

# Core excitation in breakup reactions: semi-microscopic folding models for the structure of halo nuclei

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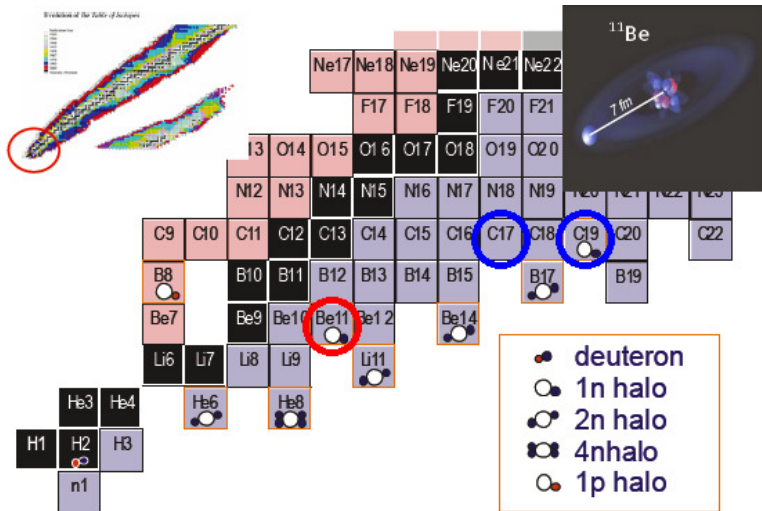
Trento, 27<sup>th</sup> May 2014



UNIVERSITÀ  
DEGLI STUDI  
DI PADOVA



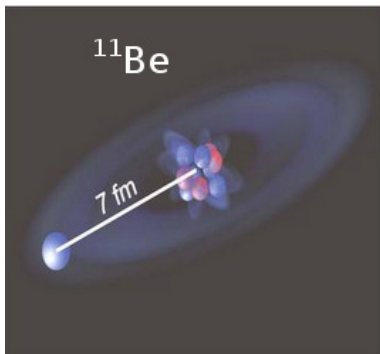
- 1 Motivation
  - Discretizing the Continuum
  - Core excitations
- 2 Structure
  - Halo nuclei with excitations of the core
  - PS discretization method
  - Semi-microscopic model
- 3 Results
  - Structure of  $^{11}\text{Be}$ ,  $^{19}\text{C}$ , and  $^{21}\text{C}$
  - More reactions with  $^{19}\text{C}$



# Simplest model for $^{11}\text{Be}$ and other halo nuclei

## Single Particle Model

- 1 Two-body model for  $^{11}\text{Be}$  ( $n+^{10}\text{Be}$ )
- 2 Ignore possible core-excited components of  $^{10}\text{Be}$ .
- 3 Reaction treated within CDCC formalism



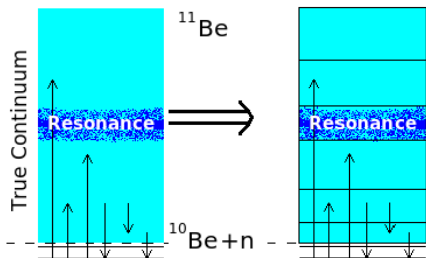
!! contribution of  $^{10}\text{Be}$  in its 1<sup>st</sup> excited state  $2^+$

# The Continuum problem in CDCC

Discretization methods:

- Binning:

$$u_i^j(k) = \sqrt{\frac{2}{\pi N_i}} \int_{k_{i-1}}^{k_i} f_i(k) \phi_I(k, r) dk.$$



# The Continuum problem in CDCC

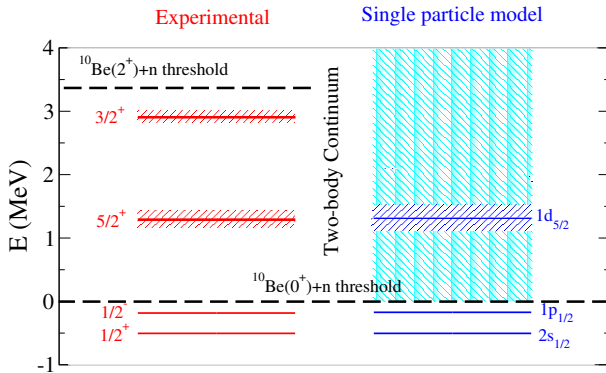
Discretization methods:

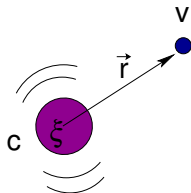
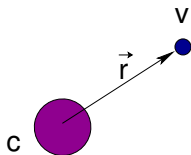
- Binning:

$$u_i^j(k) = \sqrt{\frac{2}{\pi N_i}} \int_{k_{i-1}}^{k_i} f_i(k) \phi_I(k, r) dk.$$

- Pseudo-states.

- ⇒ The Continuum is described through a basis (HO, Sturmian) of square-integrable states.
- ⇒ It improves the convergence rate in the case of narrow resonances.

$^{11}\text{Be}$  in a single particle model



✗ No core excitations

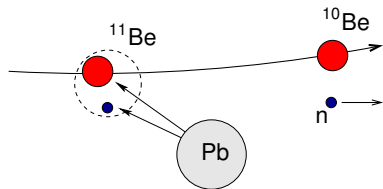
$$|^{11}\text{Be}(1/2^+)\rangle = |^{10}\text{Be}(0^+ \text{ g.s.}) \otimes \nu s_{1/2}\rangle$$

✓ Core excitations

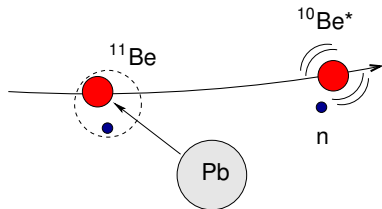
$$|^{11}\text{Be}(1/2^+)\rangle = \alpha |^{10}\text{Be}(0^+ \text{ g.s.}) \otimes \nu s_{1/2}\rangle + \beta |^{10}\text{Be}(2^+) \otimes \nu d_{5/2}\rangle + \dots$$

$|\alpha|^2, |\beta|^2 = \text{spectroscopic factors}$

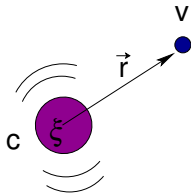




✗ Pure valence excitation



✓ Core-excitation mechanism



## Hamiltonian with core excitation

$$\mathcal{H}_p = T(\vec{r}) + h_{core}(\xi) + V_{NC}(\vec{r}, \vec{\xi})$$

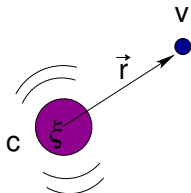
### Model for the core $h_{core}(\xi)$

- Selecting the model space  $\Rightarrow$  which states are included
- The model for core excitations will determine  $V_{NC}(\vec{r}, \vec{\xi})$

### Same formalism for different interaction models:

- Particle-Rotor model (deformed core)
- Particle-Vibration
- From microscopic transition densities
- ...

## Weak coupling limit



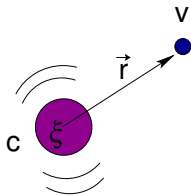
Hamiltonian with core excitation

$$\mathcal{H}_p = T(\vec{r}) + h_{core}(\xi) + V_{NC}(\vec{r}, \vec{\xi})$$

We look for a basis including core degrees of freedom

Coupling core  $\varphi_I(\vec{\xi})$  and single particle  $\mathcal{Y}_{lsj}(\hat{r})$  to the total  $J_p$ 
 $\Rightarrow n_\alpha$  different possible combinations or channels  $\alpha = \{l, s, j, I\}$

## Generalization of Pseudo-states (PS) discretization method



Hamiltonian with core excitation

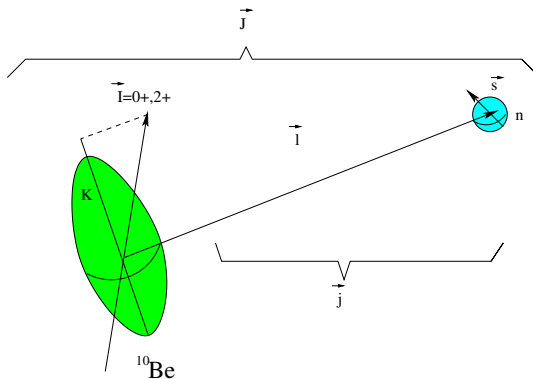
$$\mathcal{H}_p = T(\vec{r}) + h_{core}(\xi) + V_{NC}(\vec{r}, \vec{\xi})$$

Set of  $\mathcal{L}^2$  functions in this scheme:

$$|\phi_{i,J_p}(\vec{r}, \vec{\xi})\rangle = \sum_{\alpha} R_{i,\alpha}^{THO}(r) \left[ \mathcal{Y}_{l_s j}(\hat{r}) \otimes \varphi_I(\vec{\xi}) \right]_{J_p} \quad i = 1, \dots, N$$

⇒ Total number of functions: N times the number of channels

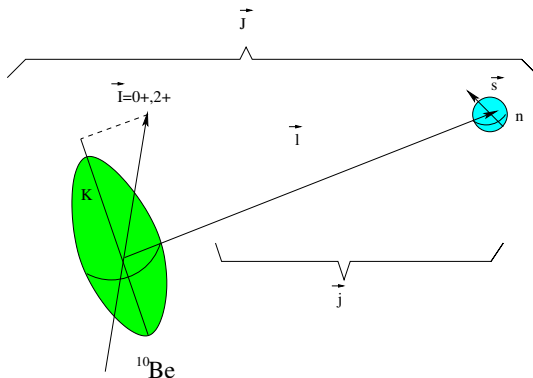
## An example is: Particle-Rotor Model



Set of  $\mathcal{L}^2$  functions in this scheme:

$$|\phi_{i,J_p}(\vec{r}, \vec{\xi})\rangle = \sum_{\alpha} R_{i,\alpha}^{THO}(r) \left[ \mathcal{Y}_{l s j}(\hat{r}) \otimes \varphi_I(\vec{\xi}) \right]_{J_p} \quad i = 1, \dots, N$$

## An example is: Particle-Rotor Model



The valence-core interaction in first order:

$$V_{NC}(\vec{r}, \vec{\xi}) = V_{NC}(r - R_0 - \Delta(\hat{\Omega})) \approx V_{Cent}(r) - \delta_2 \frac{dV_{Cent}}{dr} Y_{20}(\hat{r})$$

# Pseudo-states (PS) discretization method

- Discrete set of  $\mathcal{L}^2$  functions:  $|\phi_n\rangle$

Completeness condition:

$$\sum^N |\phi_i\rangle\langle\phi_i| \approx \mathbf{I}$$

- To diagonalize the internal Hamiltonian of a projectile  $\mathcal{H}_p$

Matrix elements:

$$\mathcal{H}_p \mapsto \sum_{n,n'} |\phi_n\rangle\langle\phi_n|\mathcal{H}_p|\phi_{n'}\rangle\langle\phi_{n'}|$$

# Pseudo-states (PS) discretization method

Eigenstates of the matrix  $N \times N$ :

$$|\varphi_n^{(N)}\rangle = \sum^N C_i^n |\phi_i\rangle$$

- $\left\{ \begin{array}{l} n_b \text{ states with } \varepsilon_n < 0 \text{ representing the bound states.} \\ N - n_b, \varepsilon_n > 0 \Rightarrow \text{discrete representation of the Continuum} \end{array} \right.$
- Orthogonal and normalizable.

**What is the most suitable basis? Lagrange, Sturmian, Harmonic Oscillator?**



# Transformed Harmonic Oscillator basis

Analytic LST from Karataglidis *et al.*, PRC71,064601(2005)

$$s(r) = \frac{1}{\sqrt{2}b} \left[ \frac{1}{\left(\frac{1}{r}\right)^m + \left(\frac{1}{\gamma\sqrt{r}}\right)^m} \right]^{\frac{1}{m}}$$

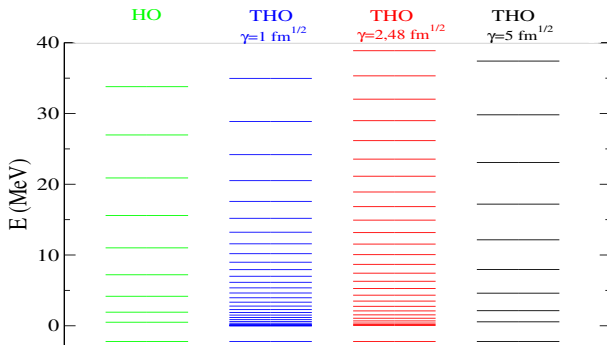
HO vs THO:

$$\phi(s) \mapsto e^{-\left(\frac{s}{b}\right)^2} \implies \phi[s(r)] \mapsto e^{-\frac{\gamma^2}{2b^2}r}$$

- Correct asymptotic behaviour for bound states.
- Range controlled by the parameters of the LST.

# THO parameters

- $b$  is treated as a variational parameter to minimize g.s. energy
- Then  $\frac{\gamma}{b}$  controls the density of states:

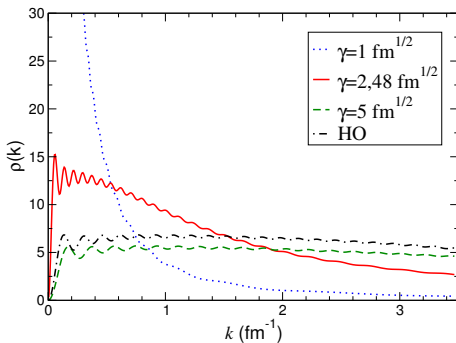


- $\gamma$  can be also used to look for resonances

# Energy distribution of pseudo-states

## Density of states

$$\rho_l^{(N)}(k) = \sum_{n=1}^N \langle \varphi_l(k) | \varphi_{n,l}^{(N)} \rangle$$



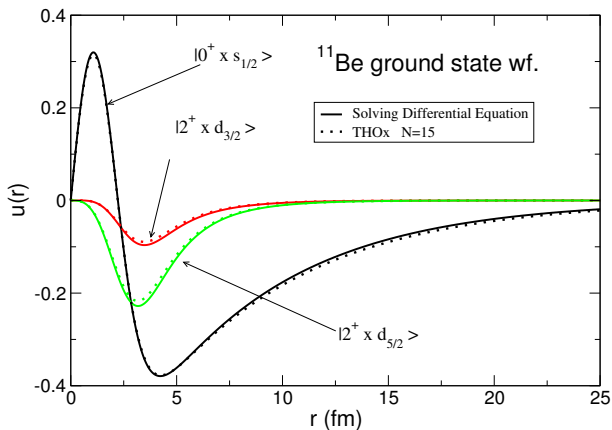
# The Smoothing Process

For any operator

$$\begin{aligned} O(\varepsilon) &= \langle k | \hat{O} | g.s. \rangle \\ &= \langle k | \sum_n | n \rangle \langle n | \hat{O} | g.s. \rangle \\ &= \sum_n \rho^{(n)}(k) \langle n | \hat{O} | g.s. \rangle \end{aligned}$$

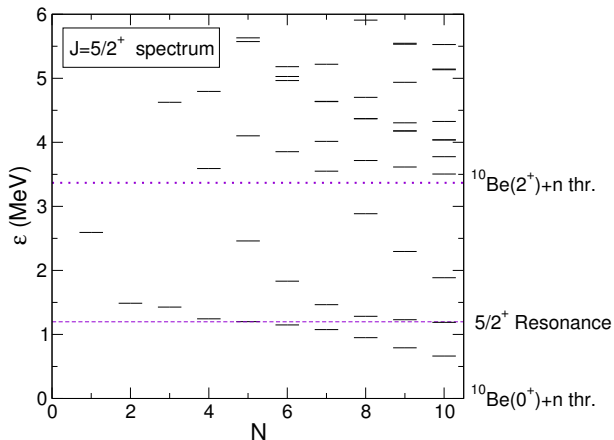
⇒ A continuous distribution in energy from discrete values of S matrix,  $B(E\lambda)$ , cross sections...

# Particle-Rotor Model

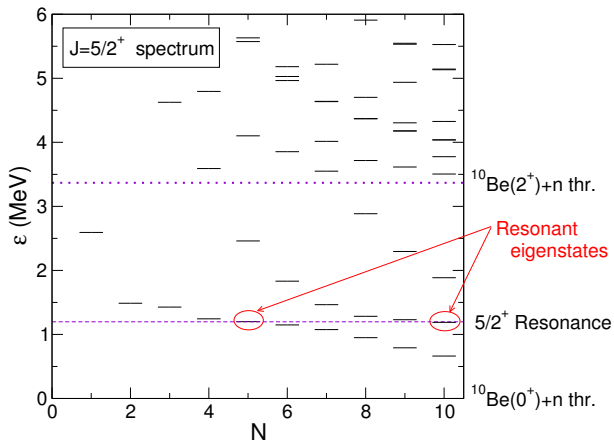


The g.s. wavefunction is well described using a small THO basis

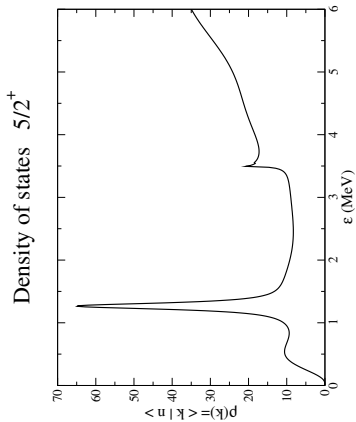
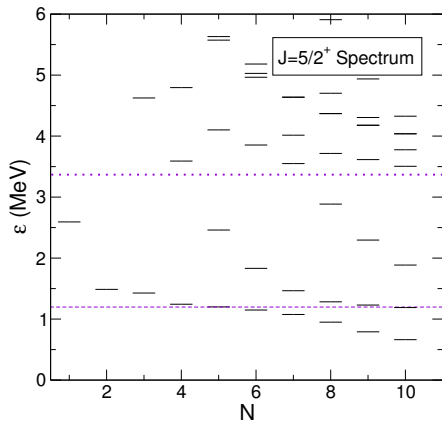
# Finding Resonances



# Finding Resonances

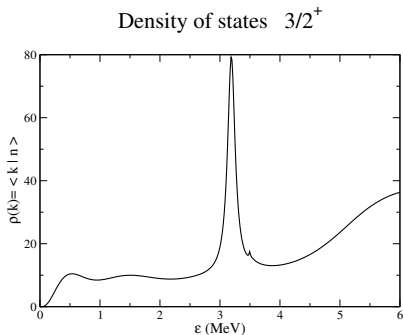
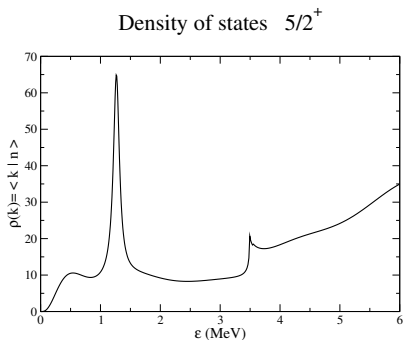


## Spectrum





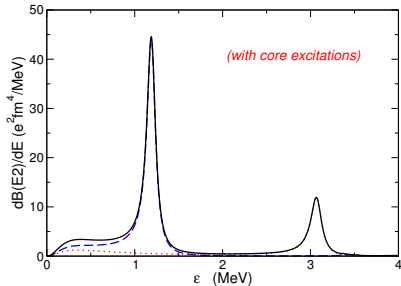
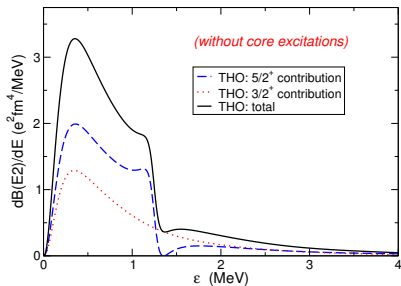
# Energy distribution of pseudo-states



⇒ Widths of 125 KeV and 140 KeV respectively.

⇒ Experimental values of  $100 \pm 10$  KeV and  $122 \pm 8$  KeV.

# Electromagnetic Transition Probabilities



$\Rightarrow$  B(E2) dominated by collective excitation of the core

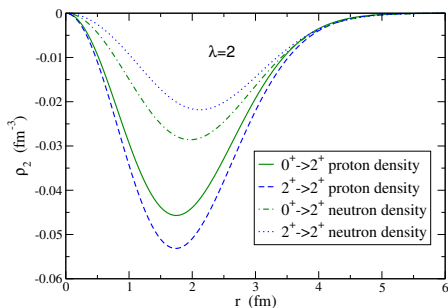
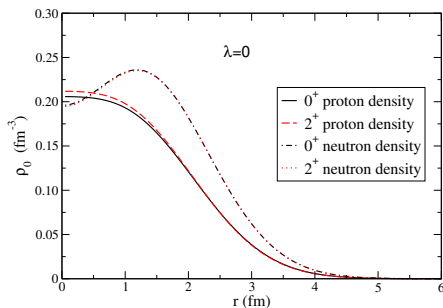
# PRM "drawbacks"

## PRM needs:

- The core to be a rotor
- A phenomenological potential based on the following parameters:

$$E(2^+), \beta_2, V_c, r, a, V_{so}, r_{so}, a_{so}$$

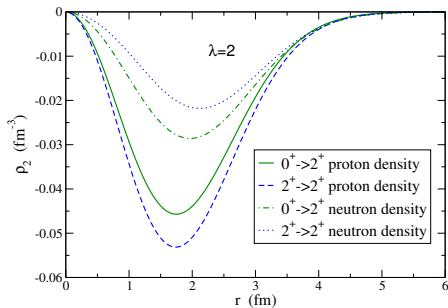
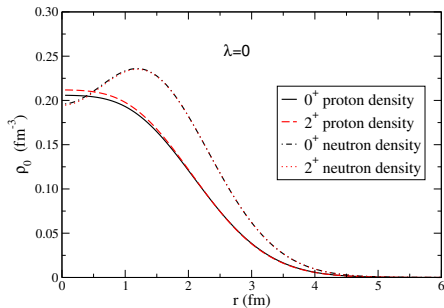
## P-AMD



## Densities from Antisymmetrized Molecular Dynamics (AMD)

Y. Kanada-En'yo *et al.* Phys. Rev. C 60, 064304 (1999)

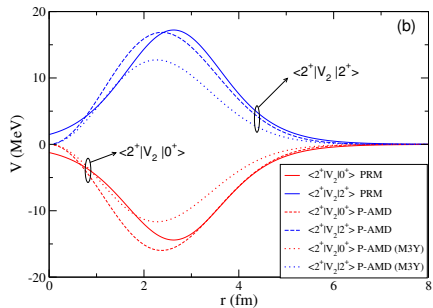
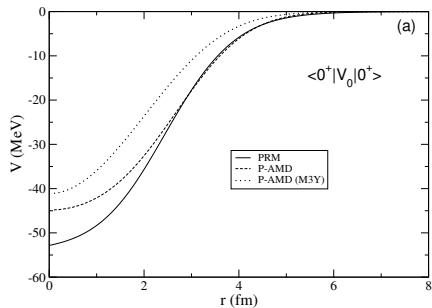
## P-AMD



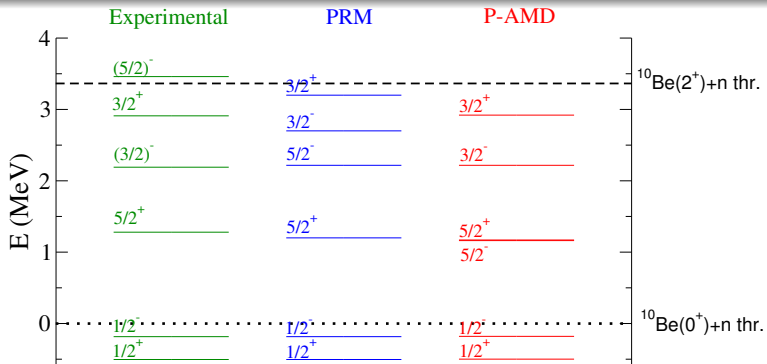
$$\langle I || V_{NC}^\lambda(r, \vec{\xi}) || I' \rangle = \int dr' \left[ \langle I || \rho_\lambda(r', \xi) || I' \rangle v_{nn}(|\vec{r} - \vec{r}'|) \right]$$

JLM interaction Phys. Rev. C 16, 80 (1977).

## P-AMD



## P-AMD



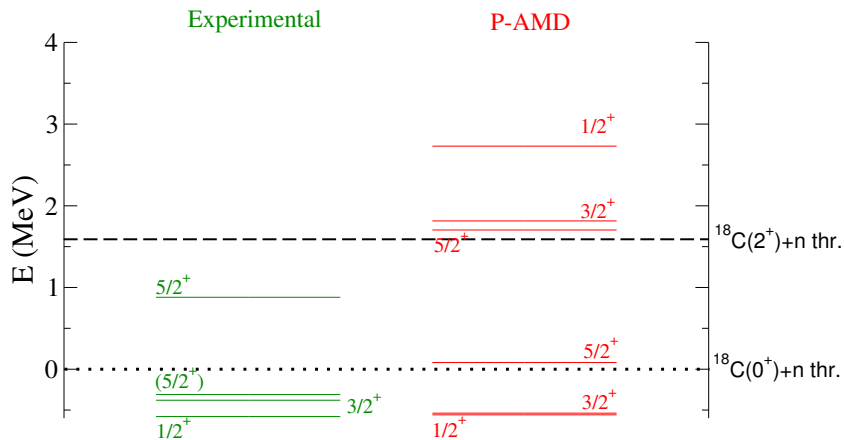
## Renormalization factors

$$\lambda_+ = 1.058 \text{ and } \lambda_- = 0.995$$

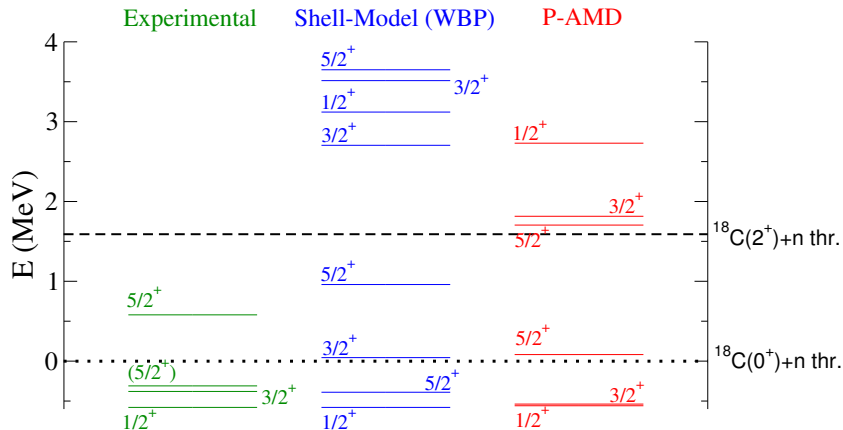
PRC 70, 054606 (2004); PRC 81, 034321 (2010); PL B 611, 239 (2005).

State	Model	$ 0^+ \otimes (ls)j\rangle$	$ 2^+ \otimes s_{1/2}\rangle$	$ 2^+ \otimes d_{3/2}\rangle$	$ 2^+ \otimes d_{5/2}\rangle$
$1/2^+$	PRM	0.857	–	0.021	0.121
	P-AMD	0.849	–	0.031	0.121
	WBT	0.762	–	0.002	0.184
$5/2^+$	PRM	0.702	0.177	0.009	0.112
	P-AMD	0.674	0.189	0.014	0.124
	WBT	0.682	0.177	0.009	0.095
$3/2^+$	PRM	0.165	0.737	0.017	0.081
	P-AMD	0.316	0.565	0.031	0.089
	WBT	0.068	0.534	0.008	0.167

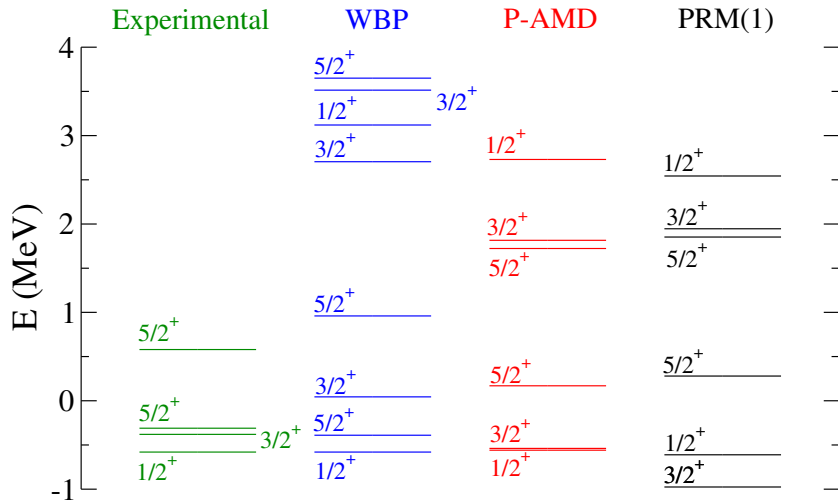


$^{19}\text{C}$  Spectrum

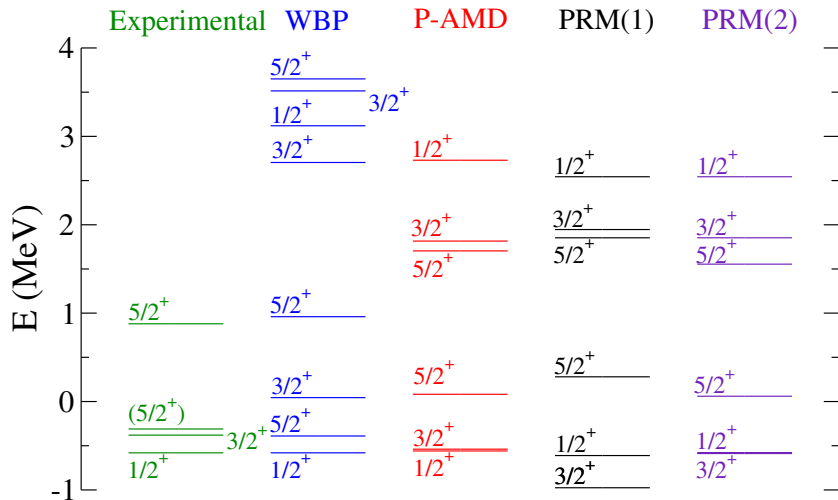
PL B 660, 320 (2008); PL B 614, 174 (2005).

$^{19}\text{C}$  Spectrum

PL B 660, 320 (2008); PL B 614, 174 (2005).

$^{19}\text{C}$  Spectrum

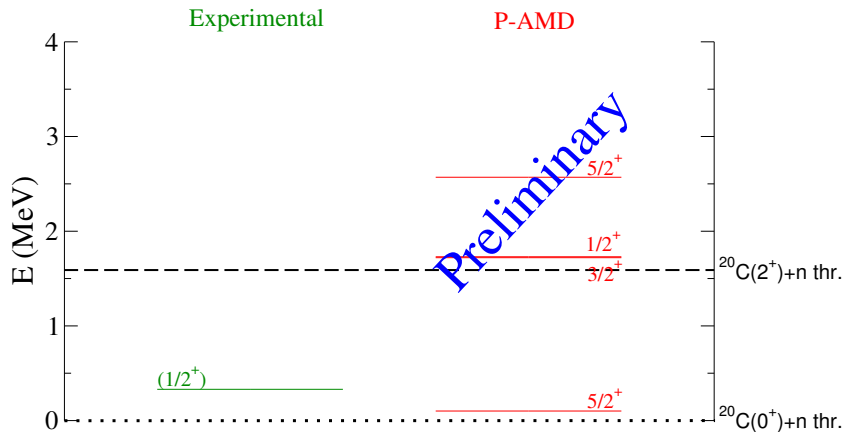
PL B 660, 320 (2008); PL B 614, 174 (2005).

$^{19}\text{C}$  Spectrum

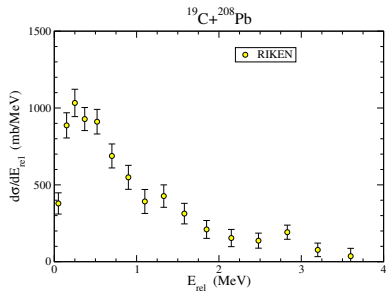
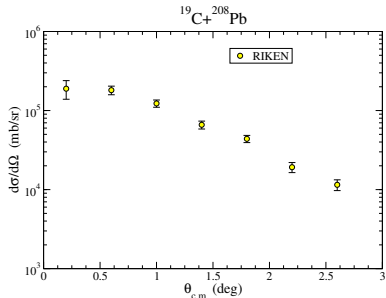
PL B 660, 320 (2008); PL B 614, 174 (2005).

Nucleus	rms (fm)	$\langle 0    \hat{\delta}_2    2 \rangle_{\text{AMD}}$ (fm)	$\langle 0    \hat{\delta}_2    2 \rangle_{\text{rot}}$ (fm)
$^{10}\text{Be}$	2.538	1.90	1.66
$^{18}\text{C}$	2.776	1.20	1.50

State	Model	$ 0^+ \otimes (\ell s)j\rangle$	$ 2^+ \otimes s_{1/2}\rangle$	$ 2^+ \otimes d_{3/2}\rangle$	$ 2^+ \otimes d_{5/2}\rangle$
$1/2_1^+$	P-AMD	0.529	–	0.035	0.436
	PRM(1)	0.517	–	0.081	0.402
	PRM(2)	0.505	–	0.033	0.462
	WBP	0.600	–	0.002	0.184
$5/2_1^+$	P-AMD	0.276	0.721	0.000	0.003
	PRM(1)	0.285	0.716	0.000	0.003
	PRM(2)	0.278	0.719	0.000	0.003
	WBP	0.383	0.015	0.000	0.751
$5/2_2^+$	P-AMD	0.200	0.142	0.002	0.657
	PRM(1)	0.217	0.178	0.004	0.602
	PRM(2)	0.207	0.100	0.002	0.690
	WBP	0.035	0.609	0.009	0.291

$^{21}\text{C}$  Spectrum

PRC86, 054604

$^{19}\text{C} + ^{208}\text{Pb}$  @ 67 MeV/u

- Coulomb Excitation

T. Nakamura *et al.*, Phys. Rev. Lett. 83, 1112 (1999).



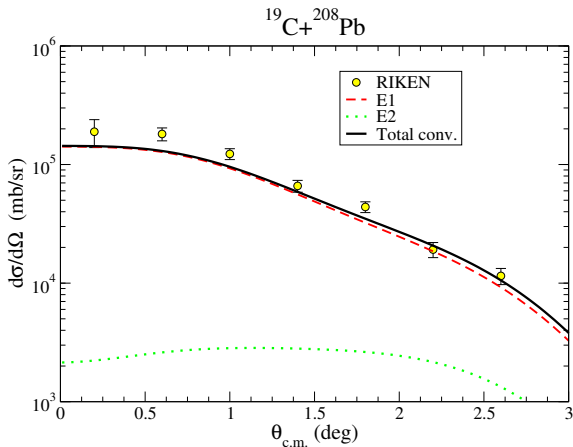
$^{19}\text{C} + ^{208}\text{Pb}$  @ 67 MeV/u

# Semiclassical calculations

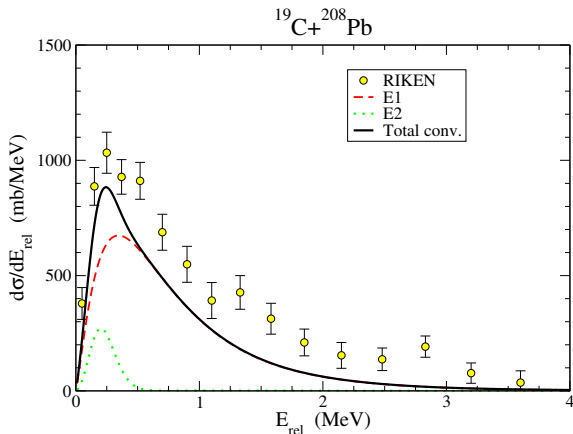
## Equivalent Photon Method

- ⇒ Only includes first order Coulomb excitation.
- ⇒ Actually this is only a first test before using the basis with DWBA<sub>x</sub>/XCDCC.

$$\left. \frac{d\sigma_\lambda}{d\Omega dE} \right|_{bu} = \frac{4\pi^3}{9} \frac{dB(E1)}{dE} \frac{dN_{E1}}{d\Omega}$$

$^{19}\text{C} + ^{208}\text{Pb}$  @ 67 MeV/uT. Nakamura *et al.*, Phys. Rev. Lett. 83, 1112 (1999).

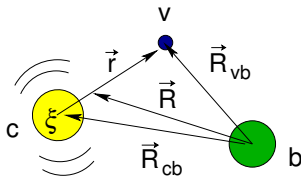
- The reaction is dominated by E1 first order Coulomb excitation as expected

$^{19}\text{C} + ^{208}\text{Pb}$  @ 67 MeV/u

- Resonant E2 contribution more important due to its low excitation energy

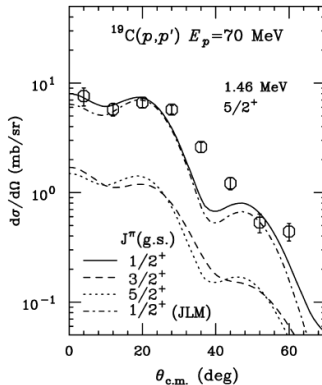
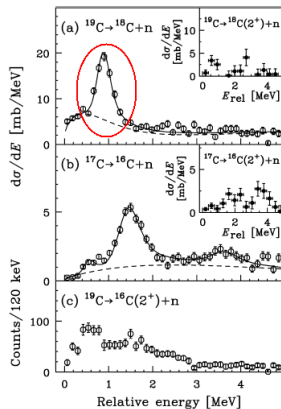
$^{19}\text{C}+p @ 67 \text{ MeV}/u$ 

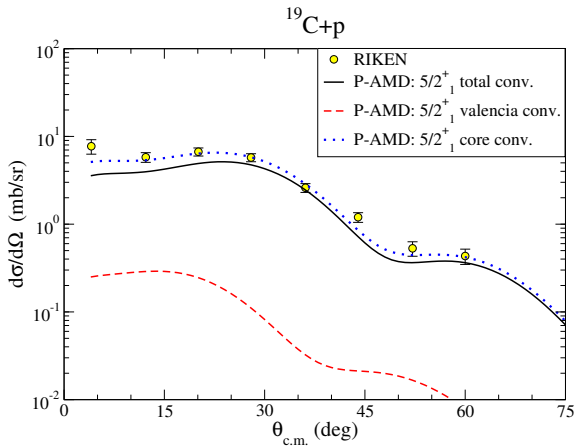
## DWBAx calculations

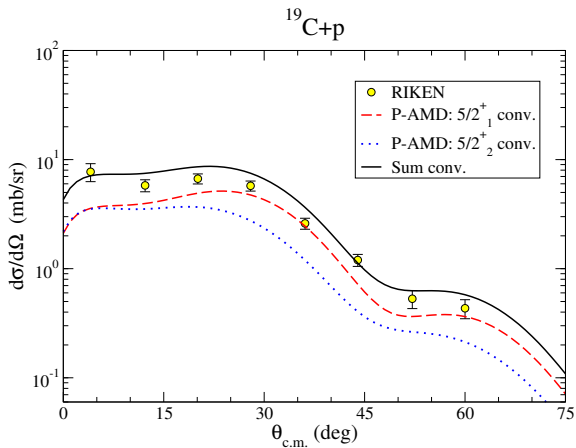


## No-recoil approach

- ⇒ Only first order excitation.
- ⇒ Core and Valence contributions explicitly separates in the calculation.  
A. M. Moro & R. Crespo, Phys. Rev. C 85, 054613 (2012).
- ⇒ Same results for these energies than XCDCC.  
A. M. Moro *et al.* AIP Conf. Proc. 1491, 335 (2012)

$^{19}\text{C} + p @ 67 \text{ MeV}/u$ Y. Satou *et al.*, Phys. Lett. B 660, 320 (2008).Microscopic DWBA calculations suggest a  $1/2^+ \Rightarrow 5/2^+$  transition

$^{19}\text{C}+p$  @ 67 MeV/u

$^{19}\text{C}+p$  @ 67 MeV/u

## Suitable discrete description of the continuum

The generalization of the **THO basis** for this case continues giving:

- A converged description of the continuum with a reduced number of functions.
- Natural and **accurate treatment of narrow resonances**.

## Particle-Rotor Model

Accurate description of even-odd halo nuclei within deformed regions like  $^{11}\text{Be}$  and  $^{19}\text{C}$ .

## P-AMD

- Accurate **semi-microscopic description** of even-odd halo nuclei.
- Predictive power for unknown halo nuclei like  $^{19,21}\text{C}$ .
- Could be able to include **core excitations from different sources**.



## Application to reactions with heavy targets

- **Core excitations** may **interfere** the extraction of  $B(E1)$  distributions.
- **Low lying quadrupole resonances** can be populated through core excitation in these reactions.

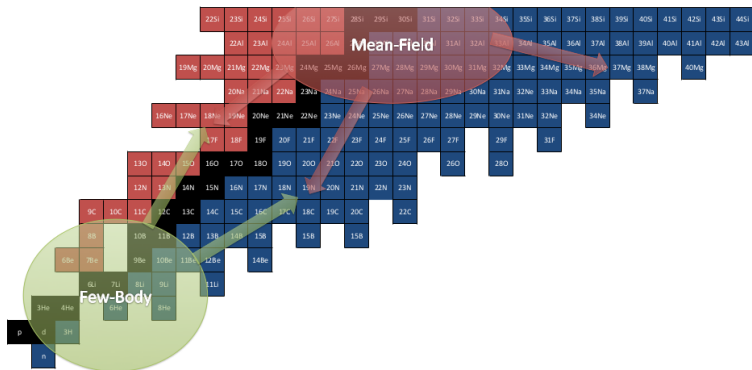
## Application to reactions within DWBAx framework

- The **interplay** between **core and valence contributions** is crucial to understand angular distribution of the break up of halo nuclei.
- **Break up reactions** are sensitive to spectroscopic factors of resonant states difficult to populate in traditional transfer reactions.

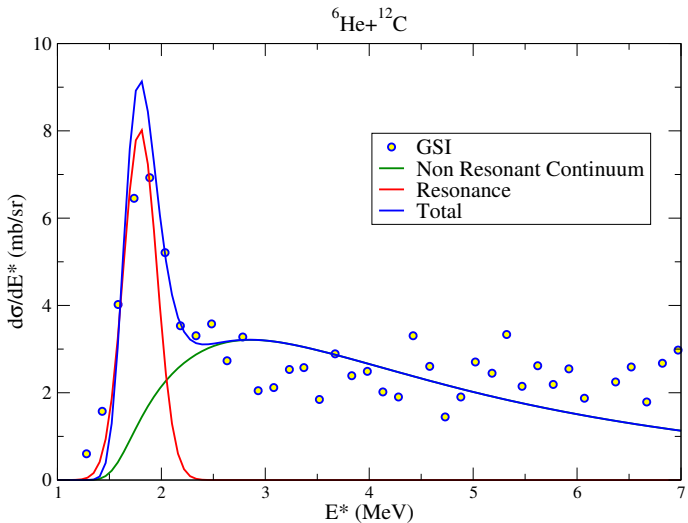
## Next steps

- Apply the P-AMD model to other halo nuclei.
- Check the use of other microscopic densities.
- Perform full PS-XCDCC calculations with the THO basis and the P-AMD model.
- Use the density distributions of the PS to obtain energy distribution of the cross sections.

- 1 To add mean field features to few-body approaches (already with a right treatment of the Continuum)
- 2 To add the knowledge about Continuum to mean-field approaches



- 3 This nuclear structure should be included in nuclear reactions frameworks in order to analyse a wider variety of experimental data



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## Spectrum

