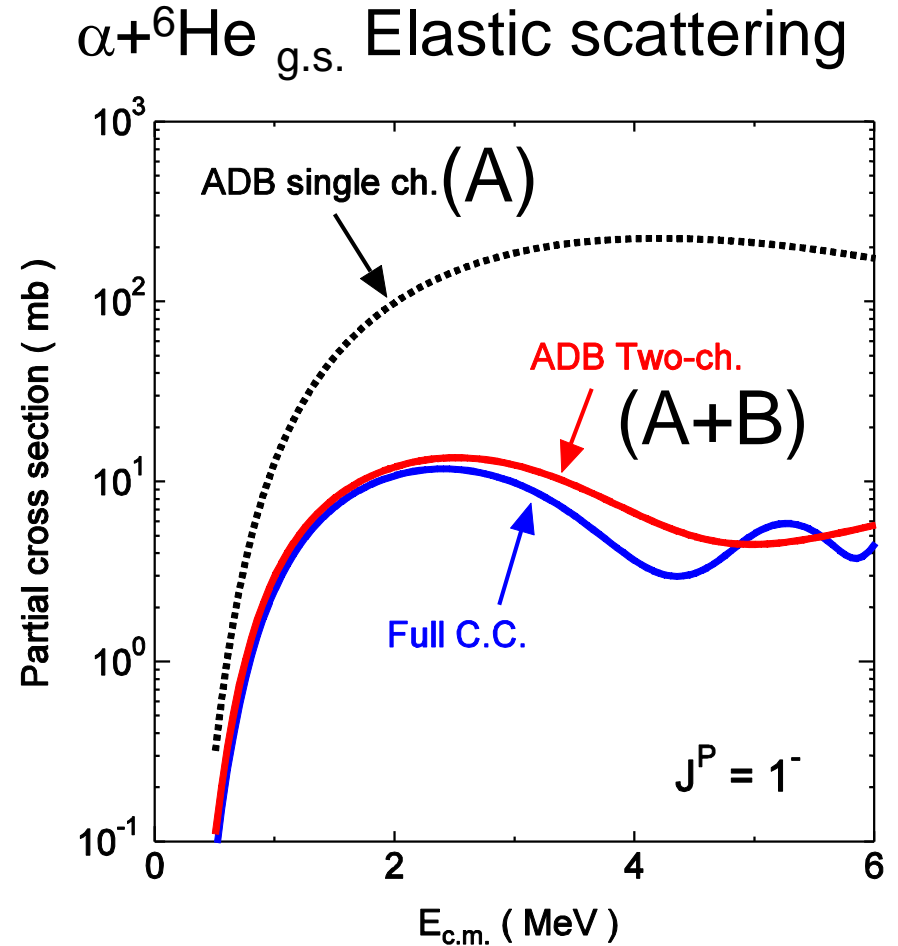
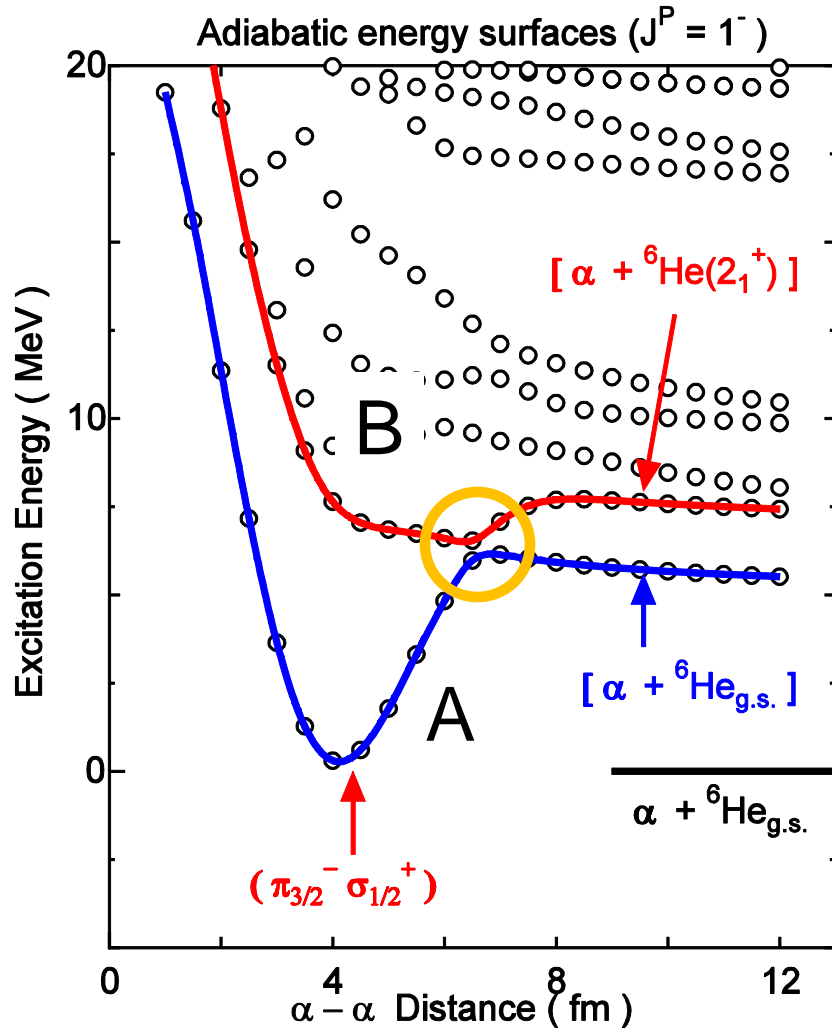


Coupled channels in $\alpha + {}^6\text{He}$ elastic scattering ($J^\pi = 0^+$)



Adiabatic approximation is good.

Coupled channels in $\alpha + {}^6\text{He}$ elastic scattering ($J^\pi = 1^-$)



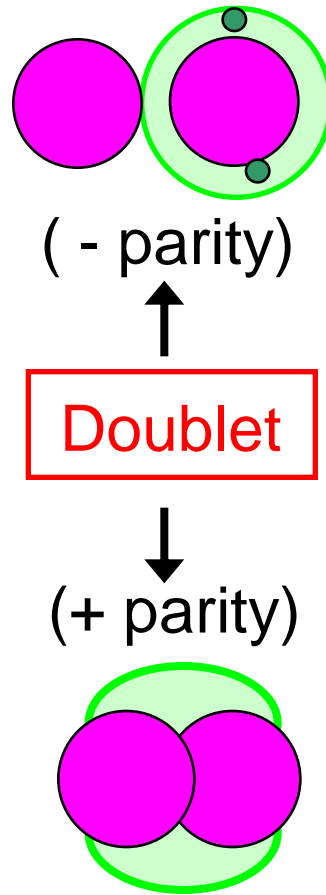
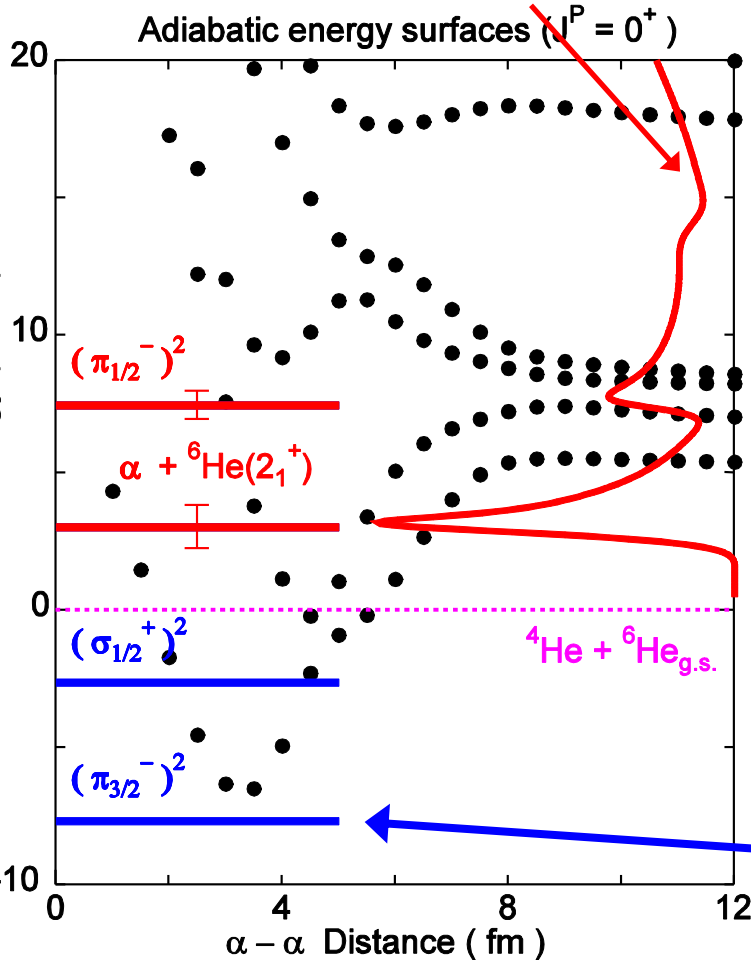
Non-adiabatic transition strongly occurs.

Enhancements in $\alpha+{}^6\text{He}$ inelastic scattering

(B. Imanisi et al., Phys. Rep. 155)

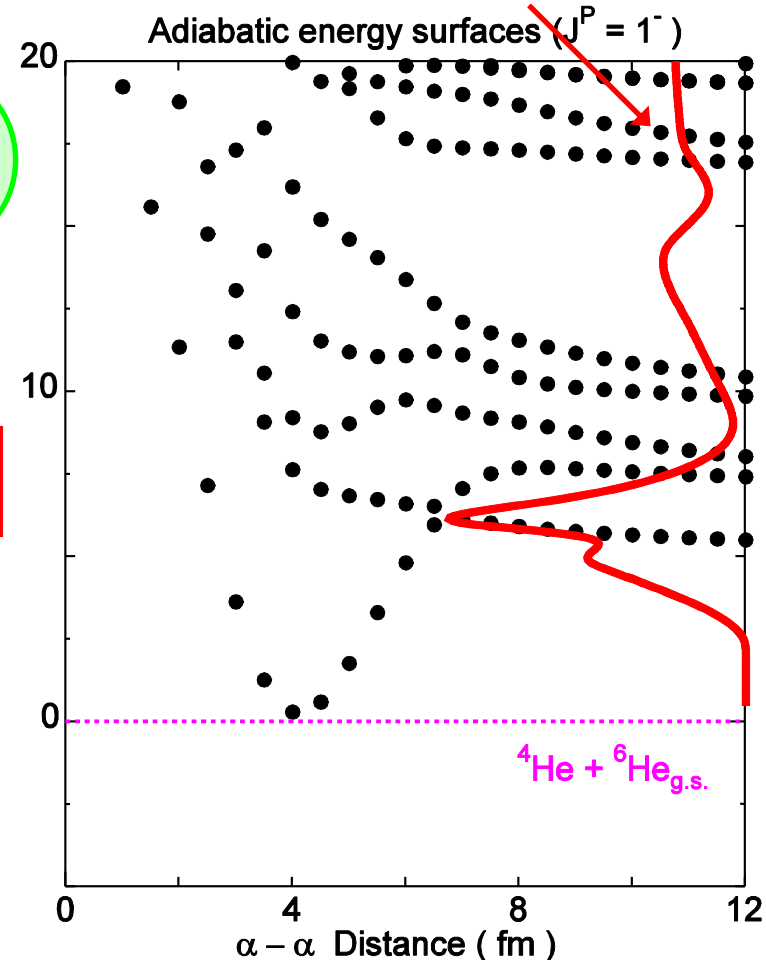
$J^\pi = 0^+$ **Resonance Poles**

$\alpha+{}^6\text{He}(0_1^+) \rightarrow \alpha+{}^6\text{He}(2_1^+)$



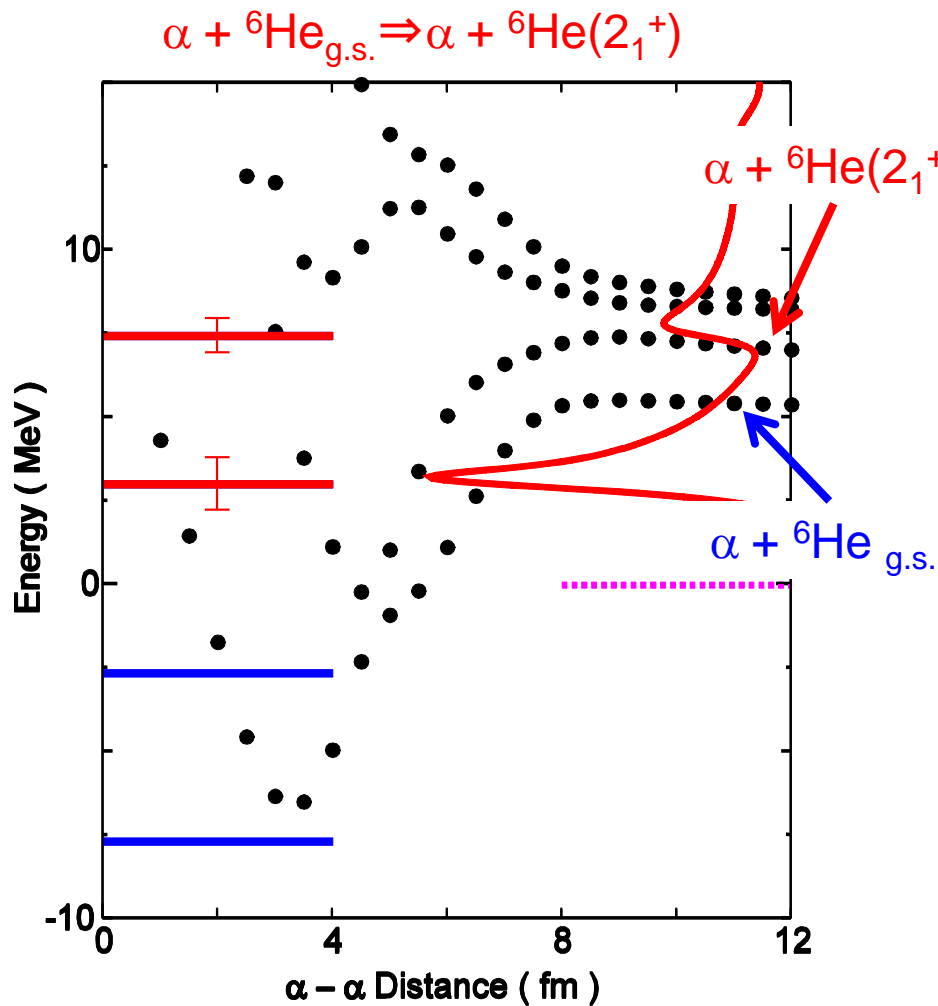
$J^\pi = 1^-$ **L-Z level crossing**

$\alpha+{}^6\text{He}(0_1^+) \rightarrow \alpha+{}^6\text{He}(2_1^+)$

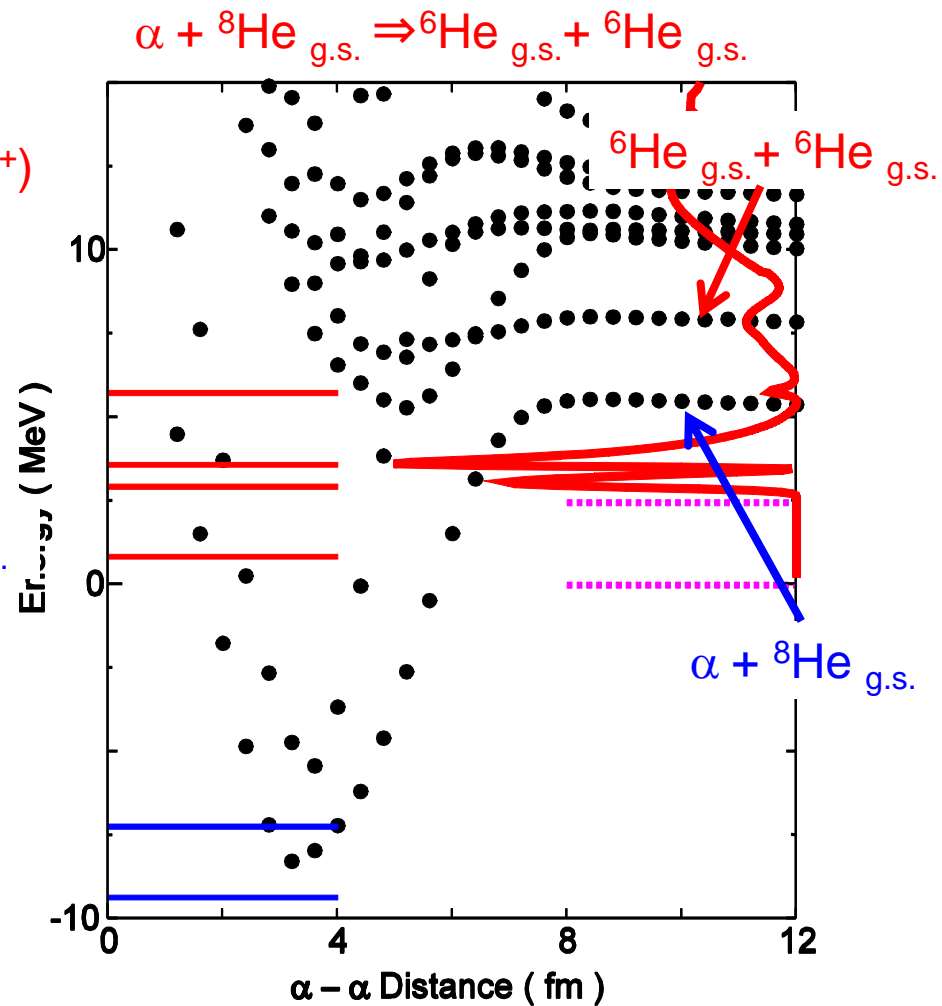


Comparison between $\alpha+{}^6\text{He}$ and $\alpha+{}^8\text{He}$ scattering

${}^{10}\text{Be}(0^+)$: Res. in Inel. Scatt.



${}^{12}\text{Be}(0^+)$: Res. in Transfer ch.



Summary

1. Studies on N-rich systems in (N,Z,E) space is very interesting.
2. Unified description of structures and reaction becomes quite important.
3. We show some enhancements in the $\alpha+{}^6\text{He}$ and $\alpha+{}^8\text{He}$ scattering.

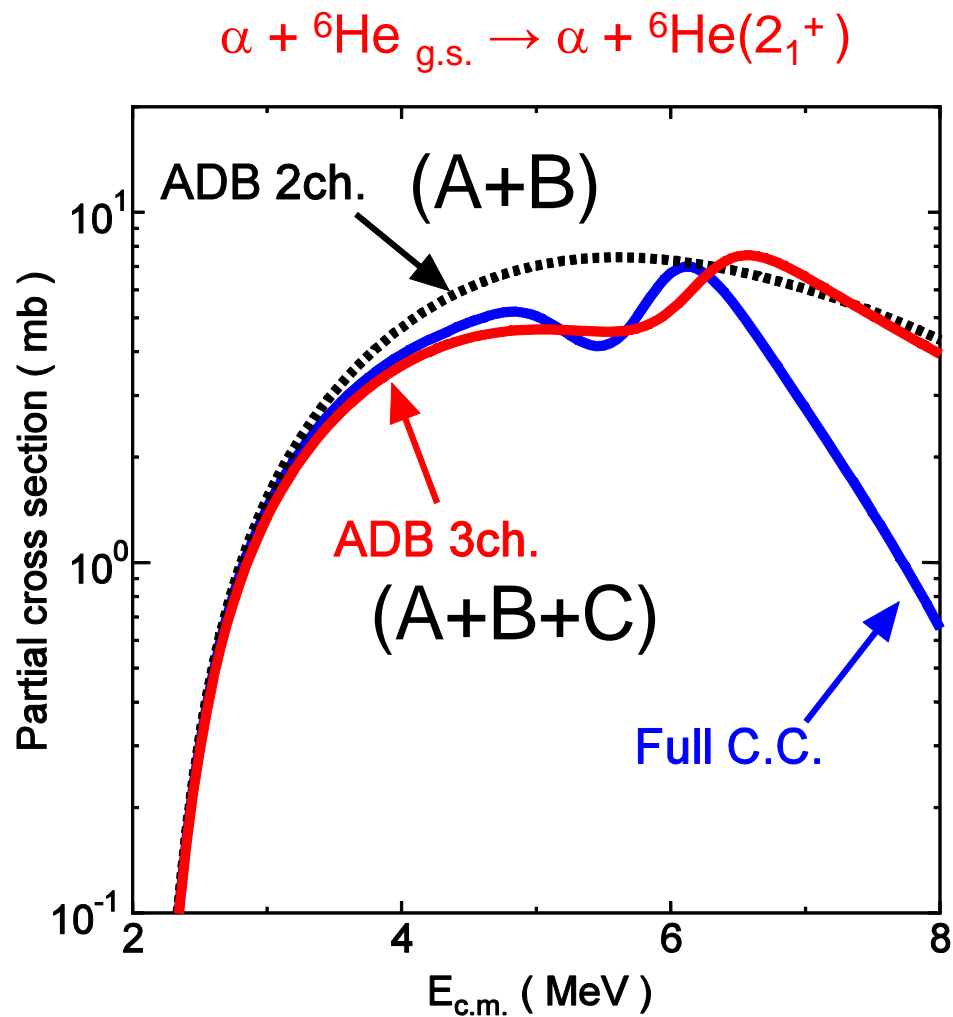
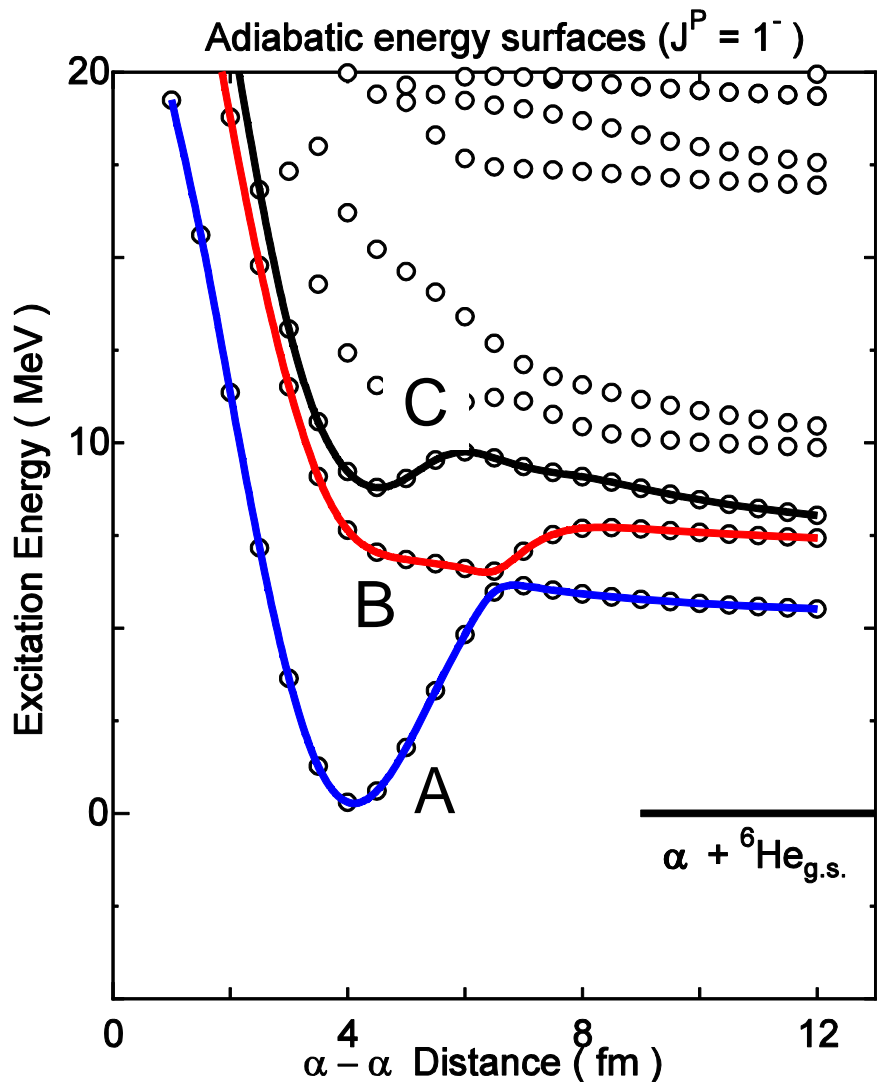
Results of ${}^{10}\text{Be}$

1. Molecular and atomic states coexist in this system.
2. Reaction process is different between the positive parity and the negative one.
3. L-Z transition is predicted in connection to the Parity doublet.

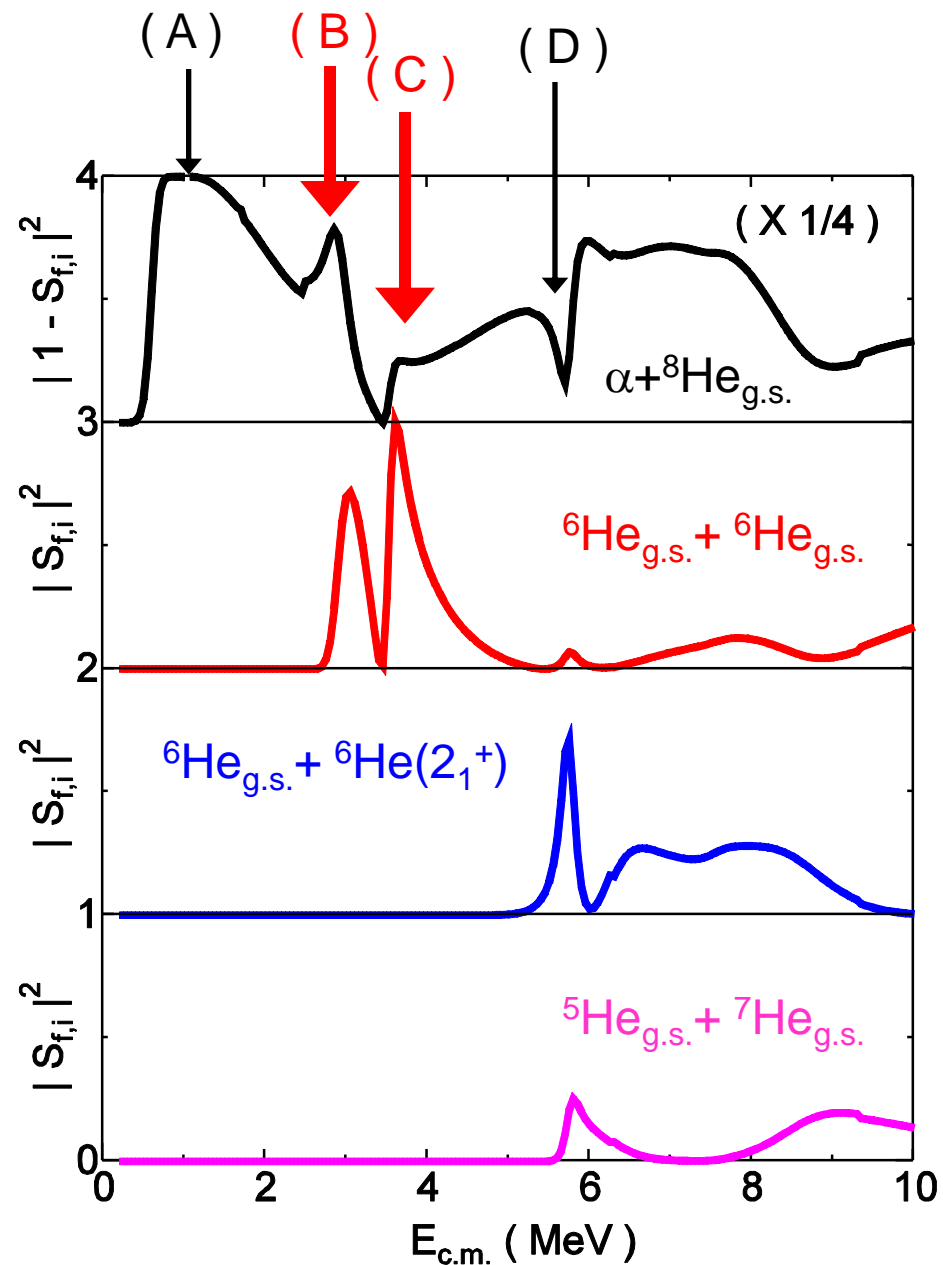
Results of ${}^{12}\text{Be}$

1. Resonances will be observed in two neutron transfers.
2. Exotic structures are excited. (I will report at YITP post symposium.)

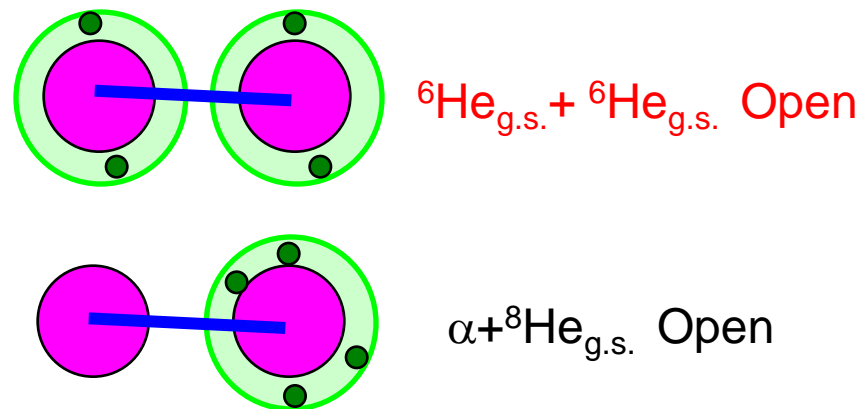
Coupled channels with the adiabatic states



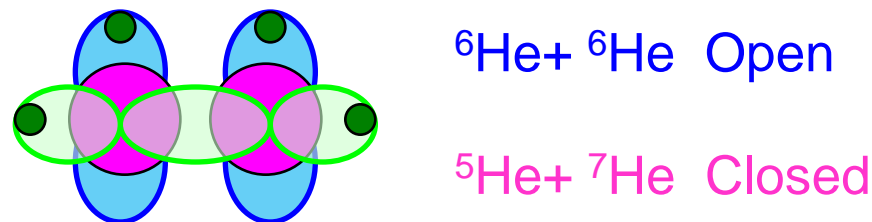
Resonances in $\alpha + {}^8\text{He} \Rightarrow {}^x\text{He} + {}^y\text{He}$



(B) Two nucleon transferred state

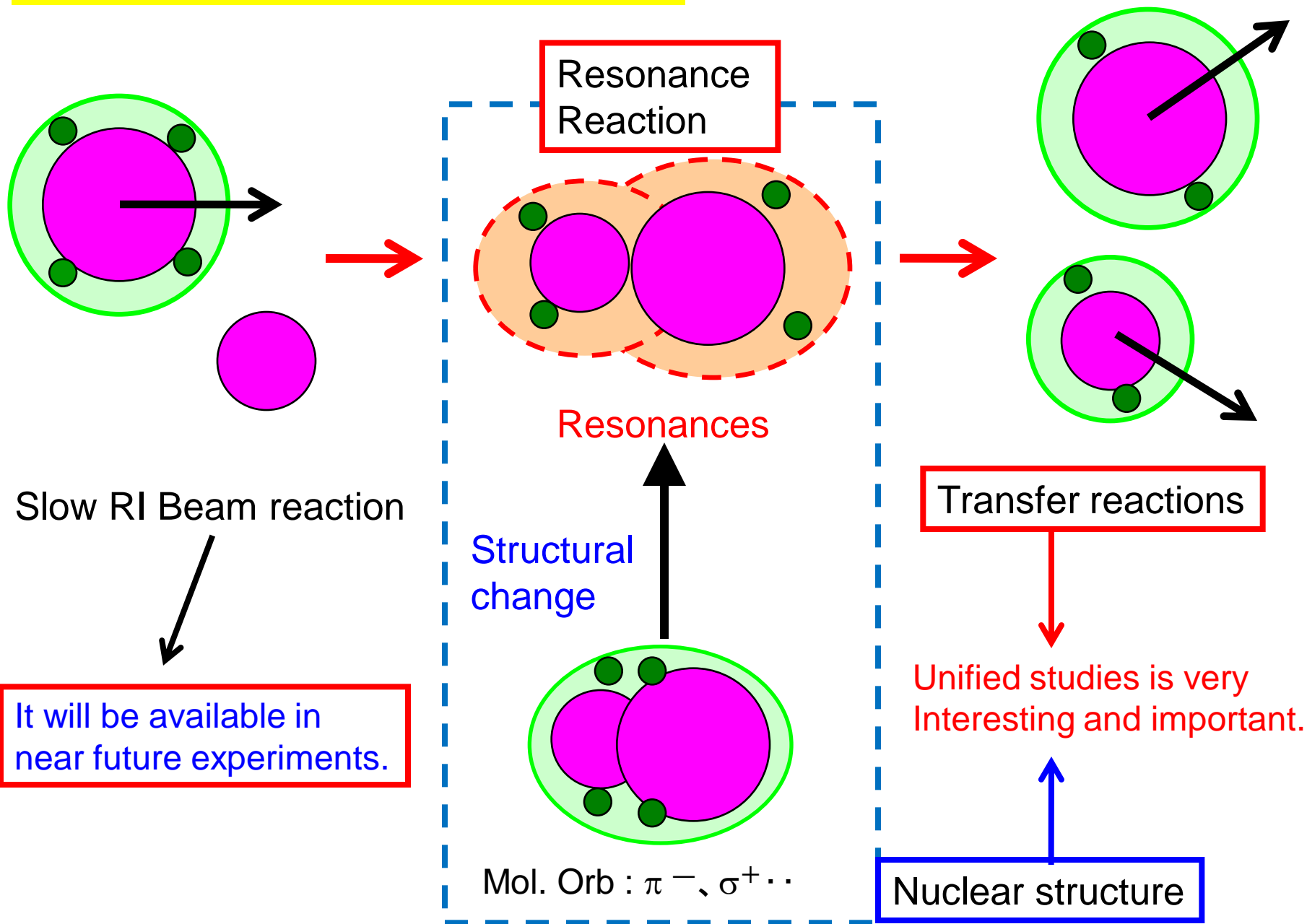


(C) Atom-Molecule hybrid state

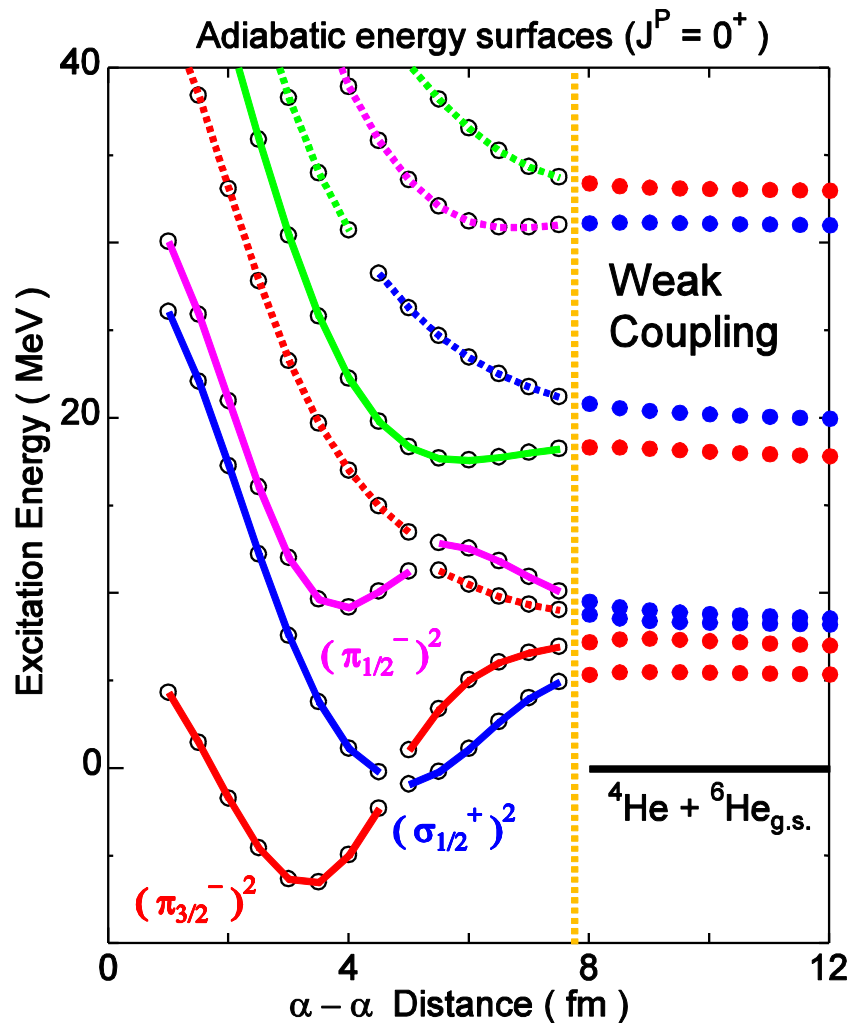


Decay scheme depends on the intrinsic structure

Unified studies in neutron-rich system



Coupled channels with the adiabatic states



$$u^{rel}(\alpha, R, S) \varphi^{in}(\alpha, \xi_\alpha)$$

$$\Phi^{AD}(i, S) = \sum_{\beta} C_{\beta}(i) u^{rel}(\beta, R_{\beta}, S) \varphi^{in}(\beta, \xi_{\beta})$$

$$u^{rel}(\beta, R_{\beta}, S) \propto \exp[-v(R_{\beta} - S)^2]$$

$$u^{rel}(\alpha, R_{\alpha}, S) \propto \exp[-v(R_{\alpha} - S)^2] \quad (R \leq R_c)$$

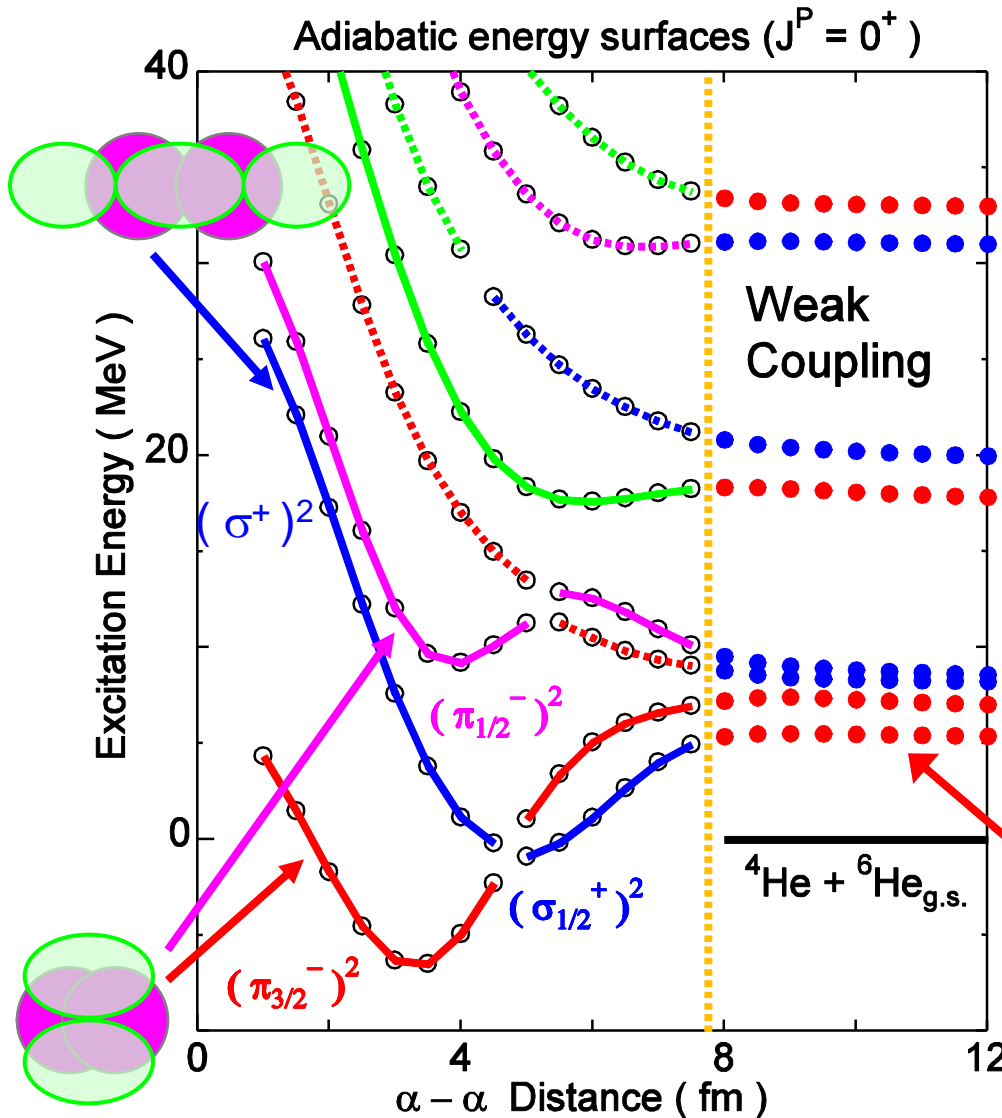
$$u_{L\alpha}^{(-)}(R_{\alpha}) - \varepsilon u_{L\alpha}^{(+)}(R_{\alpha}) \quad (R \geq R_c)$$

$$\Psi = \sum_{i,S} f(i, S) \Phi^{AD}(i, S)$$

$$(H - E)\Psi = 0$$

Generalized Two-center cluster model (GTCCM) : PLB588

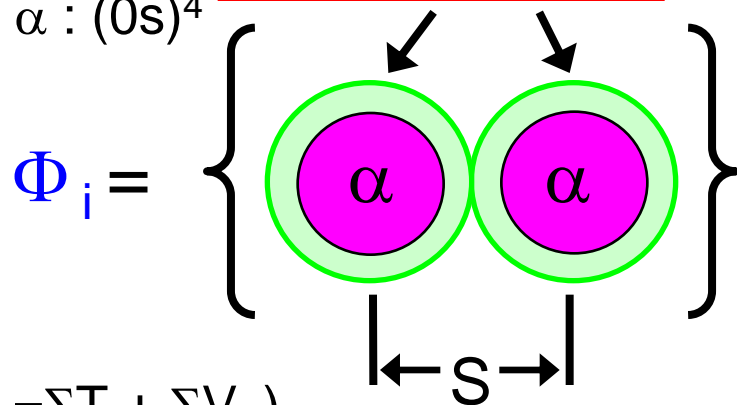
$$^{10}\text{Be} = \alpha + \alpha + \text{N} + \text{N} : J^\pi = 0^+$$



Basis : Atomic orbital (A.O.)

$0p_i$ A.O. ($i=x,y,z$)

$\alpha : (0s)^4$



$$(H = \sum T_i + \sum V_{ij})$$

Total wave function

$$\Psi^{J\pi} = P^{J\pi} A \sum C_i \Phi_i$$

J^π Proj. Anti-sym.

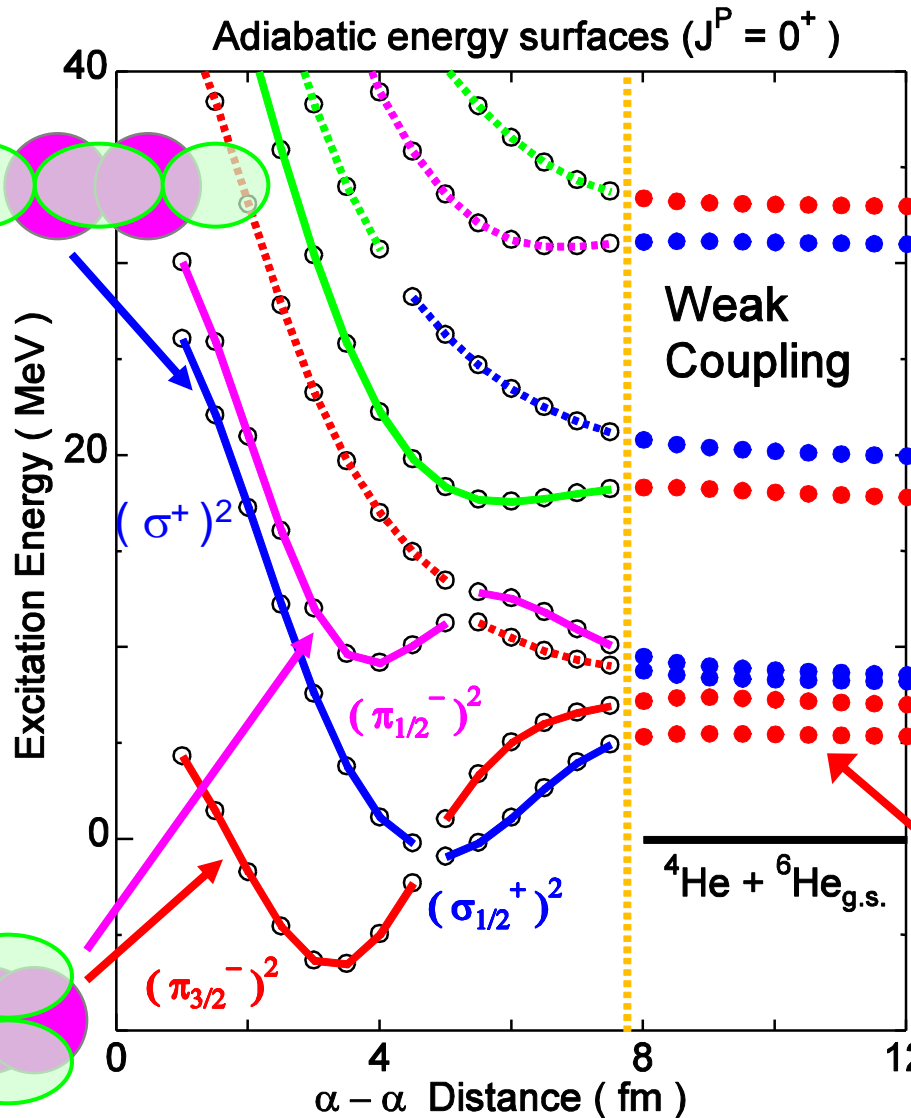
$\alpha + ^6\text{He}(0_1^+)$

● Red Dots : $[\alpha + ^6\text{He}(1)]$ LJ

● Blue Dots : $[^5\text{He}(I_1) + ^5\text{He}(I_2)]$ LJ

Generalized Two-center cluster model (GTCCM) : PLB588

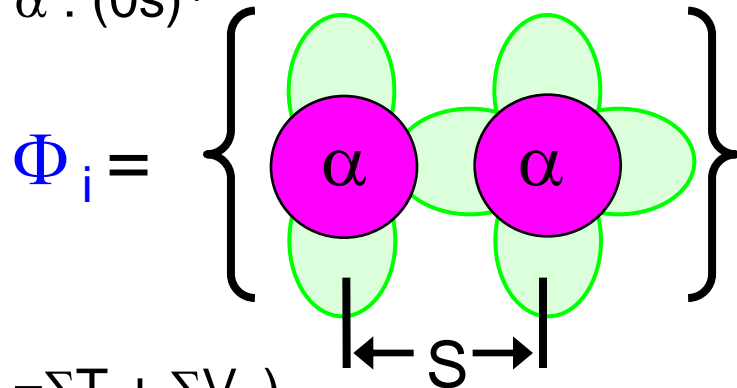
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$$(H = \sum T_i + \sum V_{ij})$$

Total wave function

$$\Psi^{J\pi} = P^{J\pi} \mathcal{A} \sum C_i \Phi_i$$

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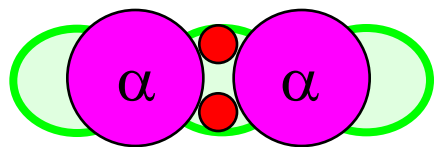
$\alpha + ^6\text{He}(0_1^+)$

● Red Dots : $[\alpha + ^6\text{He}(1)]$ LJ

● Blue Dots : $[^5\text{He}(I_1) + ^5\text{He}(I_2)]$ LJ

Formulation

Linear Combination of Atomic Orbital (LCAO)



$$(\sigma^+)^2 = (p_z(L) - p_z(R))^2$$

