

# MICROSCOPIC CALCULATION OF THE TOTAL PHOTOABSORPTION CROSS SECTION OF $A = 6$ NUCLEI

In collaboration with:

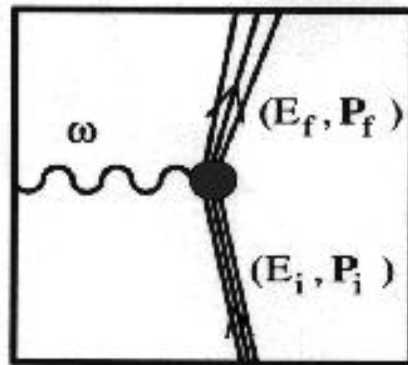
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## OUTLINE OF THE TALK:

- \* Inclusive processes: LIT method
- \* EIHH approach
- \* Lanczos algorithm
- \* Results
- \* Conclusions and future perspectives

# INCLUSIVE PHOTODISINTEGRATION OF A NUCLEUS



Cross section  $\sigma(\omega) = \frac{4\pi^2\alpha}{2J_i+1} \omega F(\omega)$

$$F(\omega) = \int d\Psi_f |\langle \Psi_f | \mathbf{E}1 | \Psi_0 \rangle|^2 \delta(E_f - E_0 - \omega)$$

A new AB INITIO calculation for the 6-body problem is allowed via the use of:

LIT Lorentz Integral Transform Method

*P.L. B 338 (1994) V.D.Efros, W.Leidemann, G.Orlandini*

EIHH Effective Interaction with HH Formalism

*P.R. C 61 (2000) N.Barnea, W.Leidemann, G.Orlandini*

## LORENTZ INTEGRAL TRANSFORM METHOD FOR INCLUSIVE PROCESSES

It enables to consider the full final state interaction of the system exactly.

1. Apply an integral transform on  $F(\omega)$  with a lorentzian shape kernel  $K(\omega)$

$$\begin{aligned} L(\sigma) &= \int d\omega K(\omega) F(\omega) \\ &= \int d\omega \frac{F(\omega)}{(\omega - \sigma_R)^2 + \sigma_I^2} = \langle \tilde{\Psi} | \tilde{\Psi} \rangle \end{aligned}$$

$$\sigma = -\sigma_R + i\sigma_I$$

2. The problem is reduced to the solution of a Schrödinger-like equation

$$(H - E_0 - \sigma_R + i\sigma_I) |\tilde{\Psi}\rangle = \hat{O} |\Psi_0\rangle$$

which can be found with bound state techniques.

3. Inversion of the transform

$$L(\sigma) \longrightarrow F(\omega) \longrightarrow \sigma(\omega)$$

## EFFECTIVE INTERACTION THEORY

Problem: Solve  $\hat{H}|\Psi\rangle = E|\Psi\rangle$  for an A-body system with  $\hat{H} = \hat{T} + \hat{V}$  expanding  $|\Psi\rangle$  on a finite set  $P$  of basis functions  $|\Phi_i\rangle$ .

Purpose: Define an effective interaction that, acting in a truncated model space reproduces the true energy spectrum.

### LEE-SUZUKI APPROACH

1. Similarity transformation  $X = e^\omega$

$$\tilde{H} = X^{-1}\hat{H}X = \hat{P}\tilde{H}\hat{P} + \hat{P}\tilde{H}\hat{Q} + \hat{Q}\tilde{H}\hat{P} + \hat{Q}\tilde{H}\hat{Q}$$

2. Decoupling condition

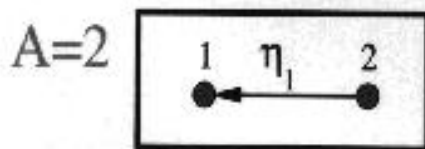
$$\hat{Q}\tilde{H}\hat{P} = \hat{Q}(X^{-1}\hat{H}X)\hat{P} = 0$$

3. Effective interaction

$$\hat{H}_{eff} = \hat{P}\tilde{H}\hat{P} = \hat{P}(X^{-1}\hat{H}X)\hat{P}$$

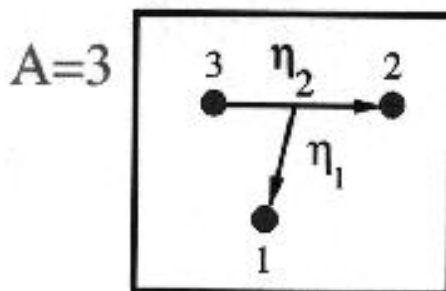
## HYPERSPHERICAL FORMALISM

It is used for the expansion of an  $A$  body internal wave function.



$$\eta_1 \rightarrow \rho_1, \theta_1, \phi_1$$

$$\Delta_{(1)} = \Delta_{\rho_1} - \frac{1}{\rho_1^2} \hat{L}_1^2$$



$$\eta_1 = \rho_1 = \rho_2 \cos \varphi_2$$

$$\eta_2 = \rho_2 \sin \varphi_2$$

$$\eta_1, \eta_2 \rightarrow \rho_2, \underbrace{\varphi_2, \theta_1, \phi_1, \theta_2, \phi_2}_{\hat{\Omega}_2}$$

$$\Delta_{(2)} = \Delta_{\rho_2} - \frac{1}{\rho_2^2} \hat{K}_2^2$$

In General

$$\eta_1, \eta_2, \dots, \eta_{A-1} \rightarrow \rho, \hat{\Omega}$$

$$H = \Delta_{\rho}^2 - \frac{\hat{K}^2}{\rho^2} + \sum_{i < j}^A \hat{V}_{ij}(\rho, \hat{\Omega})$$

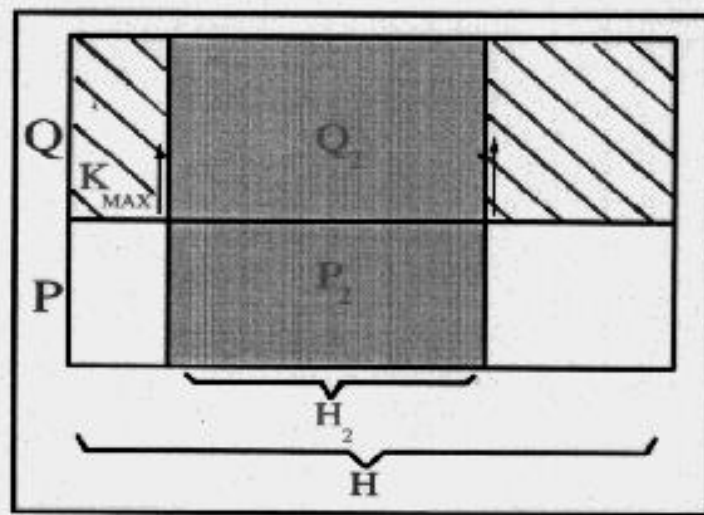
The HH are eigenfunctions of the hyperspherical angular momentum  $\hat{K}^2$ .

## EIH METHOD

$$H = \Delta_\rho^2 - \frac{\hat{K}^2}{\rho^2} + V_{12}(\rho, \hat{\Omega}) + \dots + V_{AA-1}(\rho, \hat{\Omega})$$

$$H^{(2)}(\rho) = \Delta_\rho^2 - \frac{\hat{K}^2}{\rho^2} + V_{AA-1}(\rho, \hat{\Omega})$$

$$H^{(2)}(\rho) \xrightarrow{L.S.} H_{eff}^{(2)}(\rho) \rightarrow V_{eff}^{(2)}(\rho) = H_{eff}^{(2)}(\rho) - \frac{\hat{K}^2}{\rho^2}$$



In the limit of  $K_{MAX} \rightarrow \infty$  one has  $V_{eff}^{(2)} \rightarrow V$   
 $\Rightarrow$  the solution is exact when convergence is reached

## ADVANTAGES OF EIIH

- $\rho$  is a collective coordinate
- $\hat{K}^2$  depends on the angular momenta of the residual  $A - 2$  system

⇒ medium correction ⇒ faster convergence.

## POTENTIAL

In the two-body hyperspherical Hamiltonian

$$V_{AA-1}(\rho, \hat{\Omega})$$

we use a semirealistic potential:

- Minnesota
- Malfliet Tjon

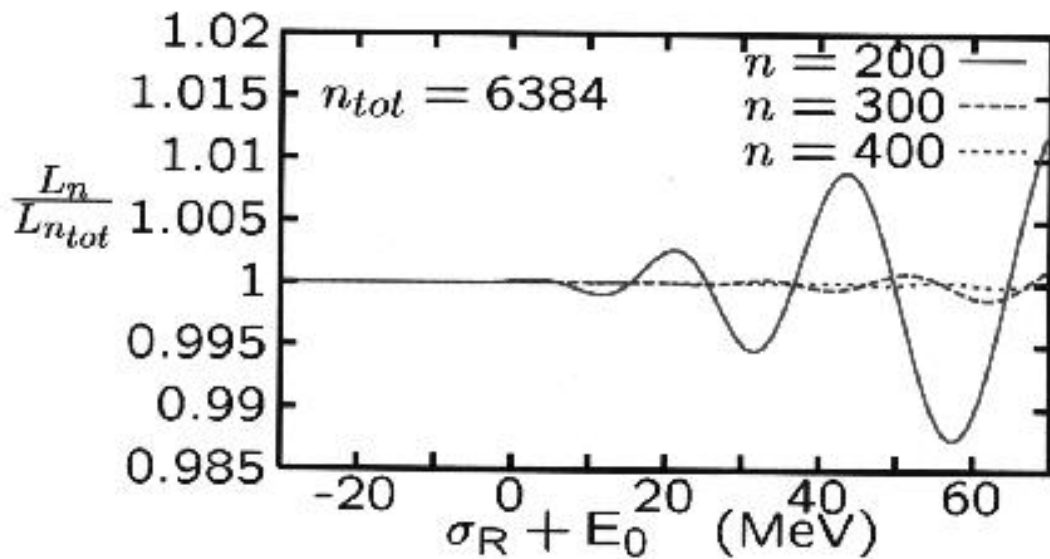
## CALCULATION OF THE LIT WITH THE LANCZOS ALGORITHM

Recursive definition of the Lanczos vectors

$$b_{n+1} |\phi_{n+1}\rangle = \widehat{H} |\phi_n\rangle - a_n |\phi_n\rangle - b_n |\phi_{n-1}\rangle$$

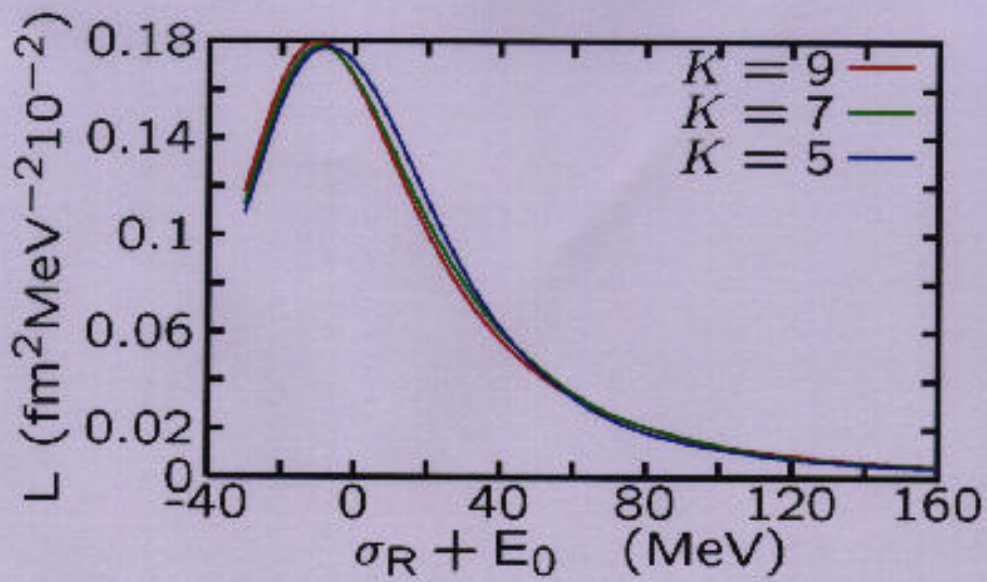
The Lorentz integral transform can be written as a function of the Lanczos' coefficients

$$L(\sigma) = \frac{1}{\sigma_I} \text{Im} \frac{\langle \psi_0 | \widehat{D}_z^\dagger \widehat{D}_z | \psi_0 \rangle}{(z - a_0) - \frac{b_1^2}{(z - a_1) - \frac{b_2^2}{(z - a_2) - b_3^2 \dots}}}$$

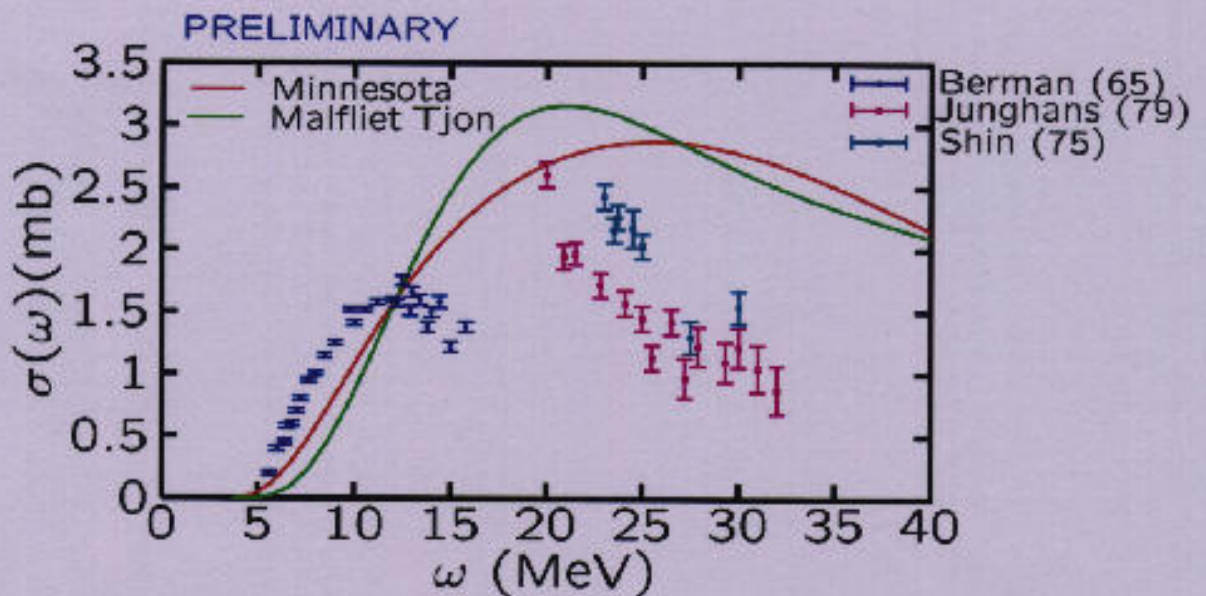




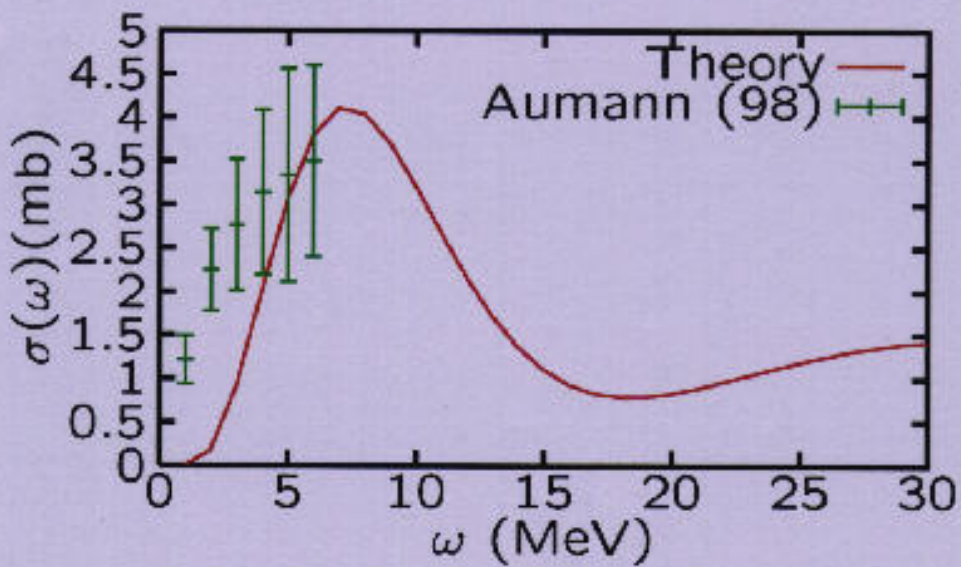
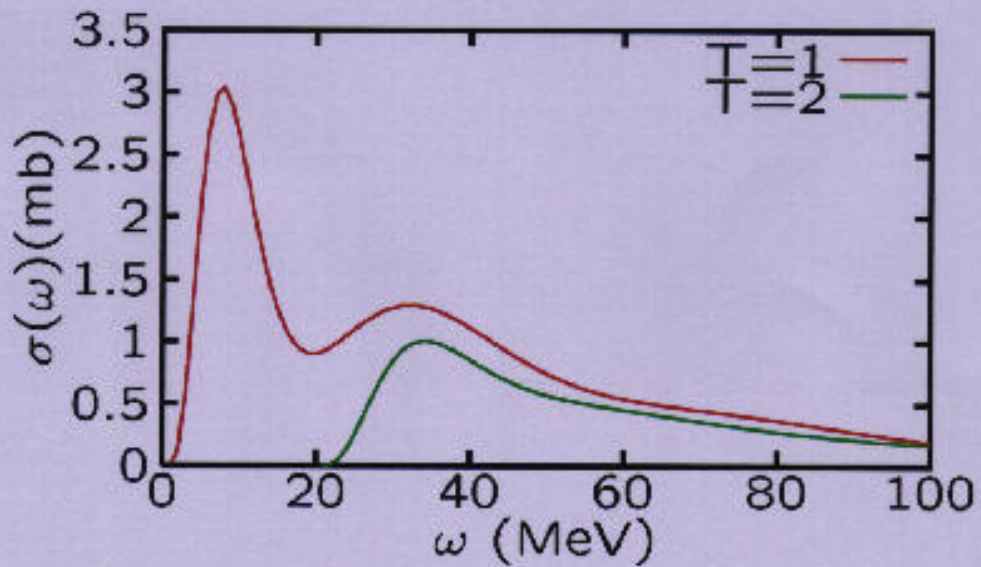
## LIT RESULTS FOR ${}^6\text{Li}$



## CROSS SECTION FOR ${}^6\text{Li}$



# CROSS SECTION FOR ${}^6\text{He}$ Minnesota



## CONCLUSIONS

- We are presenting the first microscopic calculation of inelastic reactions of  $A = 6$  nuclei with the full final state interaction
- The LIT and EIHH are two very powerful techniques for an AB INITIO calculation with  $A > 4$

## FUTURE PERSPECTIVES

- The convergence has to be improved (parallelization of the code)
- Use of more realistic potentials
- Applications with virtual photons
- Study of even larger systems



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