

## Competition between long- range collective and short range pairing correlations in two-neutron transfer reactions

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## Outline

Nuclear spectroscopy via transfer reactions between heavy ions
$>$ The $\left({ }^{18} \mathrm{O},{ }^{16} \mathrm{O}\right)$ reaction

$>$ CRC and two-step DWBA calculations
> Microscopic cluster calculations

## Nuclear spectroscopy via $\left(\mathbf{1 8}^{\mathbf{1 8}} \mathbf{0},{ }^{\mathbf{1 6}} \mathbf{0}\right)$ reaction

The $\left({ }^{18} 0,{ }^{16} \mathrm{O}\right)$ reactions are good candidates to show the role of pairing interaction thanks to
$>$ The presence of a correlated pair of neutrons in the ${ }^{18} \mathrm{O}_{\text {g.s. }}$ wave function
> The very low polarizability of the ${ }^{16} \mathrm{O}$ core
${ }^{14} \mathrm{C}$ is a good benchmark for considerations
on the reaction mechanism, ${ }^{64} \mathrm{Ni}$ and ${ }^{28}$ Si are good benchmark for studying long-range vs short-range correlations

Studies on both
${ }^{13} \mathrm{C}\left({ }^{18} \mathrm{O},{ }^{17} \mathrm{O}\right){ }^{14} \mathrm{C}$ 1n transfer and ${ }^{12} \mathrm{C}\left({ }^{18} \mathrm{O},{ }^{16} \mathrm{O}\right){ }^{14} \mathrm{C} \mathbf{2 n}$ transfer


- Presence of 2 n correlations in the ${ }^{14} \mathrm{C}_{\text {g.s. }}$ wave function
- Strong selectivity in the populated states
- Absolute cross sections reproduced without any scaling factor
M. Cavallaro et al., PRC 88 (2013) 054601


## Theoretical models and main ingredients

## Exact finite range CRC and two-step CCBA calculations

> Sao Paulo Potential (SPP) used in the optical model

$$
\text { L.C. Chamon, et al., PRL } 79 \text { (1997) } 5218
$$

> Wood-Saxon form factors were used to generate single particle and cluster wave functions. Depth were adjusted to fit the exp. separation energies
> Deformation parameters for collective excitations
> Spectroscopic Amplitudes by shell-model in the $1 p_{1 / 2}, 1 d_{5 / 2}, 2 \mathrm{~s}_{1 / 2}$ model space (zbm interaction)
A.P. Zuker, et al., PRL 17 (1969) 983

## Theoretical models and main ingredients

The CRC equations are in many cases of the form

$$
\begin{aligned}
{\left[E_{\kappa p t}-T_{\kappa L}\left(R_{\kappa}\right)-U_{\kappa}\left(R_{\kappa}\right)\right] J_{\alpha}\left(R_{\kappa}\right) } & =\sum_{\alpha^{\prime}, \Gamma>0} i^{L^{\prime}-L} V_{\alpha: \alpha^{\prime}}^{\Gamma}\left(R_{\kappa^{\prime}}\right) J_{\alpha^{\prime}}\left(R_{\kappa^{\prime}}\right) \\
& +\sum_{\alpha^{\prime}, \kappa^{\prime} \neq \kappa} i^{L^{\prime}-L} \int_{0}^{R_{m}} V_{\alpha: \alpha^{\prime}}\left(\left(R_{\kappa}\right), R_{\kappa^{\prime}}\right) J_{\alpha^{\prime}}\left(R_{\kappa^{\prime}}\right) d R_{\kappa^{\prime}}
\end{aligned}
$$

Single nucleon states are given by

$$
\begin{aligned}
& \phi_{J M}\left(\xi_{c}, \mathbf{r}\right)=\sum_{\ell j I} A_{\ell s j}^{j I J}\left[\phi_{I}\left(\xi_{c}\right) \varphi_{\ell s j}(\mathbf{r})\right]_{J M} \\
= & \sum_{\ell j I, m \mu m_{s} m_{\ell}} A_{\ell s j}^{j I J}\langle j m I \mu \mid J M\rangle \phi_{I \mu}\left(\xi_{c}\right)\left\langle\ell m_{\ell} s m_{s} \mid j m\right\rangle Y_{\ell}^{m_{\ell}}(\hat{\mathbf{r}}) \phi_{s}^{m_{s}} \frac{1}{r} u_{\ell s j I}(r)
\end{aligned}
$$

and are the solution of

$$
\left[T_{\ell}(r)+V(r)+\epsilon_{I}-E\right] u_{\ell s j I}(r)+\sum_{\ell^{\prime} j^{\prime} I^{\prime}, \Gamma>0} V_{\ell s j I \ell^{\prime} s j^{\prime} I^{\prime}}^{\Gamma}(r) u_{\ell^{\prime} s j^{\prime} I^{\prime}}(r)=0
$$

## Theoretical models and main ingredients

Independent coordinate model

$$
\begin{aligned}
\varphi_{12}\left(\mathbf{r}_{1}, \mathbf{r}_{2}\right)= & \sum_{i} c_{i}\left|\left(\ell_{1}(i), s_{1}\right) j_{1}(i),\left(\ell_{2}(i), s_{2}\right) j_{2}(i) ; J_{12} T\right\rangle \\
& \rightarrow \sum_{u} c_{i} \sum_{L \ell S j}\left|L,\left(\ell,\left(s_{1} s_{2}\right) S^{r}\right) j ; J_{12} T\right\rangle \phi_{L(\ell S) j}^{J_{12} T, i}(r, \rho) \\
\phi_{L(\ell S) j}^{J_{12} T, i}(r, \rho)= & \left\langle L,\left(\ell,\left(s_{1} s_{2}\right) S\right) j ; J_{12} T \mid\left(\ell_{1}(i), s_{1}\right) j_{1}(i),\left(\ell_{2}(i), s_{2}\right) j_{2}(i) ; J_{12} T\right\rangle \\
& \times\left\langle\left[Y_{L}(\hat{\mathbf{r}}) Y_{\ell}(\hat{\rho})\right]_{\lambda} \mid\left[\varphi_{\ell_{1} s_{1} j_{1}}\left(\mathbf{r}_{1}\right) \varphi_{\ell_{2} s_{2} j_{2}}\left(\mathbf{r}_{2}\right)\right]_{J_{12} T}\right\rangle
\end{aligned}
$$

and the radial integral overlaps are derived from using Moshinsky harmonic oscillator expansion

## Theoretical results for other channels



Presence of two-neutron pairing correlations in other ${ }^{14} \mathrm{C}$ states
M. Cavallaro et al., PRC 88 (2013) 054601

## Extreme Cluster Model

(CRC)

* Relative motion of the $2 n$ system frozen and separated by the c.m.
* Only the term with the $2 n$ coupled to $S=0$ participates to the transfer


## Sequential transfer

 (DWBA)Introducing the ${ }^{17} \mathrm{O}+{ }^{13} \mathrm{C}$ intermediate partition

## I ndependent coord.

(CRC)

## CRC - 1 n transfer

## No arbitrary scaling

## New works published in 2016-2017

What happens if we add a neutron to the ${ }^{14} \mathrm{C}$ system?

## Study of the ${ }^{13} \mathrm{C}\left({ }^{18} \mathrm{O},{ }^{16} \mathrm{O}\right){ }^{15} \mathrm{C}$ reaction at 84 MeV incident energy

D. Carbone et al., PRC 95, 034603 (2017)

## ${ }^{15} \mathrm{C}$ energy spectrum


> Same states populated in the (t,p) reactions
$>$ Strong population of states with ${ }^{13} C+2 n$ configurations
$>$ Population of the Giant Pairing Vibration above $S_{2 n}$

- F. Cappuzzello et al., Nat. Commun. 6, 6743 (2015)
- D. Carbone, EPJ Plus (2015) 130:143

$$
9^{\circ}<\theta_{\text {lab }}<10^{\circ}
$$

Energy resolution ~ 200 keV

## ARTICLE

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## Signatures of the Giant Pairing Vibration in the ${ }^{14} \mathrm{C}$ and ${ }^{15} \mathrm{C}$ atomic nuclei

F. Cappuzzello ${ }^{1,2}$, D. Carbone ${ }^{2}$, M. Cavallaro ${ }^{2}$, M. Bondi ${ }^{11,2}$, C. Agodi ${ }^{2}$, F. Azaiez ${ }^{3}$, A. Bonaccorso ${ }^{4}$, A. Cunsolo ${ }^{2}$,
L. Fortunato ${ }^{5,6}$, A. Foti ${ }^{1,7}$, S. Franchoo ${ }^{3}$, E. Khan ${ }^{3}$, R. Linares ${ }^{8}$, J. Lubian ${ }^{8}$, J.A. Scarpaci ${ }^{9}$ \& A. Vitturi ${ }^{5,6}$



Table 1 | Main spectroscopic features of the populated states.

| S.No. | Excitation energy <br> (MeV) (present <br> work) | Excitation energy <br> (MeV) (values <br> from ref. 38) | $\boldsymbol{J}^{\boldsymbol{\pi}} \mathbf{(}^{\star}$ ) |
| :--- | :---: | :---: | :---: |
| 15 C states |  |  |  |
| 1 | $0.00 \pm 0.02$ | 0 | $1 / 2^{+}$ |
| 2 | $0.73 \pm 0.02$ | 0.7400 | $5 / 2^{+}$ |
| 3 | $3.12 \pm 0.02$ | 3.103 | $1 / 2^{-}$ |
| 4 | $4.21 \pm 0.02$ | 4.220 | $5 / 2^{-}$ |
| 5 | $4.65 \pm 0.02$ | 4.657 | $3 / 2^{-}$ |
| 6 | $5.87 \pm 0.02$ | 5.866 | $1 / 2^{-}$ |
| 7 | $6.85 \pm 0.02$ | 6.841 | $7 / 2^{-}$ |
| 8 | $7.36 \pm 0.02$ | 7.352 | $9 / 2^{-}$ |
| 9 | $8.47 \pm 0.02$ | 8.47 | $1 / 2^{+}, 3 / 2^{+}, 5 / 2^{+}$ |
|  |  |  | (from ref. 39) |
| 10 | $9.06 \pm 0.02$ | 9.00 | $1 / 2^{-}$(present work) |
| 11 | $13.7 \pm 0.1$ |  |  |
| 14 C states |  |  | $0^{+}$ |
| 1 | $0.00 \pm 0.02$ | 0 | $1^{-}$ |
| 2 | $6.10 \pm 0.02$ | 6.0938 | $3^{-}$ |
| 3 | $6.71 \pm 0.02$ | 6.7282 | $2^{+}$ |
| 4 | $7.00 \pm 0.02$ | 7.0120 | $2^{-}$ |
| 5 | $7.36 \pm 0.02$ | 7.314 | $2^{+}$ |
| 6 | $8.33 \pm 0.02$ | 8.3179 | $0^{+}$ |
| 7 | $9.81 \pm 0.02$ | 9.7460 | $2^{+}, 3^{-}$ |
| 8 | $10.43 \pm 0.02$ | $10.425,10.498$ | $4^{+}$ |
| 9 | $10.73 \pm 0.02$ | 10.736 | $3^{-}$ |
| 10 | $12.88 \pm 0.02$ | 12.963 |  |
| 11 | $13.96 \pm 0.02$ | 14.05 | $6^{+}$(from ref. 40) |
| 12 | $16.42 \pm 0.02$ | 16.43 | $6^{-}$(from ref. 40) |
| 13 | $16.74 \pm 0.02$ | 16.715 | $0^{+}$(present work) |
| 14 | $16.9 \pm 0.1$ |  |  |

*Values from ref. 38, except those explicitly indicated.




Supplementary Figure 7 - Comparison with calculations Discretized continuum scheme calculations for the $L=0$ case (red line) and experimental cross section angular distribution for the ${ }^{14} \mathrm{C}$ resonance at $16.9 \pm 0.1 \mathrm{MeV}$. No scaling factors are used.

## CRC and DWBA calculations



## Extreme cluster model

- Relative motion of the $2 n$ frozen and separated by the c.m.
- Only the term with the $2 n$ coupled to $S=0$ participates to the transfer
- S.A. $=1$ for all configurations


## I ndependent coordinate model

- The transfer is described taking into account spectroscopic information obtained by shell model calculations

| Coupling scheme | *) ${ }^{\text {誉高 }}$ |  |
| :---: | :---: | :---: |
|  |  |  |

## Sequential transfer (DWBA)

- Introducing the ${ }^{17} \mathrm{O}+{ }^{14} \mathrm{C}$ intermediate partition

Coupling scheme


## CRC and DWBA calculations



Extreme cluster model overestimate the cross section (S.A. = 1) I ndependent coordinate model describes quite well the cross section Sequential transfer (DWBA) underestimate the cross section

## Microscopic cluster calculations

Wave functions for two particles in an harmonic oscillator common potential ( $j-j$ coupling)
wave functions in terms of the relative and centre of
mass coordinates of the two particles ( $\angle S$ coupling)
( $\mathrm{n}, \mathrm{I}$ ) internal cluster state
$(\mathrm{N}, \mathrm{L})$ cluster motion relative

$$
\begin{aligned}
& \text { to the core } \\
& \hat{a}=\sqrt{2 a+1} \\
& \text { Moshinsky coefficients }
\end{aligned}
$$

Two neutron amplitudes - zbm interaction

| Initial state | $\mathrm{j}_{1} \mathrm{j}_{2}$ | $\mathrm{J}_{12}$ | Final state | Spectr. Amp. | n | 1 | N | L | $\wedge$ | S | Spec. Amp. (c.m.) |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| ${ }^{13} \mathrm{C}_{\mathrm{g} .5}\left(1 / 2^{-}\right)$ | $\left(p_{1 / 2} \mathrm{~s}_{1 / 2}\right)$ | 0 | ${ }^{15} \mathrm{C}_{\text {g. }}\left(1 / 2^{+}\right)$ | -0.641 | 1 | 0 | 2 | 1 | 1 | 1 | -0.292 |
|  |  |  |  |  | 1 | 1 | 1 | 2 | 1 | 1 | 0.338 |
|  |  |  |  |  | 1 | 1 | 2 | 0 | 1 | 1 | -0.075 |
|  | $\left(p_{1 / 2} s_{1 / 2}\right)$ | 1 |  | -1.110 | 1 | 0 | 2 | 1 | 1 | 0 | 0.292 |
|  |  |  |  |  | 1 | 1 | 1 | 2 | 1 | 0 | -0.338 |
|  |  |  |  |  | 1 | 1 | 2 | 0 | 1 | 0 | 0.075 |
|  |  |  |  |  | 1 | 0 | 2 | 1 | 1 | 1 | -0.413 |
|  |  |  |  |  | 1 | 1 | 1 | 2 | 1 | 1 | 0.477 |
|  |  |  |  |  | 1 | 1 | 2 | 0 | 1 | 1 | -0.107 |

## Microscopic cluster calculations



## Extreme cluster model

## Microscopic cluster 1s

- Taking into account configurations with $\mathrm{n}=1 \mathrm{I}=0$

Microscopic cluster 1s + 1p

- Taking into account configuration with $\mathrm{n}=1 \mathrm{I}=0,1$

Coupling scheme

> Transitions to ground and 3.103 MeV states reproduced rather well with $1 s+1 p$ waves
> Transition to 0.74 MeV state probably needs more configurations

## Microscopic cluster calculations



## New works published in 2016-2017

Test of model space for the $<{ }^{18} \mathrm{O} \mid{ }^{16} \mathrm{O}>$ projectile overlaps

## Study of the ${ }^{18} \mathrm{O}\left({ }^{16} \mathrm{O},{ }^{18} \mathrm{O}\right)^{16} \mathrm{O}$ reaction at 84 MeV incident energy zbm vs psdmod interactions

| Model space | valence orbitals |
| :--- | :--- |
| zbm $\left({ }^{12} \mathrm{C}\right.$-core $)$ | $1 \mathrm{p}_{1 / 2}, 1 \mathrm{~d}_{5 / 2}, 2 \mathrm{~s}_{1 / 2}$ |
| psdmod $\left({ }^{4} \mathrm{He}\right.$ core $)$ | $1 \mathrm{p}_{3 / 2}, 1 \mathrm{p}_{1 / 2}, 1 \mathrm{~d}_{5 / 2}, 2 \mathrm{~s}_{1 / 2}, 1 \mathrm{~d}_{3 / 2}$ |

## Experimental results



M. J. Ermamatov et al., PRC 94 (2016) 024610

## Results of theoretical calculations


g.s. only $S=0(A)$
$2+S=0(A)$ or (P)
Extreme cluster model works
M. J. Ermamatov et al., PRC 94 (2016) 024610


For the lower states of projectile overlaps the zbm model- space is enough. The study of the higher excited states is in progress

## New works in progress (some results)

## Study of the ${ }^{18} \mathrm{O}\left({ }^{64} \mathrm{Ni},{ }^{66} \mathrm{Ni}\right){ }^{16} \mathrm{O}$ reaction at 84 MeV incident energy

| Model space | valence orbitals |
| :--- | :--- |
| protons | $1 \mathrm{p}_{1 / 2}, 1 \mathrm{~d}_{5 / 2}, 2 \mathrm{~s}_{1 / 2}$ |
| neutrons | $1 \mathrm{p}_{3 / 2}, 1 \mathrm{p}_{1 / 2}, 1 \mathrm{~d}_{5 / 2}, 2 \mathrm{~s}_{1 / 2}, 1 \mathrm{~d}_{3 / 2,1}, 1 \mathrm{~g}_{7 / 2}$ |



## Results of theoretical calculations





Microscopic results: g.s.: IC results are better, specially in the bell-shaped region. Same order: one and two step. $2^{+}$: Long-range correl. (coll.) dominates over the short-range (pairing)

## Results of theoretical calculations




Microscopic results:
g.s.: IC results are better, specially in the bell-shaped region. $2^{+}$: Long-range correl. (coll.) dominates over the short-range (pairing)


For details, see R. Magana poster
B. Paes et al PRC 96.044612 (2017) -yesterday

# Results of theoretical calculations 

| Nucleus B(E2); $0^{+}!$ | $2^{+}\left(\mathrm{e}^{2} \mathrm{~b}^{2}\right)$ |
| :---: | :---: | :---: |
| ${ }^{14} \mathrm{C}$ | 0.0018 |
| ${ }^{18} \mathrm{O}$ | 0.0045 |
| ${ }^{28} \mathrm{Mg}$ | 0.035 |
| ${ }^{66} \mathrm{Ni}$ | 0.060 |
| ${ }^{76} \mathrm{Ge}$ | 0.270 |

Small for ${ }^{14} \mathrm{C}{ }^{18} \mathrm{C}$
Big for ${ }^{28} \mathrm{Mg}^{66} \mathrm{Ni}{ }^{76} \mathrm{Ge}$

## New works in progress (some results)

## Study of the ${ }^{18} \mathrm{O}\left({ }^{28} \mathrm{Si},{ }^{30} \mathrm{Si}\right){ }^{16} \mathrm{O}$ reaction at 84 MeV incident energy

Model space ( ${ }^{4} \mathrm{He}$ core)

## valence orbitals (similar to Ni)

$1 p_{3 / 2}, 1 p_{1 / 2}, 1 d_{5 / 2}, 2 s_{1 / 2}, 1 d_{3 / 2}$

$$
\begin{equation*}
1 p_{3 / 2}, 1 p_{1 / 2}, 1 d_{5 / 2}, 2 s_{1 / 2}, 1 d_{3 / 2} \tag{b}
\end{equation*}
$$

$\mathbf{E}(\mathrm{MeV}) ; \mathbf{J}^{\pi}$



## Results of theoretical calculations



Cluster model is not good for ${ }^{28,30} \mathrm{Si}$

Microscopic results: g.s.: Two-step DWBA results are better. Same order: one and two step. $2^{+}$: Long-range correl. (coll.) dominates over the short-range (pairing) Si* the same results as the $2^{+}$state

## Results of theoretical calculations



## Does our theoretical calculations describe other observables?

- Elastic scattering
- Inelastic scattering


## Conclusions and outlooks

$>{ }^{12,13} \mathrm{C}\left({ }^{18} \mathrm{O},{ }^{16} \mathrm{O}\right){ }^{15} \mathrm{C},{ }^{16} \mathrm{O}\left({ }^{18} \mathrm{O},{ }^{16} \mathrm{O}\right){ }^{18} \mathrm{O},{ }^{64} \mathrm{Ni}\left({ }^{18} \mathrm{O},{ }^{16} \mathrm{O}\right){ }^{66} \mathrm{Ni}$, ${ }^{28} \mathrm{Si}\left({ }^{18} \mathrm{O},{ }^{16} \mathrm{O}\right){ }^{30} \mathrm{Si}$, at 84 MeV incident energy
> Four models were used to calculate the cross section:
$\checkmark$ Extreme cluster
$\checkmark$ Independent coordinate
$\checkmark$ DWBA
$\checkmark$ Microscopic cluster (only for ${ }^{13}$ C)
$>$ no need for any "unhappiness" factor to reproduce the absolute cross sections
$>$ In ${ }^{13} \mathrm{C}$ importance of a two-neutron correlation in the nuclear wave function, the extra neutron does not destroy the correlations observed in the ${ }^{14} \mathrm{C}$ case
> Dominance of the 1 s and 1 p waves in the two-neutron cluster internal wave function
$>$ Adequacy of zbm interaction for low-lying overlaps of the projectile were established for the projectile.
$>$ Dominance of long-range correlations for the excited $2^{+}$state of ${ }^{66} \mathrm{NI}$ over the short-range pairing correlations. The opposite for the g.s. $>$ Dominance of long-range correlations in all states of ${ }^{30} \mathrm{Si}$.

## Conclusions and outlooks

## Outlooks:

$>$ Include other waves in the microscopic cluster calculations
$>$ Enlarge the model space for higher energy transitions $\left(\mathrm{d}_{3 / 2}\right)$
$>$ Describe high excited states of the projectile.
$>$ Include the deformed target ( ${ }^{28} \mathrm{Si}$ ) to study the mixing of collective and single particle configurations.
$>$ Study the $2 p$ and $n p$ transfers to study the pairing correlations in collaboration with the structure group of Genova of Prof. Santopinto.


REACTIONS SCHEME CONCERNING THE ${ }^{116} \mathrm{Cd}\left({ }^{(20} \mathrm{Ne},{ }^{20} \mathrm{O}\right){ }^{116} \mathrm{Sn}$


## Working group

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## Thank you

