

# Continuum effects: Structure and reactions of ${}^6\text{He}$

M. Rodríguez-Gallardo<sup>1,2,3 a</sup>, J. M. Arias<sup>2</sup>, J. Gómez-Camacho<sup>2</sup>, R. C. Johnson<sup>3</sup>, A. M. Moro<sup>2</sup>, I. J. Thompson<sup>3</sup>, and J. A. Tostevin<sup>3</sup>

<sup>1</sup> Centro de Física Nuclear, Universidade de Lisboa, Av. Prof. Gama Pinto 2, 1649-003 Lisboa, Portugal

<sup>2</sup> Departamento de Física Atómica, Molecular y Nuclear, Universidad de Sevilla, Apartado 1065, 41080 Sevilla, Spain

<sup>3</sup> Department of Physics, University of Surrey, Guildford GU2 7XH, United Kingdom

Received: date / Revised version: date

**Abstract.** A description of three-body weakly bound systems using the Transformed Harmonic Oscillator (THO) method is addressed. First, relevant structure observables are presented for the Borromean nucleus  ${}^6\text{He}$ . Then, the THO method is applied to the study of  ${}^6\text{He}$  scattering within the Continuum-Discretized Coupled Channels (CDCC) framework.

**PACS.** 21.45.+v – 21.10.-k – 27.20.+n – 24.10.-i – 24.10.Eq – 25.60.-t – 25.60.Bx

## 1 Introduction

A particularly interesting example of weakly bound systems is that of Borromean nuclei, that is, three-body composite systems with no binary bound subsystems. For an appropriate description of the reactions induced by a Borromean nucleus, like  ${}^6\text{He}$ , the introduction of the continuum part of the spectrum is necessary. For this purpose it is required to discretize the three-body continuum. This work has been developed by the group of M. Kamimura [1, 2] using a discretization method based on Gaussian functions. Here, an alternative discretization method, based on Hyperspherical Harmonics and orthogonal polynomials, is presented.

## 2 Three-body continuum discretization method: Application to ${}^6\text{He}$

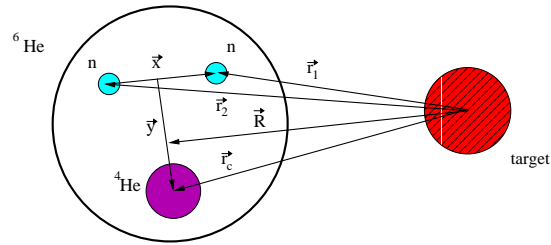
This formalism is a generalization of the method presented in Ref. [4] for a two-body system. For a three-body case the hyperspherical coordinates  $\{\rho, \alpha, \hat{x}, \hat{y}\}$  will be used. They are obtained from the Jacobi coordinates  $\{\mathbf{x}, \mathbf{y}\}$  (see, for instance, Ref. [8]). The Jacobi coordinates are presented in Fig. 1.

First, the wave functions of the system are expanded in Hyperspherical Harmonics (HH) as

$$\Psi_{\beta j \mu}(\rho, \Omega) = R_{\beta j}(\rho) \sum_{m \sigma} \langle l m S_x \sigma | j \mu \rangle \Upsilon_{K l m}^{l_x l_y}(\Omega) \chi_{S_x}^{\sigma} \quad (1)$$

where  $R_{\beta j}(\rho)$  is the hyperradial wave function,  $\chi_{S_x}^{\sigma}$  the spin wave function of the two particles related by the co-

<sup>a</sup> Present address: Centro de Física Nuclear, Universidade de Lisboa



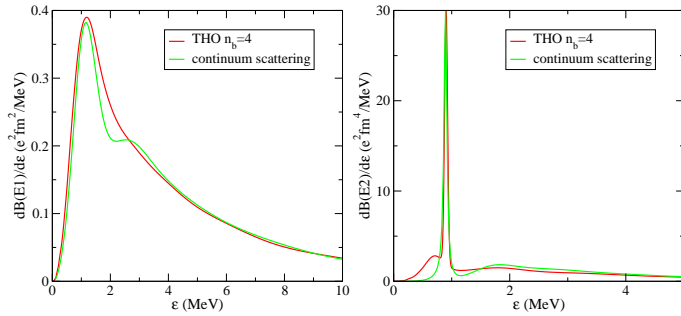
**Fig. 1.** Jacobi coordinates;  ${}^6\text{He}$  scattered by a target.

ordinate  $\mathbf{x}$ , and  $\Upsilon_{K l m}^{l_x l_y}(\Omega)$  are the hyperspherical harmonics that depend on the angular variables  $\Omega \equiv \{\alpha, \hat{x}, \hat{y}\}$ . The set of quantum numbers  $\beta \equiv \{K, l_x, l_y, l, S_x\}$  defines each channel. Here,  $K$  is the hypermomentum,  $l_x$  and  $l_y$  are the orbital angular momenta associated with the Jacobi coordinates  $\mathbf{x}$  and  $\mathbf{y}$ ,  $l = l_x + l_y$  is the total orbital angular momentum,  $S_x$  is the spin of the particles related by the coordinate  $\mathbf{x}$  and  $j = l + S_x$  is the total angular momentum.

Then the Transformed Harmonic Oscillator method (THO) [3] is used to obtain  $R_{\beta j}(\rho)$ . The basic idea of the THO method [4–7] is to convert the bound ground-state wave function of the system into the ground-state wave function of the Harmonic Oscillator (HO), defining a Local Scale Transformation (LST). Considering that the ground-state wave function can be written as a linear combination of the basis functions given in Eq. (1), the equation that defines the LST for every channel  $\beta$  is

$$\int_0^{\rho} d\rho' \rho'^5 |R_{B\beta}(\rho')|^2 = \int_0^s ds' s'^5 |R_{0K}^{HO}(s')|^2, \quad (2)$$

where  $R_{0K}^{HO}(s)$  is the hyperradial wave function of the HO. Finally, the THO basis is constructed for each channel



**Fig. 2.** (left)  $B(E1)$  distribution for  $n_b = 4$ . (right)  $B(E2)$  distribution for  $n_b = 4$ .

applying the LST,  $s_\beta(\rho)$ , to the HO basis

$$R_{i\beta}^{THO}(\rho) = R_{B\beta}(\rho)L_i^{K+2}(s_\beta(\rho)^2), \quad (3)$$

where  $L_i^\lambda(t)$  are Laguerre polynomials. In this way, a complete and orthonormal basis is obtained. Here the quantum number  $i$  denotes the hyperradial excitations. Note that as  $i$  increases, the functions  $R_{i\beta}^{THO}(\rho)$  become more oscillatory and explore larger distances.

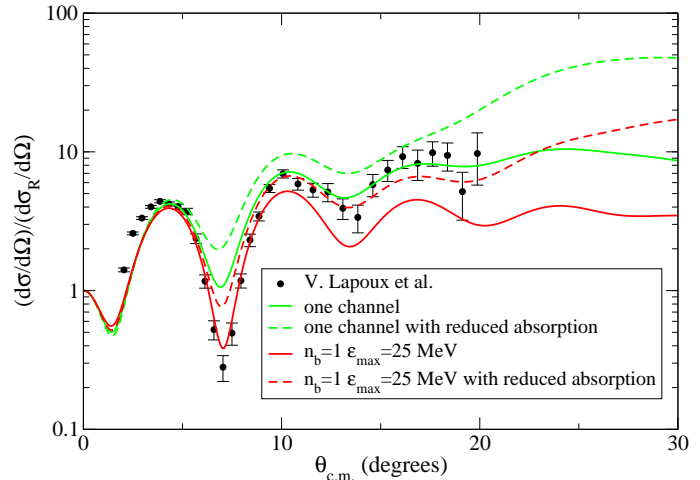
Next we apply the method to the Borromean nucleus  ${}^6\text{He}$ . In that case the Hamiltonian (H) is detailed in [8] and includes two-body interactions plus an effective three-body potential and a treatment of Pauli principle using a pseudopotential. H is diagonalized in a finite THO basis with  $N$  states (including  $n_b$  hyperradial excitations in every channel so  $N = (n_b + 1) \times nchan$ , where  $nchan$  is the number of channels for each state).

For  $j = 0^+$  we get almost exactly the ground-state of the system. For  $j = 2^+$  and  $n_b < 5$ , we obtain a very stable state around the experimental energy of the  $2^+$  resonance for  ${}^6\text{He}$ .

In Fig. 2 we present the  $B(E1)$  and  $B(E2)$  distributions for  $n_b = 4$  as a function of the excitation energy on the continuum. For comparison, the distribution calculated with the continuum scattering wave functions, taken from Ref. [9], is also shown. For both cases,  $B(E1)$  and  $B(E2)$ , the two distributions are in good agreement with the same total strength. For the  $B(E2)$  distribution the most relevant feature is the presence of a narrow low-lying resonance.

### 3 Reactions induced by ${}^6\text{He}$

In Fig. 1 we show a scheme of the scattering of  ${}^6\text{He}$  by a target, where  $\mathbf{R}$  is the position of the center of mass of  ${}^6\text{He}$  from the target. In order to perform CDCC [10] calculations, we take as the projectile basis, the eigenstates obtained from the diagonalization of H in a THO basis with a given numbers of states ( $N$ ). Then it is necessary to calculate the coupling potentials  $V_{Lnj, L'n'j'}^J(\mathbf{R}) = \langle LnjJM | V_{1t}(\mathbf{r}_1) + V_{2t}(\mathbf{r}_2) + V_{ct}(\mathbf{r}_c) | L'n'j'JM \rangle$ . These potentials are read externally in the program FRESKO [11], that solves the system of coupled equations as in the conventional CDCC method. As a preliminary result we show



**Fig. 3.** Differential elastic cross section relative to Rutherford cross section for the reaction  ${}^6\text{He}+{}^{12}\text{C}$  at 229.8 MeV.

in Fig. 3 the differential elastic cross section relative to Rutherford cross section for the reaction  ${}^6\text{He}+{}^{12}\text{C}$  at 229.8 MeV. We can see how the calculation with a very small THO basis reproduces quite well the experimental data of V. Lapoux et al. [12]. In order to reproduce satisfactory the normalization of the data for angles beyond  $10^\circ$  we had to reduce the imaginary part of the coupling potentials by around 40%. A similar reduction was also required in the work [1].

The purpose, in the next future, is the application of this procedure to analyze scattering cross sections of halo nuclei.

### Acknowledgments

This work was supported in part by the DGICYT under projects FIS2005-01105 and FPA 2005-04460. M.R.G. acknowledges a FCT grant POCTI/ISFL/2/275 and FCT under the grant POCTI/FIS/43421/2001.

### References

1. T. Matsumoto et al., Phys. Rev. C **70** (2004) 061601.
2. T. Matsumoto et al., Phys. Rev. C **73** (2006) 051602(R).
3. I. Zh. Petkov and M. V. Stoitsov, Compt. Rend. Bulg. Acad. Sci. **34** (1981) 1651.
4. F. Pérez-Bernal et al., Phys. Rev. A **63** (2001) 052111.
5. I. Martel et al., Phys. Rev. A **65** (2002) 052708.
6. M. Rodríguez-Gallardo et al., Phys. Rev. C **69** (2004) 034308.
7. M. Rodríguez-Gallardo et al., Phys. Rev. C **72** (2005) 024007.
8. I. J. Thompson et al., Comput. Phys. Commun. **161** (2004) 87.
9. I. J. Thompson et al., Phys. Rev. C **61** (2000) 24318.
10. N. Austern et al., Phys. Rep. **154** (1987) 125.
11. I. J. Thompson, Comput. Phys. Rep. **7** (1988) 167.
12. V. Lapoux et al., Phys. Rev. C **66** (2002) 034608.