

Studies of nuclear halo formation with (e,e',p) , (e,e',d) reactions

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Motivation

- (e,e',p) , (e,e',d) experiments may be used to study the halo formation in exotic nuclei
- Necessity to develop proposals for new experiments
- Key factor: the interplay between nuclear structure and scattering experiments
- New ways towards modern physics
- Outlook

Topics

- Review of the theory for scattered electron in continuum
- Proton and deuteron knock-out reaction
- Microscopic correlation in the dynamic of even nuclei
- Example: ${}^6\text{Li}$
- Momentum distribution in exotic nuclei
- Effect of the halo in the scattering cross-section
- Correlated three-particle system

Cross-section for electron scattering in continuum

$$\frac{d^2 \sigma}{d\Omega dE_2} = \frac{\sigma_{Mott}}{M} \left(W_2 + 2W_1 \tan^2 \frac{\theta}{2} \right)$$

where

$$W_1 = 2\pi M F_T^2$$

$$W_2 = 4\pi M \left(\frac{\Delta^2}{q^2} \right) \left[\left(\frac{\Delta^2}{q^2} \right) F_C^2 + \frac{1}{2} F_T^2 \right]$$

with

$$\sigma_{Mott} = \frac{\alpha^2}{4k_1^2} \cdot \frac{\cos^2 \frac{\theta}{2}}{\sin^4 \frac{\theta}{2}} \cdot \frac{1}{M}$$

$$\vec{q} = \vec{k}_1 - \vec{k}_2$$

$$\Delta^2 = -\omega^2 + q^2$$

$$\omega = k_1^0 - k_2^0 \equiv E_1 - E_2$$

→

q momentum transfer

ω energy transfer

M mass of the target nucleus

F_T transversal form factor

F_C longitudinal form factor

Inverse relations and dimensionless response functions

Inverse relations

$$F_T^2 = \frac{2W_1}{4\pi M}$$

$$F_C^2 = \frac{1}{4\pi M} \left(\frac{\vec{q}^2}{\Delta^2} \right) \left[\left(\frac{\vec{q}^2}{\Delta^2} \right) W_2 - W_1 \right]$$

The dimensionless functions are defined by:

$$S_T = MF_T^2 \quad S_T \equiv S_{Transversal}$$

$$S_L = MF_C^2 \quad S_L \equiv S_{Longitudinal}$$

Scattering cross-section and nuclear model

$$\frac{d^2\sigma}{d\Omega_2 dE_2} = \sigma_M K \left(A + B \tan^2 \frac{\theta}{2} \right)$$

with: $K \equiv |f(\Delta^2)|^2 \left(\frac{V}{4\pi^3} \right)$

$$\left(\frac{V}{4\pi^3} \right) n(\vec{p}) \xrightarrow{\text{to be replaced by}} |\Psi_{DCM}(\vec{p})|^2$$

Taken from the Fermi gas model

Totally correlated n-particle momentum distribution

$$F_T^2 = \frac{K}{4\pi M} B$$

$$F_C^2 = \frac{1}{4\pi} \left(\frac{\vec{q}^2}{\Delta^2} \right)^2 \left[A - \frac{1}{2} B \right]$$

Calculation of A and B

For the computation of A and B we use Ueberall:
 „Electron scattering from complex nuclei“

$$\frac{1}{2} A(q, \omega) = I_0 - \frac{\omega}{m} \left(I_0 + \frac{2I_1}{q} \right) + \frac{1}{m^2} \left[\left(\frac{G}{2} - I_2 \right) \left(1 - \frac{\omega^2}{q^2} \right) + I_2 \right] - \frac{\omega^2}{m^2} \left(2.63I_0 - \frac{I_1}{q} \right) + 1.17I_0 \frac{q^2}{m^2}$$

$$\frac{1}{2} B(q, \omega) = \frac{1}{m^2} \left[(G + 5.77q^2 I_0) - I_2 \right]$$

To calculate the integrals $I_0, I_1, I_2 \dots$ we are not using the momentum distribution of the Fermi model but the distribution calculated in the BDCM model.

Momentum integrals

$$I_l = \int d^3 p |\Psi_{DCM}(p)|^2 \left[1 - \left| \Psi_{DCM} \left(\left| \vec{p} + \vec{q} \right| \right) \right|^2 \right] (p \cos \phi)^l \delta \left(\omega - \varepsilon \left(\left| \vec{p} + \vec{q} \right| \right) + \varepsilon(p) \right)$$

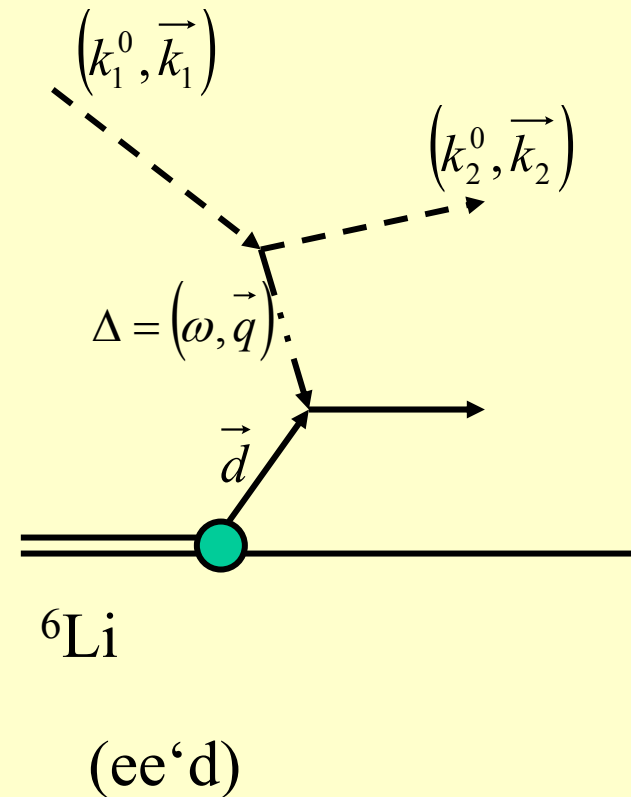
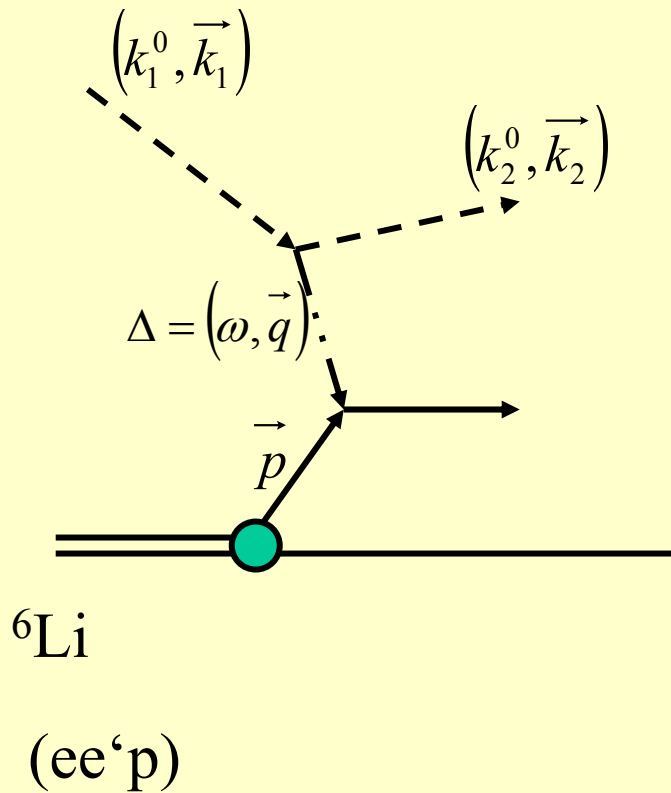
$$G = \int d^3 p |\Psi_{DCM}(p)|^2 \left[1 - \left| \Psi_{DCM} \left(\left| \vec{p} + \vec{q} \right| \right) \right|^2 \right] p^2 \delta \left(\omega - \varepsilon \left(\left| \vec{p} + \vec{q} \right| \right) + \varepsilon(p) \right)$$

$$\phi = \text{angle}(\vec{p}, \vec{q})$$

$\varepsilon(p)$ the energy of two particle excitation as a function of its momentum

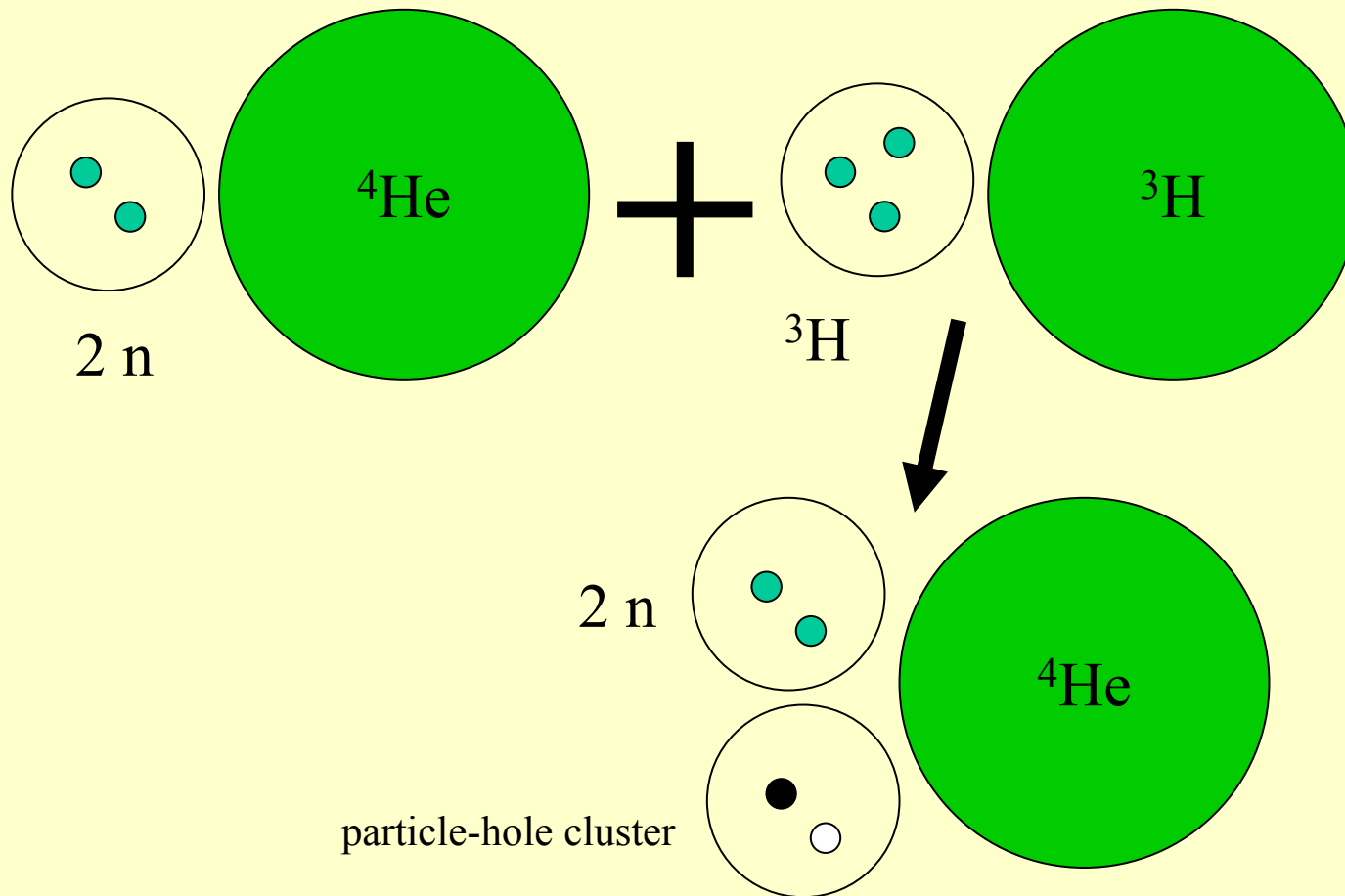
An effective mass m^* for the bound nucleons has been introduced.

Electron scattering on ${}^6\text{Li}$



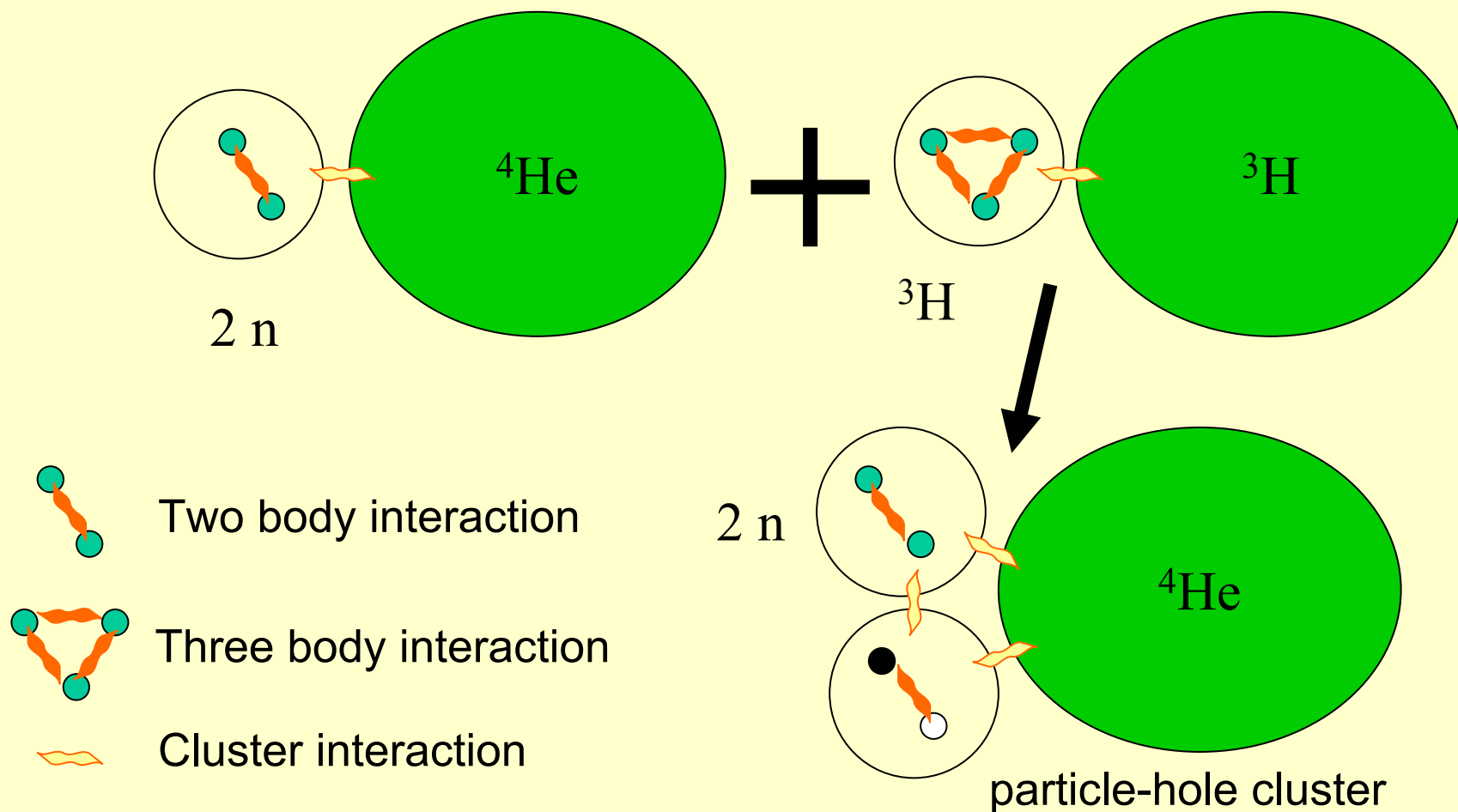
Shape coexistence in the vacuum of BDCM: Example ${}^6\text{He}$

In the ground state of ${}^6\text{He}$ two different clusters coexist:





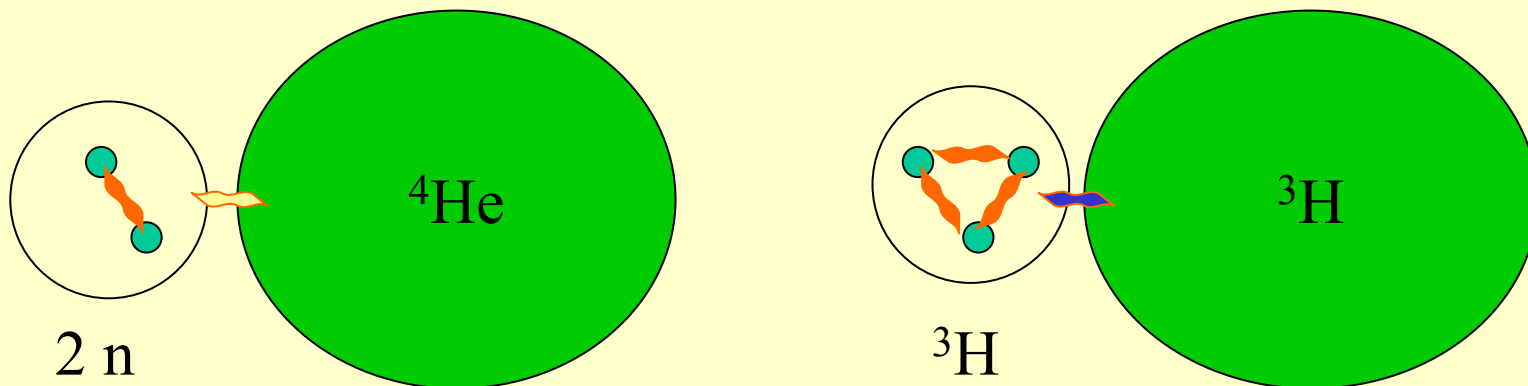
Interaction between clusters: deformation of the vacuum

In the ground state of ${}^6\text{He}$ two different clusters coexist:



From interacting clusters to DCM

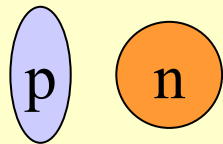
-  Macroscopic cluster interaction \longrightarrow ${}^2\text{H}$ cluster move
 in the optical potential formed by the alpha particle optical model.
-  Microscopic cluster interaction \longrightarrow N-particles which were
 partitioned in small and big clusters interact.



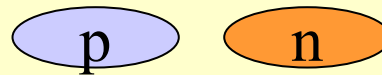
Can the dynamic correlation model reproduce the interacting cluster picture?

Prolate or oblate proton distributions?

- In AMD (Antisymetrized Molecular Dynamics) protons and neutrons are treated as separated clusters. By increasing the neutron number proton distribution can vary between two shapes.



${}^7\text{Li}$



${}^{11}\text{Li}$

Consequence: strong variation of proton charge radius as the neutron number increases.
Reason: the neutron average potential (mean field) is strongly deformed.

Introducing correlation in nuclei via the Unitarity Model Operator (UMO)

$$e^{iS} \prod_{i=1}^N a_i^+ |0\rangle = (1 + S_1 + S_2 + \dots) \prod_{i=1}^N a_i^+ |0\rangle$$

$$S_1 \propto a_1^+ a_2$$

$$S_2 \propto a_1^+ a_2 a_3^+ a_4$$

.....

Disadvantage:

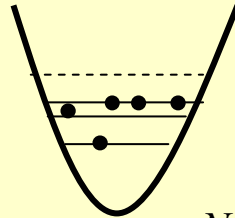
1. Perturbative
2. No Pauli Principle
3. Works with effective operator

$$H_{eff} = e^{-iS^+} H e^{iS}$$

$$O_{eff} = e^{-iS^+} O e^{iS}$$

Nuclear model based on Dynamic Correlation Model (DCM)

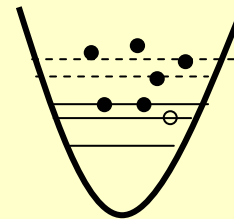
$$e^{iS} |\Psi_N\rangle \propto e^{iS} \prod_{i=1}^N a_i^+ |0\rangle$$



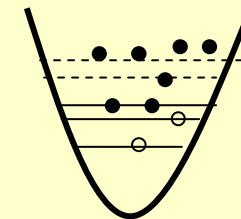
$$= \alpha |\Psi_N\rangle + \beta |\Psi_{N+1,1}\rangle + \gamma |\Psi_{N+2,2}\rangle + \dots = \alpha \prod_{i=1}^N a_i^+ |0\rangle + \beta \prod_{i=1}^{N+1} a_i^+ a_1 |0\rangle + \gamma \prod_{i=1}^{N+2} \prod_{i'=1}^2 a_i^+ a_{i'} |0\rangle \dots$$

$\alpha, \beta, \gamma, \dots$ are calculated from the following comutator chain:

$$[H, \prod_{i=1}^N a_i^+] |0\rangle \propto \prod_{i=1}^N a_i^+ |0\rangle + \prod_{i=1}^{N+1} a_i^+ a_1 |0\rangle$$



$$[H, \prod_{i=1}^{N+1} a_i^+ a_1] |0\rangle \propto \prod_{i=1}^{N+1} a_i^+ a_1 |0\rangle + \prod_{i=1}^{N+2} \prod_{i'=1}^2 a_i^+ a_{i'} |0\rangle$$



Boson Dynamic Correlation Model (BDCM)

$$[H, \prod_{i=1}^{N=2} a_i^+] |0\rangle \propto \prod_{i=1}^{N=2} a_i^+ |0\rangle + \prod_{i=1}^{N=3} a_i^+ a_1 |0\rangle$$

$$[H, \prod_{i=1}^{N=3} a_i^+ a_1] |0\rangle \propto \prod_{i=1}^{N=3} a_i^+ a_1 |0\rangle + \prod_{i=1}^{N=4} \prod_{i'=1}^{N=2} a_i^+ a_{i'} |0\rangle$$

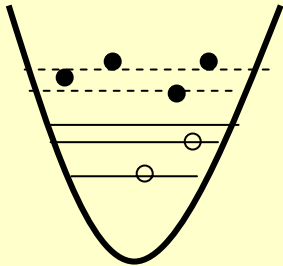
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Advantage:

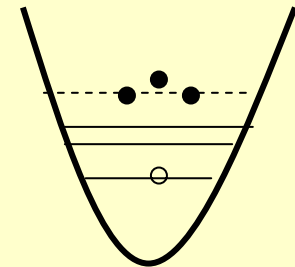
1. The commutator chain reduced to an eigenvalue problem by introducing dynamic linearisation approximation
2. Pauli Principle
3. Microscopic calculation without effective operators

Example for linearisation approximation

- In this work we are mainly concerned with calculating admixture coefficients for the ground state wavefunction.
 - That means that vacuum boiling configuration of higher complexity are poorly admixed.
- Under these considerations we introduce the following approximations:



$$\prod_{i=1}^{N=4} \prod_{i'=1}^{N'=2} a_i^+ a_{i'} |0\rangle \propto \sum_l \prod_{j=1}^{N=3} a_{lj}^+ a_l |0\rangle$$



With the linearisation the 4 particle – 2 holes configuration are approximated with effective 3 particle – 1 hole configurations

Effect of linearisation on commutator chain

- Use the linearisation approximation defined in the previous transparency
- Collect the resulting terms



$$\begin{vmatrix} E - E_{2p} + \langle 2p \| V \| 2p \rangle & \langle 2p \| V \| 3p1h \rangle \\ \langle 3p1h \| V \| 2p \rangle & E - E_{3p1h} + \langle 3p1h \| V \| 3p1h \rangle \end{vmatrix} \begin{vmatrix} \chi_{2p} \\ \chi_{3p1h} \end{vmatrix} = 0$$

Dynamic eigenvalue equations for mixed mode amplitudes
2 particles \Rightarrow 3 particles – 1 hole

Dynamics eigenvalue equation for one dressed boson which is solvable self-consistently

Perturbation theory

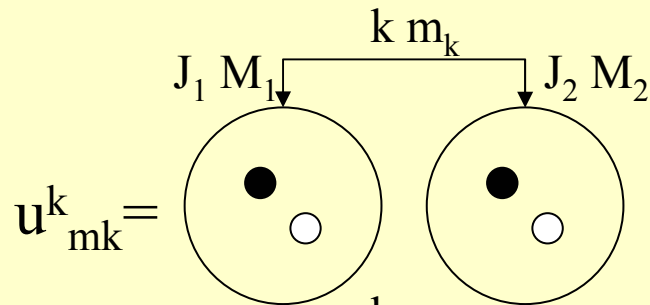
$$\begin{array}{c} \uparrow \uparrow \\ \bullet \end{array} = \begin{array}{c} \uparrow \uparrow \\ \text{wavy} \end{array} + \begin{array}{c} \uparrow \uparrow \\ \text{wavy} \\ \bullet \end{array} + \begin{array}{c} \uparrow \uparrow \\ \text{wavy} \\ \bullet \bullet \end{array} + \dots$$

Collective Hamiltonian and dressed bosons

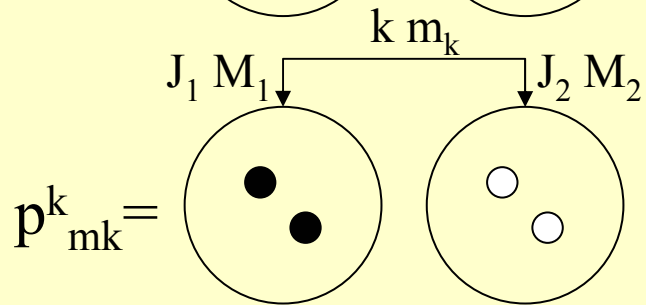
$$\left(\begin{array}{c} \text{H}_{\text{coll}} \end{array} \right) \left(\begin{array}{c} \uparrow \uparrow \\ \uparrow \uparrow \downarrow \\ \uparrow \uparrow \downarrow \downarrow \end{array} \right) \longrightarrow \left(\begin{array}{ccc} E_i & 0 & 0 \\ 0 & E_i & 0 \\ 0 & 0 & E_i \end{array} \right) \left(\begin{array}{c} \uparrow \uparrow \\ \uparrow \uparrow \text{wavy} \\ \uparrow \uparrow \text{wavy} \bullet \end{array} \right)$$

Symmetry properties of the configuration mixing wavefunction (CMWF)

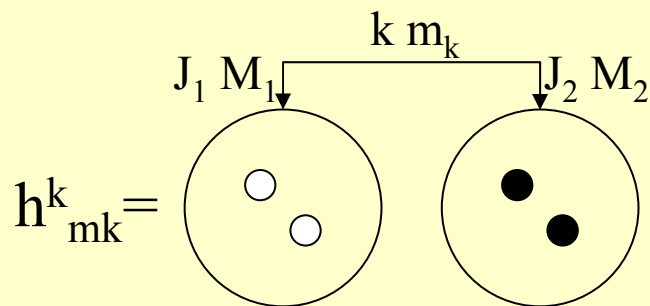
Define operators:



Destroy a hole-particle pair
and create a particle-hole pair



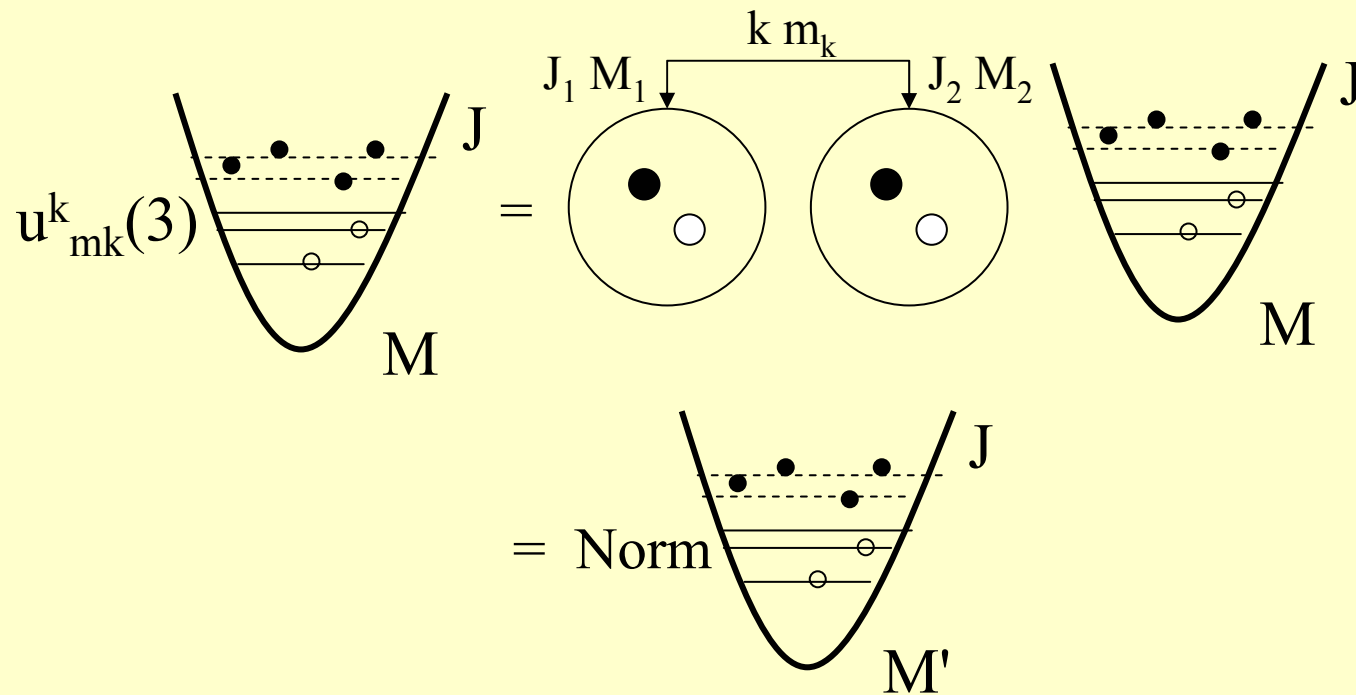
Destroy a hole-hole pair
and create a particle-particle pair



Destroy a particle-particle pair
and create a hole-hole pair

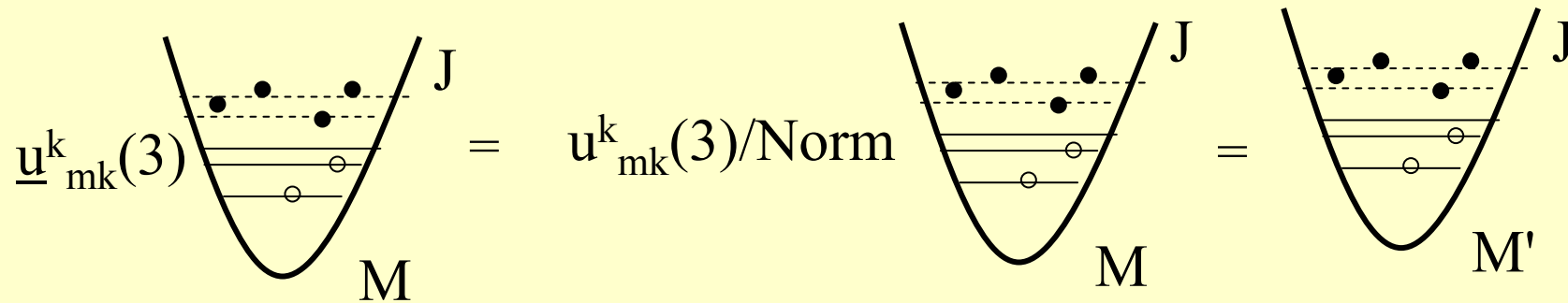
Group property of the defined operators

Action of the operator on totally antisymmetric wavefunction



The same action is valid for p^k_{mk} and h^k_{mk}

Unitary operator

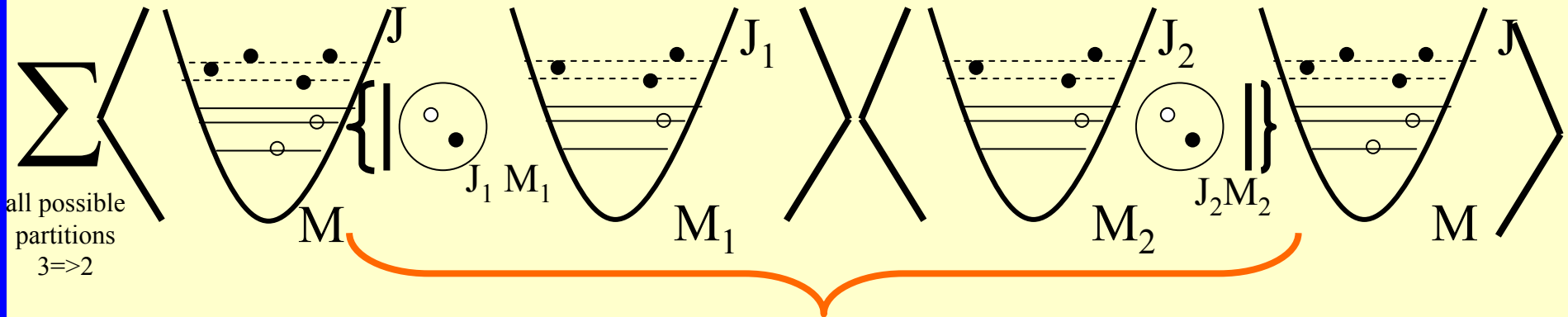
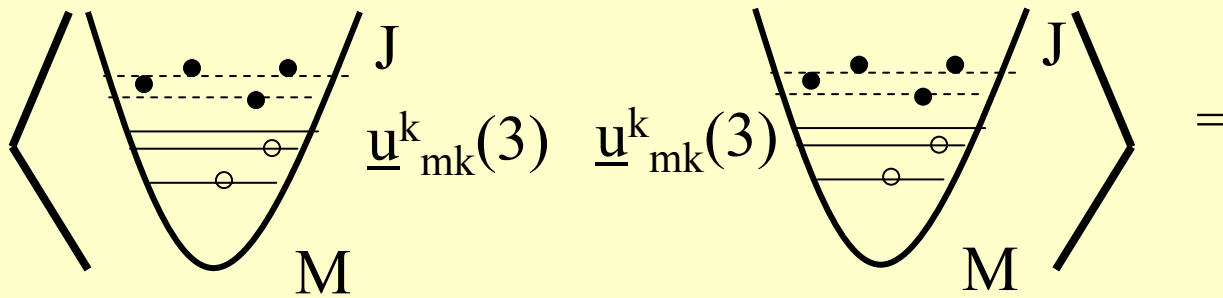


Commutator relation :

$$\underline{u}^{k1}_{mk1}(3) \underline{u}^{k2}_{mk2}(3) = \text{geometric factor} * \underline{u}^{k3}_{mk3}(3)$$

Therefore the unitary operators are generators of the $SU_{2J+1}(3)$

Cluster transformation coefficients



Cluster Transformation Coefficients

Factorisation of the model CMWFs (for electrons and nucleons) in terms of cluster coefficients

configuration	associated matrix elements / potentials	configuration	associated matrix elements / potentials	configuration	associated matrix elements / potentials
2p	<p>realistic G matrix</p>	4p	<p>realistic G matrix</p>	4p-2h	<p>realistic G matrix and phenomenological particle-hole potential</p>
1p-1h	<p>phenomenological particle-hole matrix element</p>	$= \sum_{ijkl} B_{ijkl}^1$	$= \sum_{ijklmn} \left[D_{ijklmn}^1 \right]$		
3p	<p>realistic G matrix</p> $= \sum_{ijk} A_{ijk}^1$	3p-1h	<p>realistic G matrix and phenomenological particle-hole potential</p> $= \sum_{ijkl} \left[B_{ijkl}^2 + B_{ijkl}^3 \right]$	$+ D_{ijklmn}^2$	
2p-1h	<p>realistic G matrix and phenomenological particle-hole potential</p> $= \sum_{ijk} \left[A_{ijk}^2 + A_{ijk}^3 \right]$	4p-1h	<p>realistic G matrix and phenomenological particle-hole potential</p> $= \sum_{ijklm} \left[C_{ijklm}^1 + C_{ijklm}^2 \right]$	$+ D_{ijklmn}^3$	

Distribution in BDCM

$$\begin{aligned}
 \rho_{CM}(r) &= \frac{1}{A} \langle \psi_{dressed}^{(2n-or-pn)} | \sum_{i=1,6} \delta(\vec{r} - \vec{r}_i) | \psi_{dressed}^{(2n-or-pn)} \rangle \\
 &= \frac{1}{A} \langle \psi(\vec{r}_1, \vec{r}_2) | \sum_{i=1,6} \delta(\vec{r} - \vec{r}_i) | \psi(\vec{r}_3, \vec{r}_4) \rangle \\
 &+ C_{cluster}^1 \langle \psi^a(\vec{r}_i, \vec{r}_j) \psi^p(\vec{r}_k, \vec{r}_l) | \sum_{i=1,6} \delta(\vec{r} - \vec{r}_i) | \psi(\vec{r}_3, \vec{r}_4) \rangle \quad (1) \\
 &+ C_{cluster}^2 \langle \psi(\vec{r}_1, \vec{r}_2) | \sum_{i=1,6} \delta(\vec{r} - \vec{r}_i) | \psi^a(\vec{r}_k, \vec{r}_l) \psi^p(\vec{r}_k, \vec{r}_l) \rangle \\
 &+ C_{cluster}^1 * C_{cluster}^2 \langle \psi^a(\vec{r}_i, \vec{r}_j) \psi^p(\vec{r}_k, \vec{r}_l) | \sum_{i=1,6} \delta(\vec{r} - \vec{r}_i) | \psi^a(\vec{r}_k, \vec{r}_l) \psi^p(\vec{r}_k, \vec{r}_l) \rangle
 \end{aligned}$$

a=active, p=passive

for the pairs we use the Moshinski transformation:

$$\psi_{n_1 l_1}(\vec{r}_1), \psi_{n_2 l_2}(\vec{r}_2) = \sum_{nlNL\lambda} \langle n_1 l_1 n_2 l_2 \lambda | \} nlNL\lambda \rangle \phi_{nl}(\vec{r}) \phi_{NL}(\vec{R}) \quad (2)$$

Distribution in BDCM

Before performing the Moshinski transformation we go from jj to ls coupling scheme.

JJ→LS transformation:

$$|\Phi(n_1 l_1 j_1) \Phi(n_2 l_2 j_2); JT\rangle =$$

$$\hat{j}_1 \hat{j}_2 \sum_{\lambda S J_i n l N L} \langle n_1 l_1 n_2 l_2 \lambda | \{ n l N L \lambda \rangle \quad (1)$$

$$9j(j_1 l_1 \frac{1}{2}; j_2 l_2 \frac{1}{2}; J \lambda S) 6j(L l \lambda; S J J_i) |\psi_{nl}(\vec{r}), \psi_{NL}(\vec{R})\rangle |\chi_1^{\frac{1}{2}} \chi_2^{\frac{1}{2}} S\rangle |\tau_1^{\frac{1}{2}} \tau_2^{\frac{1}{2}} T\rangle$$

This should be used also for the $\langle bra|$.

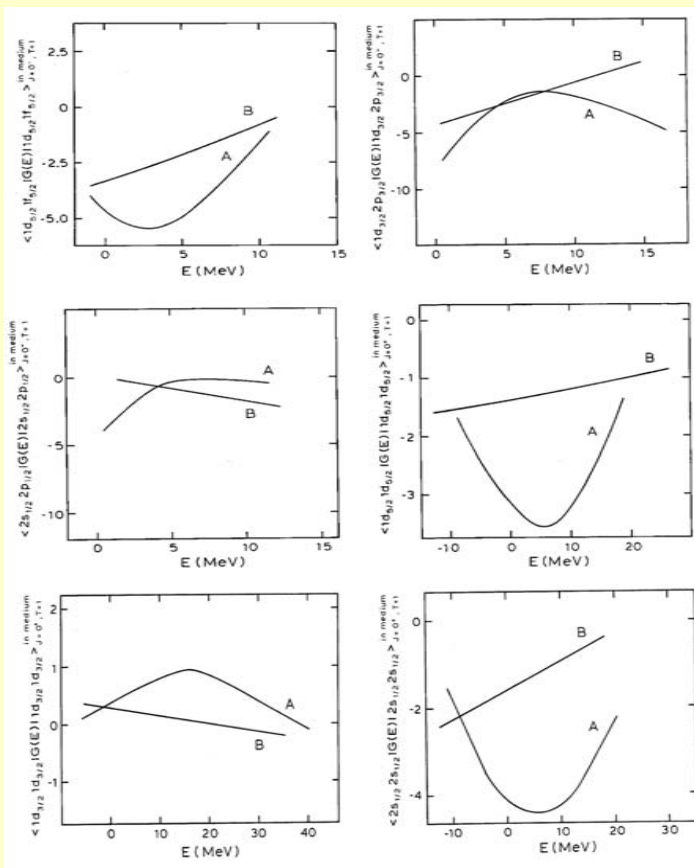
Distribution in BDCM

For one pair of particle the distribution is then:

$$\begin{aligned}
 & \langle \Phi(n'_1 l'_1 j'_1) \Phi(n'_2 l'_2 j'_2); JT | \delta(\vec{r} - \vec{r}_i) | \Phi(n_1 l_1 j_1) \Phi(n_2 l_2 j_2); JT \rangle = \\
 & \sum_{i \lambda S J_i n l N L \lambda' n' l' N' L' \lambda'}^{\hat{j}_1 \hat{j}_2 \hat{j}'_1 \hat{j}'_2} \langle n_1 l_1 n_2 l_2 \lambda | \{ n l N L \lambda \} \langle n'_1 l'_1 n'_2 l'_2 \lambda' | \{ n' l' N' L' \lambda' \} \rangle \\
 & 9j(j_1 l_1 \frac{1}{2}; j_2 l_2 \frac{1}{2}; J \lambda S) 6j(L l \lambda; S J J_i) 9j(j'_1 l'_1 \frac{1}{2}; j'_2 l'_2 \frac{1}{2}; J \lambda' S) 6j(L' l' \lambda'; S J J_i) \\
 & \langle \psi_{n'l'}(\vec{r}), \psi_{N'L'}(\vec{R}_i) | \langle \chi_1^{\frac{1}{2}} \chi_2^{\frac{1}{2}} S | \langle \tau_1^{\frac{1}{2}} \tau_2^{\frac{1}{2}} T \rangle | \delta(\vec{R} - \vec{R}_i) | \psi_{nl}(\vec{r}), \psi_{NL}(\vec{R}_i) \rangle | \chi_1^{\frac{1}{2}} \chi_2^{\frac{1}{2}} S \rangle | \tau_1^{\frac{1}{2}} \tau_2^{\frac{1}{2}} T \rangle
 \end{aligned} \tag{1}$$

where; $\hat{j}_1 = \sqrt{2j_1 + 1}$, $9j$ are the $9j$ symbols and $6j$ the $6j$ symbols.

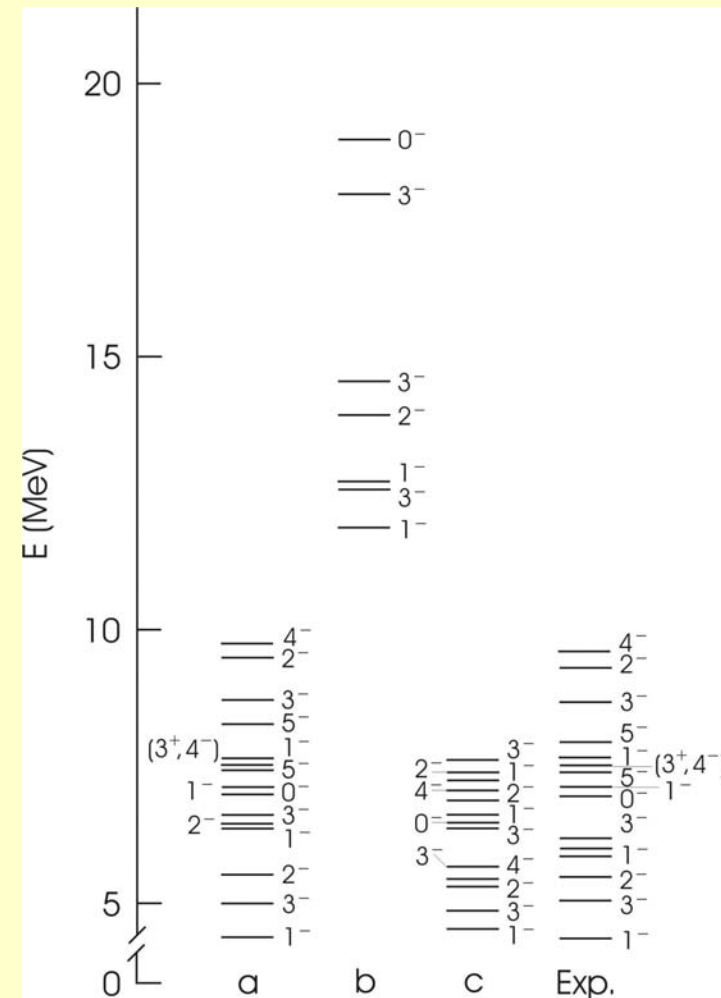
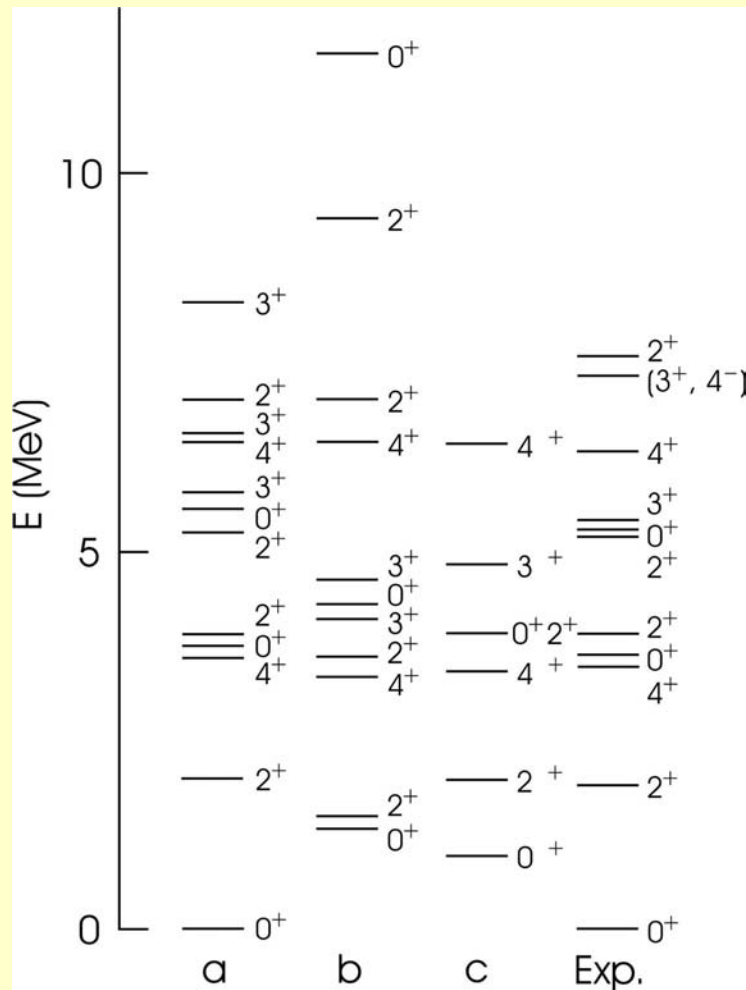
In medium G-matrix elements and corresponding two body potentials for ^{18}O



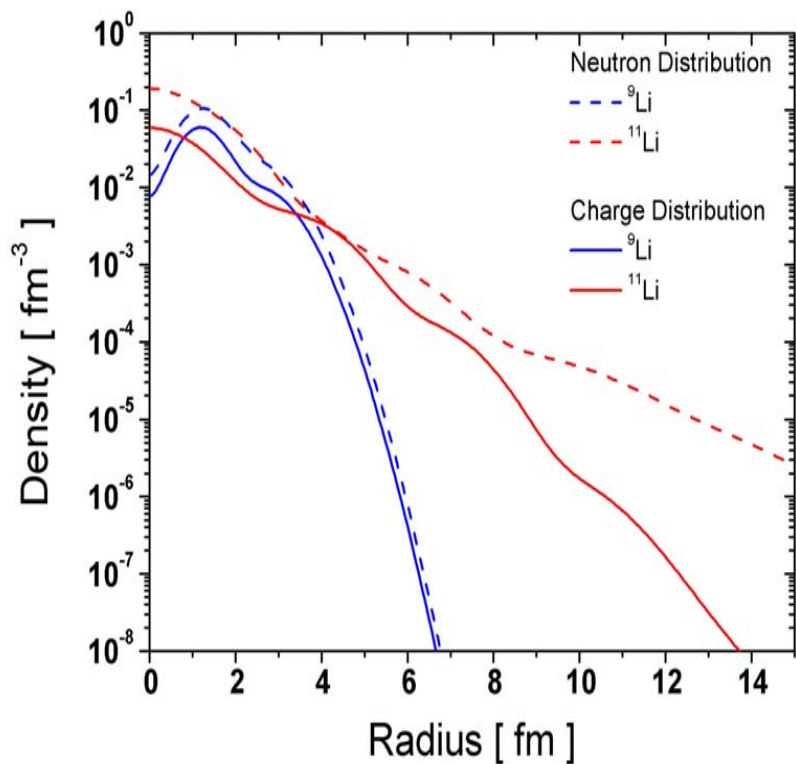
Matrix elements	$\langle 44 V 44 \rangle$	$\langle 55 V 55 \rangle$	$\langle 66 V 66 \rangle$
Bare G-matrix	-0.68	-4.49	2.59
G in medium (E=0. MeV)	-2.01	-1.47	-1.30
G in medium (E=3.65 MeV)	-2.63	-1.59	-0.91
G in medium (E=5.33 MeV)	-2.91	-1.92	-0.89
Folded G-matrix and Paris potential	-2.22	-1.61	-0.95
Folded G-matrix and Bonn A potential	-2.77	-2.05	-1.28
Empirical	-2.82	-2.12	-2.18

Matrix elements for the first three positive 0^+ levels.

Spectrum of the positive and negative parity state of ^{18}O

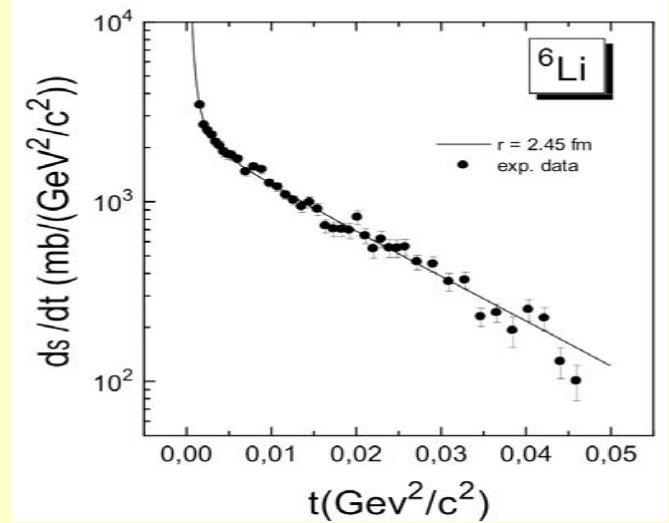
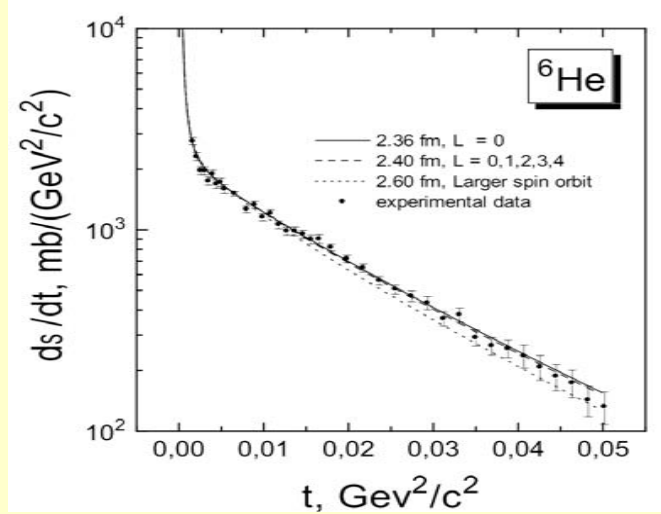
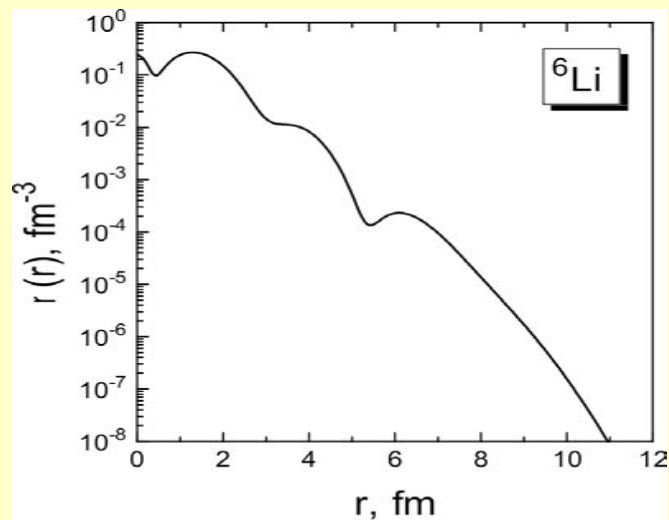
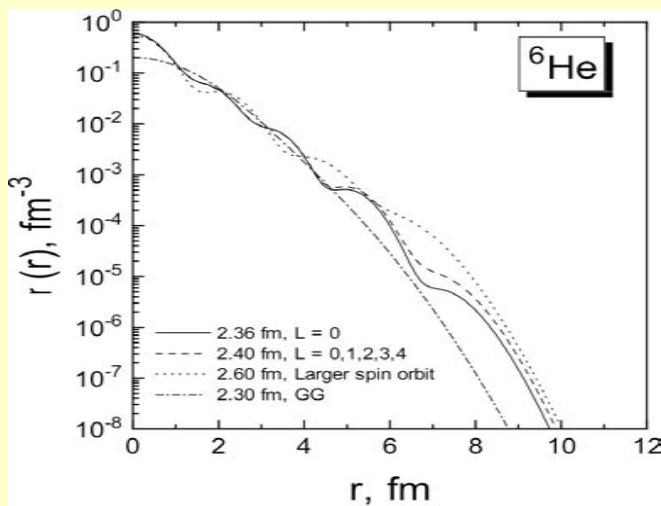


Nuclear results for Li isotopes



	$R_{\text{matter}}^{\text{calc}}$	$R_{\text{matter}}^{\text{exp}}$	$R_{\text{charge}}^{\text{calc}}$	$R_{\text{charge}}^{\text{exp}}$	μ^{calc}	μ^{exp}
${}^6\text{Li}$	2.41	2.45(7)	2.55	2.55(4)	0.82	0.82204(6)
${}^7\text{Li}$	2.36	2.35(3)	2.41	2.39(20)	3.25	3.2564
${}^9\text{Li}$	2.42	2.43(7)	2.42	?	3.44	3.43(6)
		2.32(2)				
${}^{11}\text{Li}$	2.30(2)	3.62(19)	2.67	?		
	3.64	3.12(16)				
	3.53(10)					

Distribution and proton scattering cross-section



Spectroscopic factor and groundstate wavefunction of ${}^6\text{Li}$

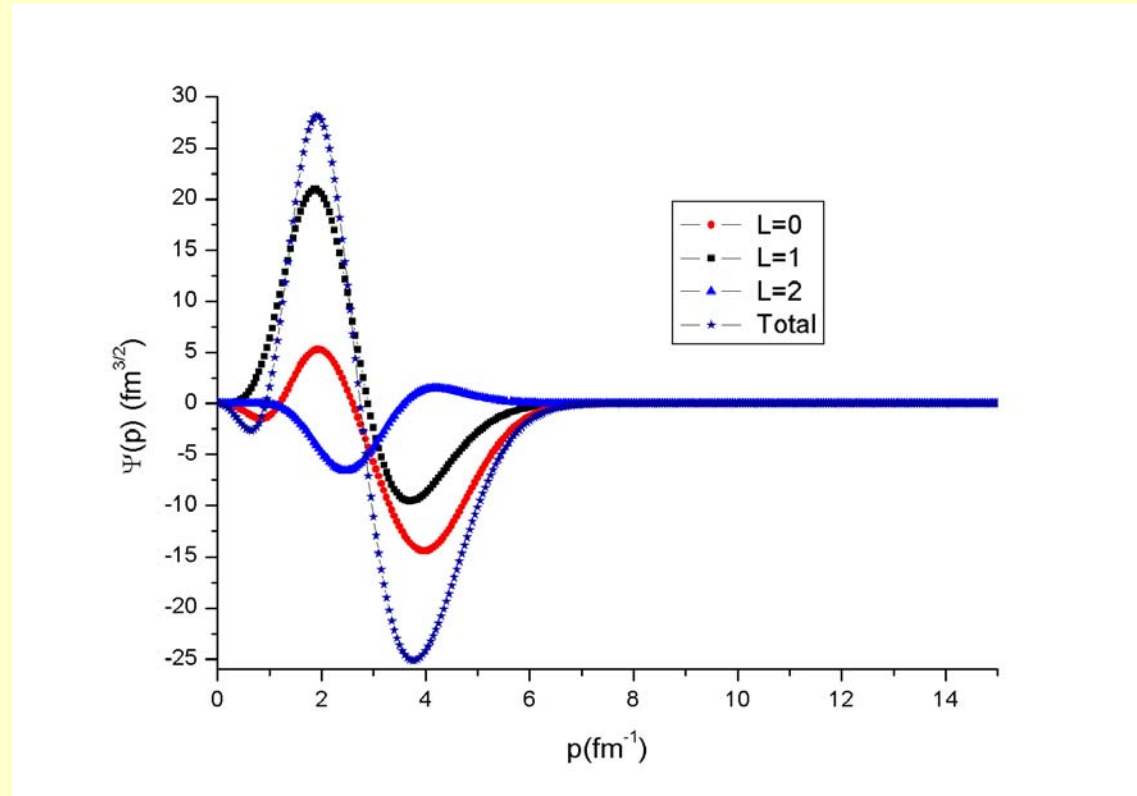
Spectroscopic factor for ${}^6\text{Li}$

pn in $p_{\frac{3}{2}}$ shell	pn in (psd) shell	pn in (psd) shell and 3p1h
1.0	0.6657	0.6564

Two particle components of the ground state wave function of ${}^6\text{Li}$

	$1p_{\frac{3}{2}}1p_{\frac{3}{2}}$	$1p_{\frac{3}{2}}1p_{\frac{1}{2}}$	$1p_{\frac{3}{2}}2p_{\frac{3}{2}}$	$1p_{\frac{3}{2}}1f_{\frac{5}{2}}$	$1p_{\frac{3}{2}}2p_{\frac{1}{2}}$	$1p_{\frac{1}{2}}1p_{\frac{1}{2}}$	$1p_{\frac{1}{2}}2p_{\frac{3}{2}}$	$1p_{\frac{1}{2}}2p_{\frac{1}{2}}$
pn	.8159	.5576	.0133	.0091	.0154	-.1133	-.0179	-.0106
pn+3p1h	.8102	.5598	.0142	.0091	.0158	-.1165	-.0175	-.0110
	$1d_{\frac{5}{2}}1d_{\frac{5}{2}}$	$1d_{\frac{5}{2}}1d_{\frac{3}{2}}$	$2s_{\frac{1}{2}}2s_{\frac{1}{2}}$	$2s_{\frac{1}{2}}1d_{\frac{3}{2}}$	$1d_{\frac{3}{2}}1d_{\frac{3}{2}}$	$1f_{\frac{7}{2}}1f_{\frac{7}{2}}$	$1g_{\frac{9}{2}}1g_{\frac{9}{2}}$	-

Momentum components of the in medium proton neutron wavefunction in ${}^6\text{Li}$



Summary of Charge Radii

Ref.	this	[8]	[1]	[1]	[2]	[3]	[4]	[5]	[6]	[7]
	Exp.	Exp.	Exp.+Th.	Exp.+Th.	Theory	Theory	Theory	Theory	Theory	Exp.+Th.
Li -	rms R_c	rms R_c	rms R_p	rms R_c	rms R_p	rms R_p		rms R_p	rms R_c	rms R_c
6	2.55	2.55 (4)	2.32 (3)	2.46 (2)	2.045	2.39			2.55	
7	2.46	2.37 (3)	2.27 (2)	2.40 (2)	1.941	2.25		2.27	2.41	
8	2.37		2.26 (2)		1.946	2.09		2.18		
9	2.30		2.18 (2)		1.986		2.04	2.10	2.42	
11			2.88 (2)						2.67	2.235

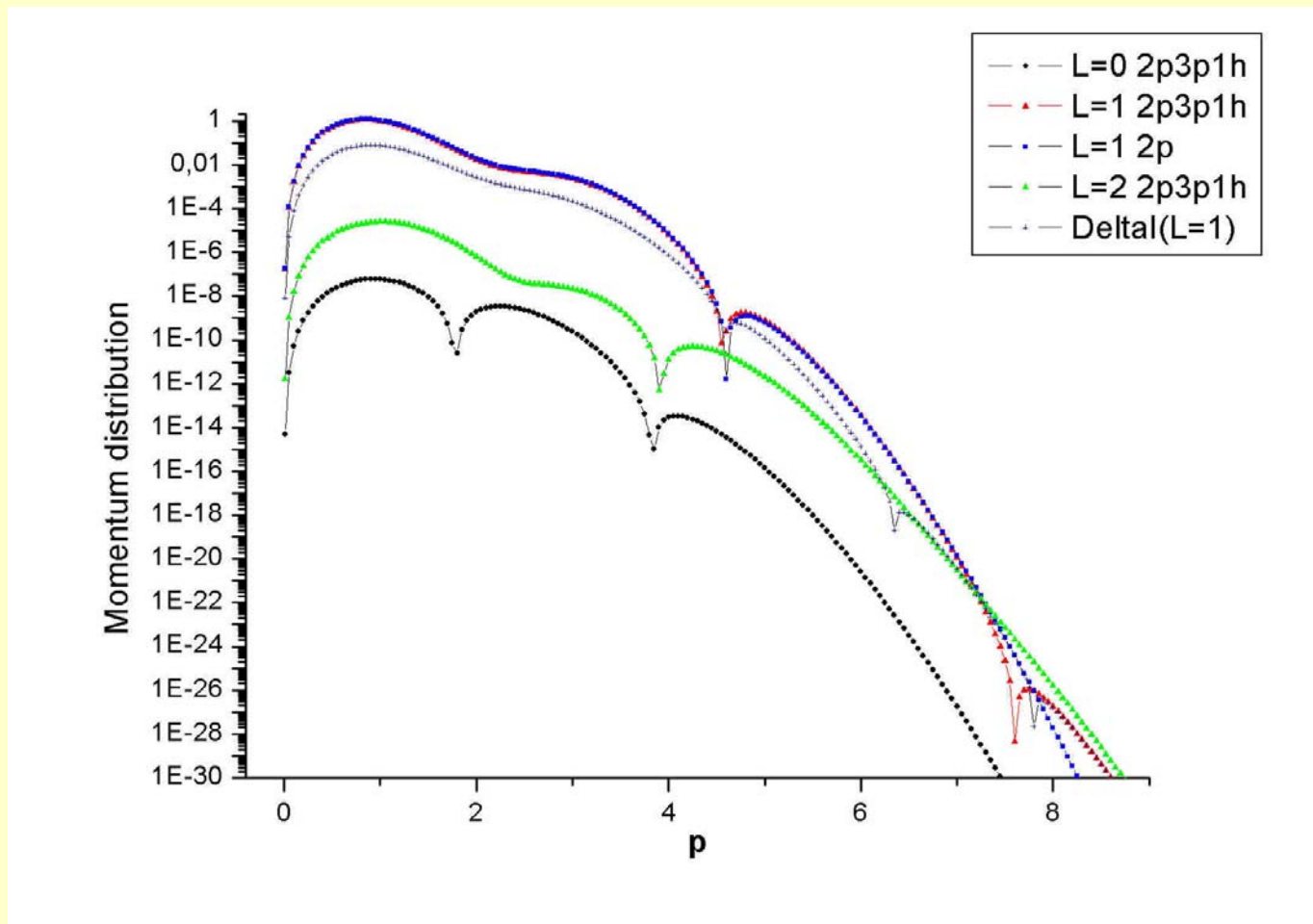
References:

- [1] I. Tanihata, Phys. Lett B 206,592 (1988)
- [2] P. Navratil, PRC 57,3119 (1998)
- [3] S. Pieper, Annu.Rev.Nucl.Part.Sci. 51, 53 (2001)
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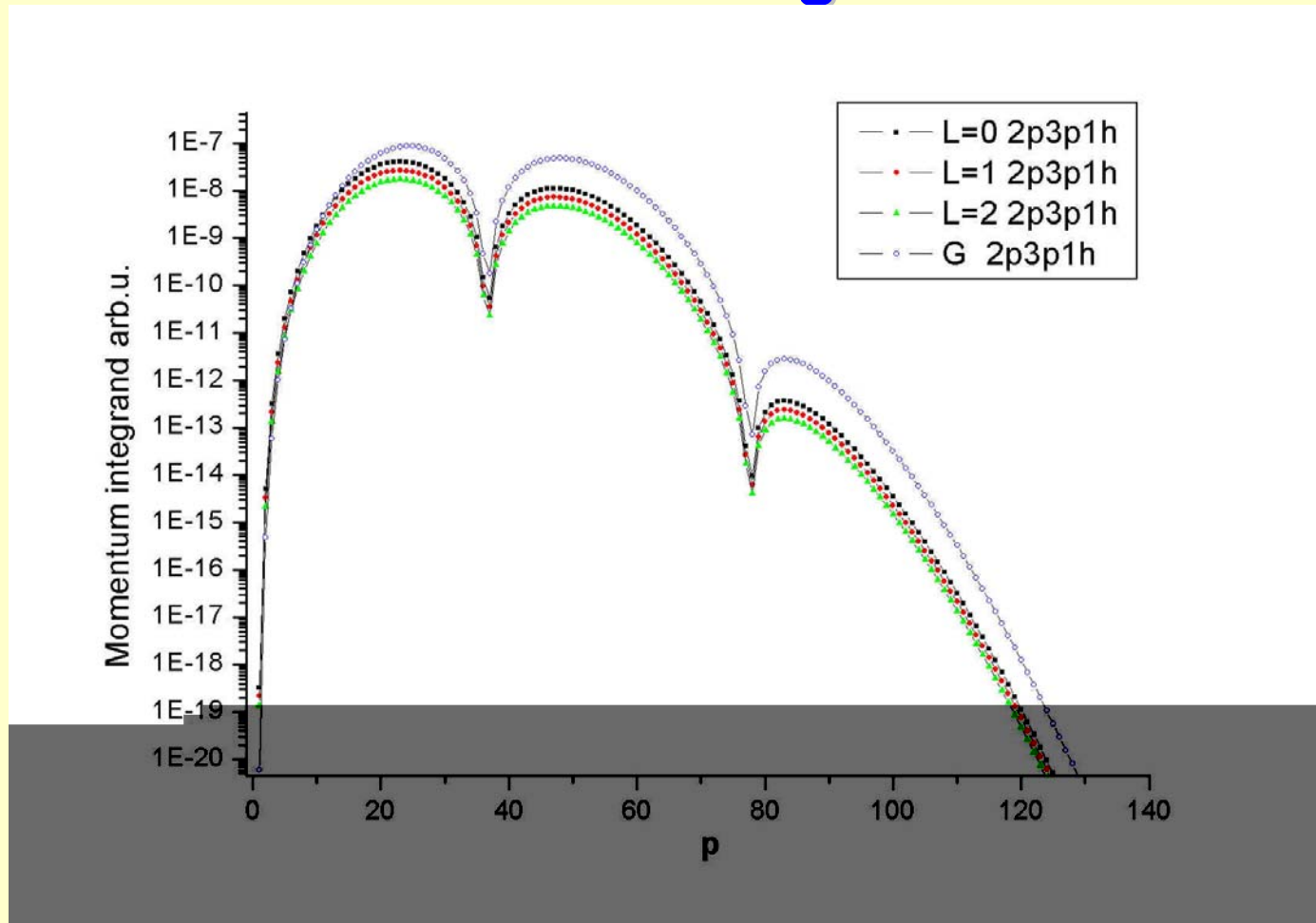
Method:

- Interaction Cross Sections with Glauber model, HO distributions
- Large-basis shell-model calculations
- Greens Function Monte Carlo AV18/IL2
- Greens Function Monte Carlo AV18/IL2
- Stochastic Variational Multicluster Method on a correlated gaussian basis
- Dynamic Correlation model
- coupled channel calculations, double-folding optical potential, M3Y effective interaction
- Electron Scattering

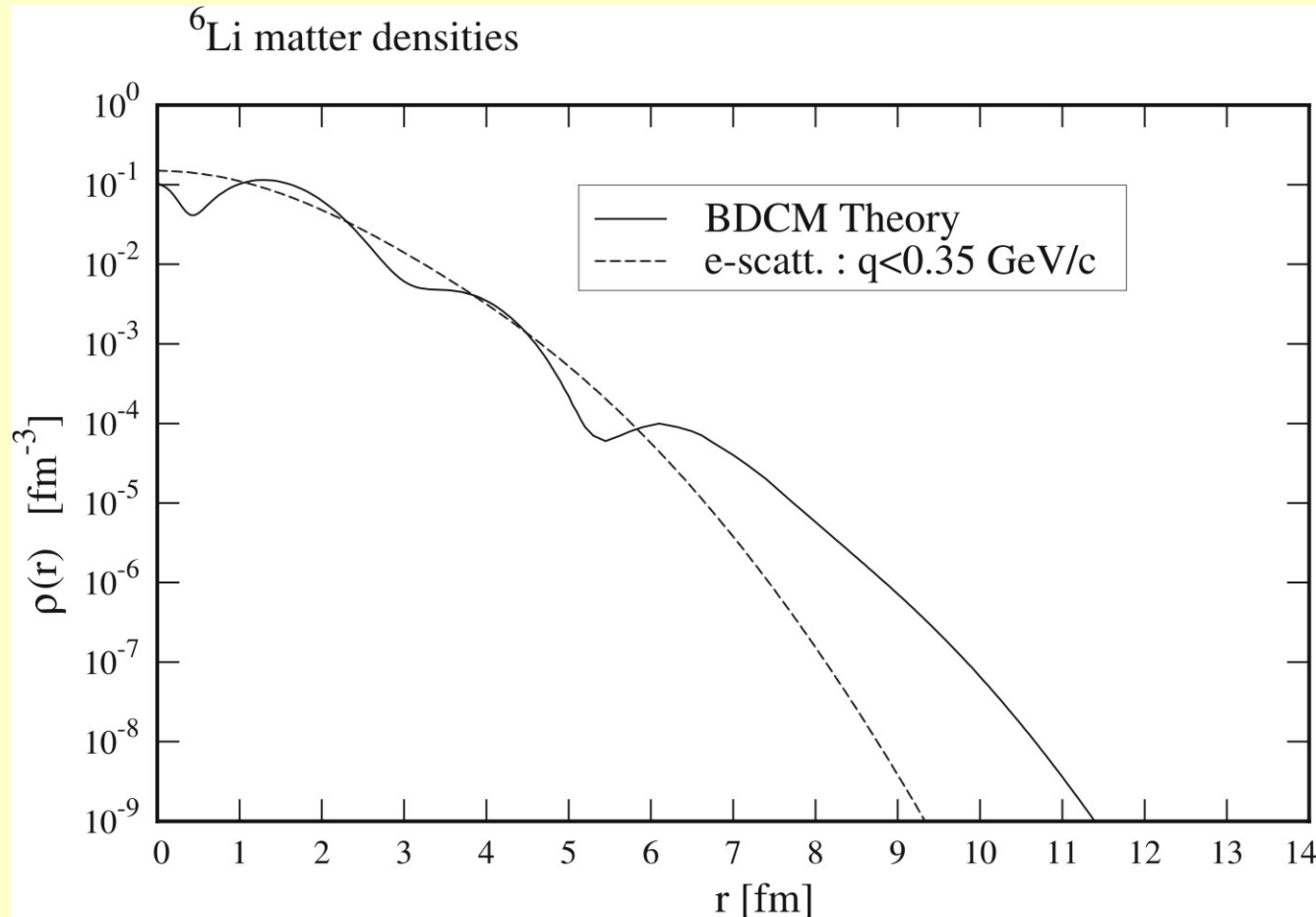
Nuclear results for Li isotopes: momentum distribution



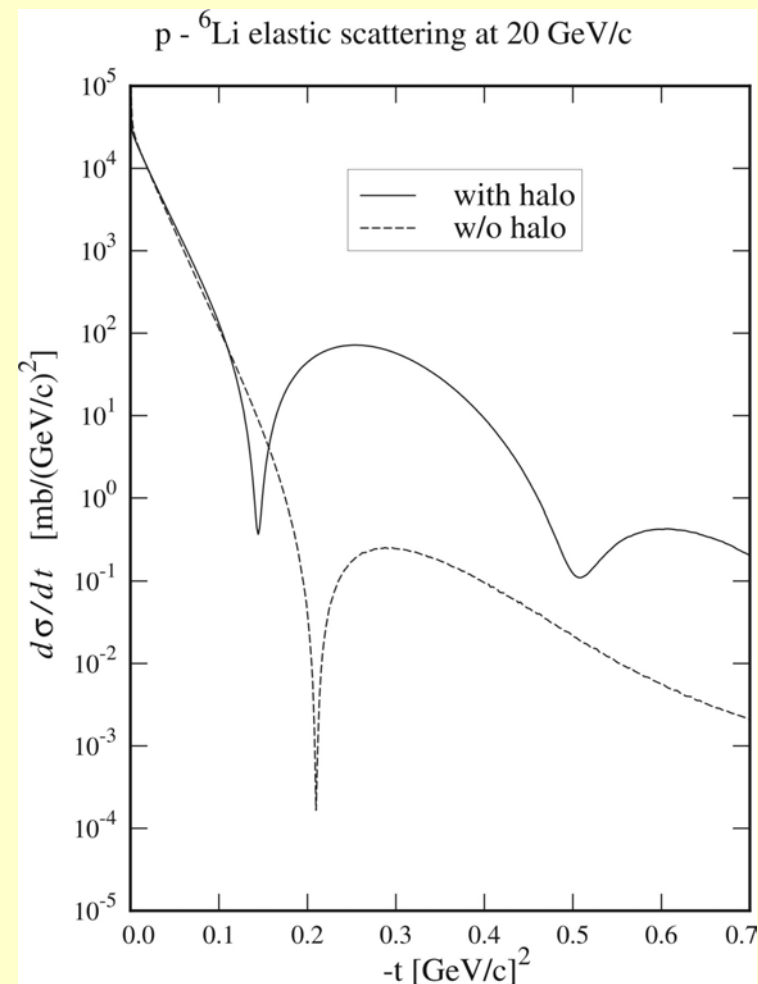
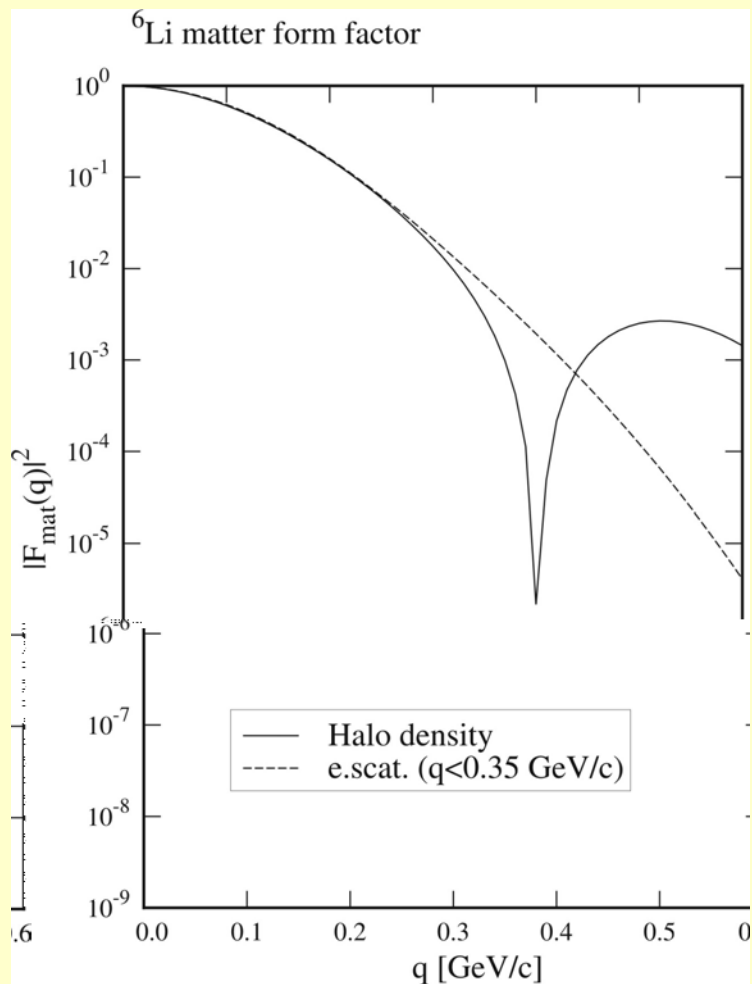
Nuclear results for Li isotopes: Momentum integrands



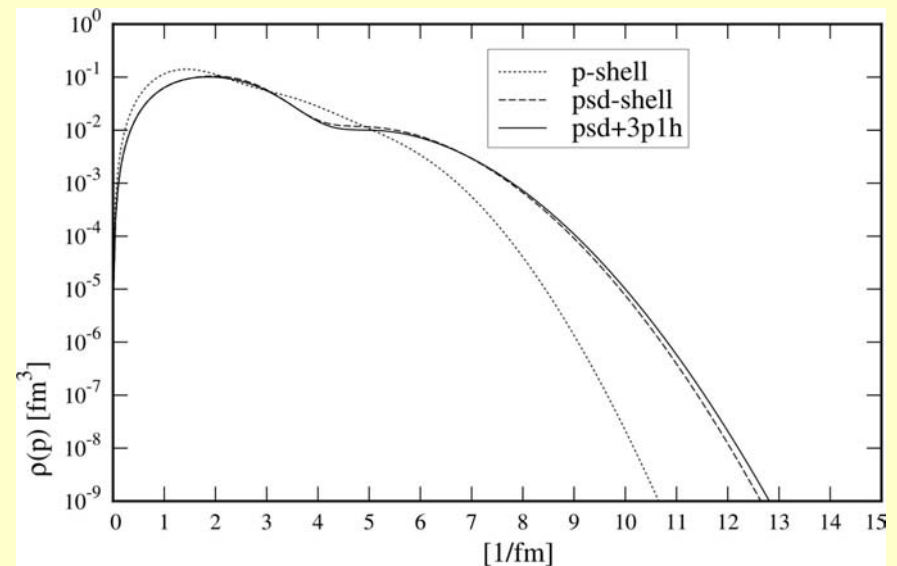
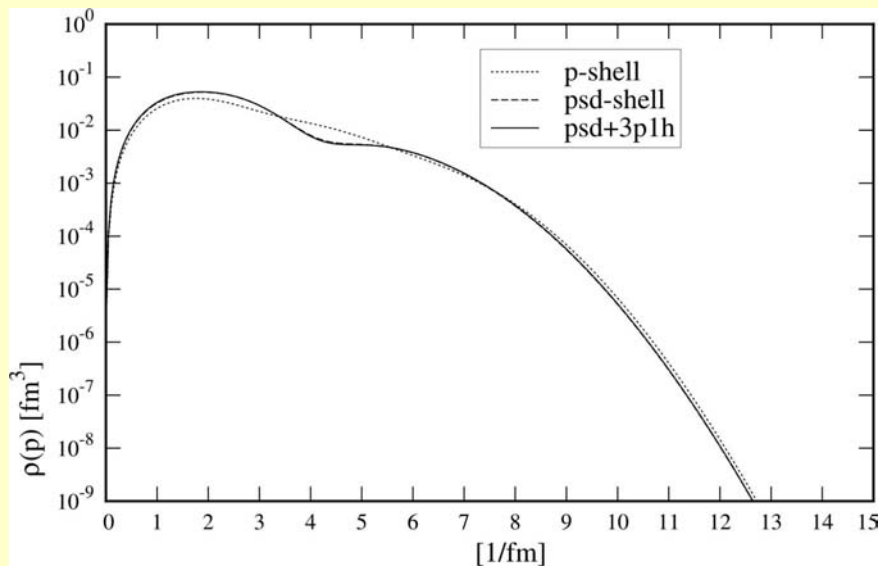
Nuclear results for Li isotopes



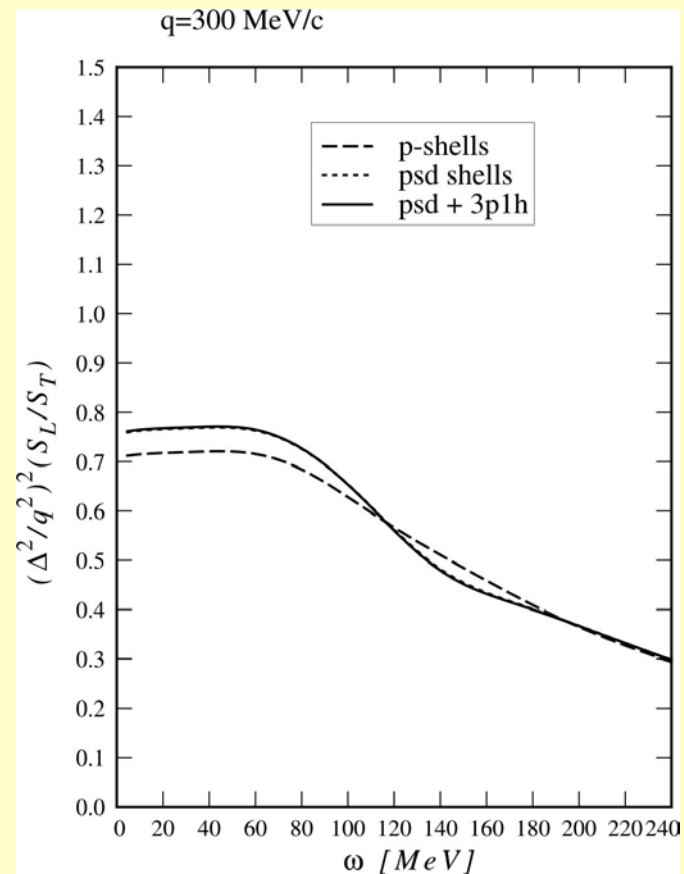
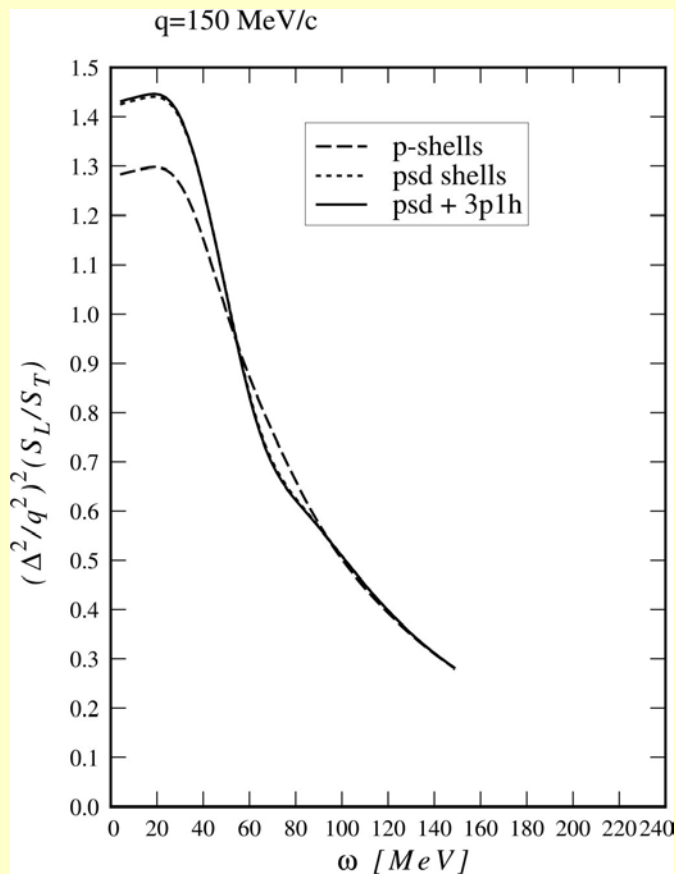
Nuclear results for Li isotopes



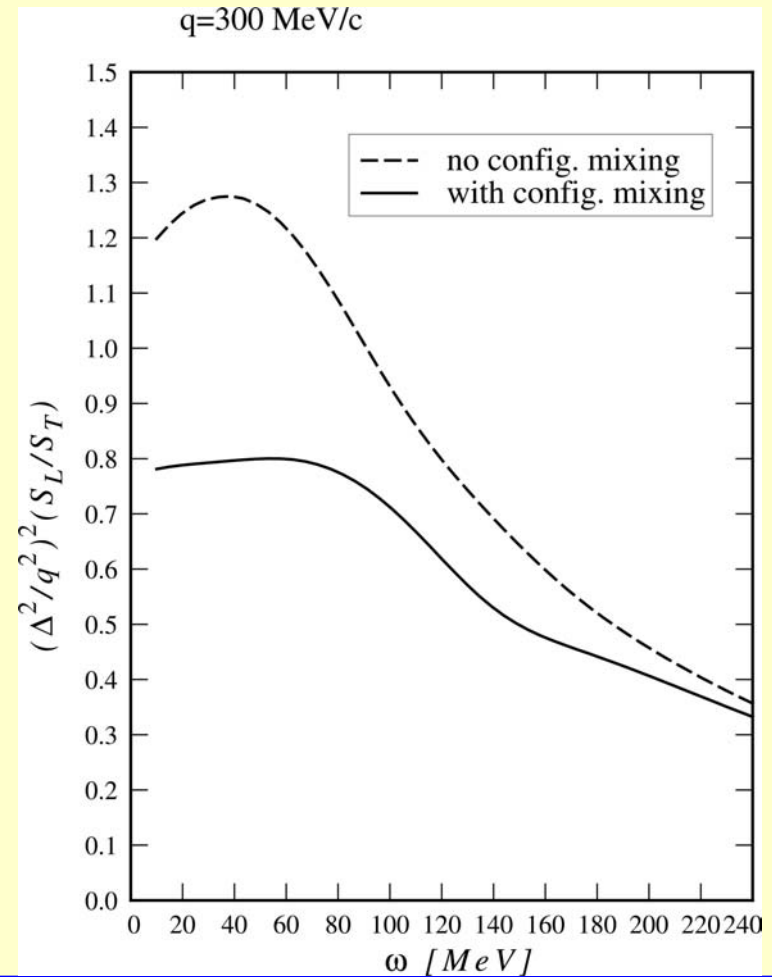
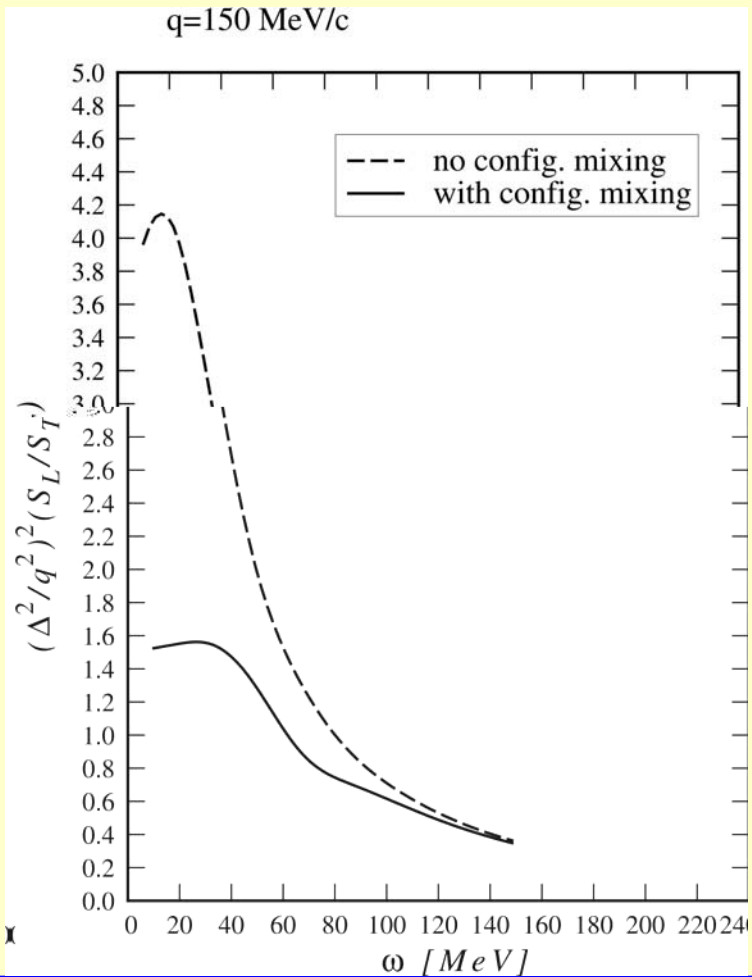
Momentum distribution for p and d in ${}^6\text{Li}$



Calculated singularity-free longitudinal to transverse response ratio R for ee'p



Calculated singularity-free longitudinal to transverse response ratio R for ee'd



Conclusions

- $(ee'p)$ and $(ee'd)$ give a good test for the halo structure of exotic nuclei
- Microscopic structure of the nucleus play an important role
- Result can be used as prediction for new experiments at the future GSI
- Presented method should be systematically applied to nuclei with pronounced halo like ^{11}Li
- Effective method to predict also the proton halo