

Study of single-nucleon transfer reactions in the ¹⁸O+⁴⁸Ti collision at 275 MeV



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Experimental setup: The MAGNEX facility @ INFN-LNS



Motivation-Introduction



NUclear Matrix Elements for Neutrinoless double beta decay $(\mathbf{0}\mathbf{v}\mathbf{\beta}\mathbf{\beta})$

✓ **NUMEN** [1] is an innovative project aiming at accessing information on the Nuclear Matrix Elements (NMEs) of the Ovßß decay through the study of heavy-ion induced Double Charge Exchange (DCE) reactions.

 \checkmark DCE reactions and $0\nu\beta\beta$ decay are mediated by different interactions BUT they exhibit some interesting similarities. Among them, both processes probe the same initial and final nuclear wave functions and the operators connecting them have a similar spin – isospin mathematical structure [2,3].

✓ Transfer reactions: Competing processes leading to the same final states as DCE reactions. Their study can provide information on the degree of competition between successive nucleon transfer and charge exchange reactions, the latter being of particular interest in the context of single and double beta decay studies [4,5].

✓ The ¹⁸O+⁴⁸Ti reaction was studied at the energy of 275 MeV for the first time under the NUMEN and NURE [6] experimental campaigns with the aim of investigating the complete net of reaction channels potentially involved in the ⁴⁸Ti \rightarrow ⁴⁸Ca DCE transition.



Optical characteristics of MAGNEX	
Max. magnetic rigidity (Tm)	1.8
Solid angle (msr)	50
Momentum acceptance	-14%, +10%
Energy resolution ($\Delta E/E$)	~1/1000
Angular resolution	~ 0.3°
Mass resolution	~ 1/160

•Target: TiO₂ evaporated on a thin ²⁷Al foil. Background estimation: 2 additional runs with a WO_3 and a ²⁷Al target. •The reaction ejectiles were detected by the MAGNEX Focal Plane Detector (FPD).







reaction.

DWBA framework

The main ingredients of the DWBA calculation are: > The distorted waves at the entrance and exit channels —> Calculated adopting the São Paulo Potential (SPP) [8].

results of a DWBA and a CCBA calculation. **CCBA framework**

A CCBA calculation is performed in the same way as the DWBA one considering also the following additions:

Coulomb deformations: Introduced in terms

of of the reduced transition probabilities $B(E\lambda)$. >The overlap functions: Single particle solutions of a Woods-Saxon potential weighted by the >Nuclear couplings: The nuclear coupling potentials were treated in the first-order corresponding spectroscopic amplitudes provided approximation as described in Ref. [9]. by many body shell-model calculations.



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Concluding Remarks and perspectives

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✓ The theoretical cross-sections are found to be in very good agreement with the experimental data suggesting the validity of the adopted optical potentials and the shell-model description of the involved nuclear states within the adopted model space. ✓ The same technique will be also tested through the data set of the two-proton transfer channel.

✓ The results of the present analysis in the ¹⁸O+⁴⁸Ti reaction will clarify the degree of competition between the direct SCE mechanism (¹⁸O,¹⁸F) and the sequential nucleon transfer.

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