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Introduction-Outline

Outline

- Basic reaction theory
 - Potentials
 - Time evolution
- Single Channel
- Coupled Channels

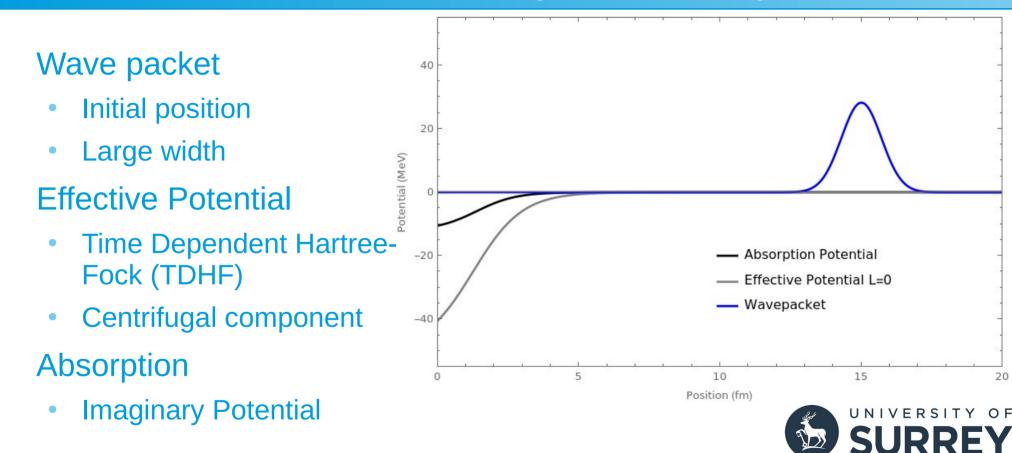


University of Warwick/Mark Garlick, https://www.eso.org/public/images/eso1733s/





Basic Problem-Wave packet In Space



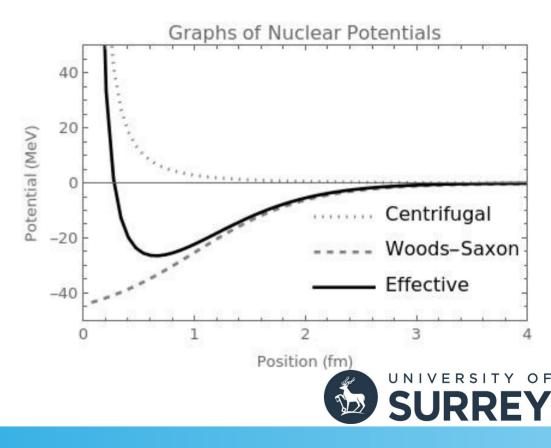
Basic Problem-Potential

Internal Potential

$$V_{WS} = -\frac{U_{WS}}{1 + \exp\frac{r - r_{ws}}{a_{WS}}}$$

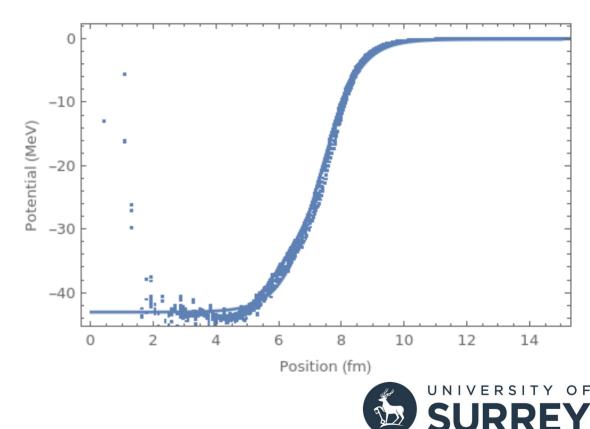
Centrifugal Potential

$$V_C = \frac{\hbar^2 l(l+1)}{2\mu r^2}$$



Basic Problem-TDHF

- Static TDHF
 - Many internal wavepackets
 - Allowing them to evolve
 - Mapping x,y,z into radial.
- Least Squares
 - Effective Woods Saxon
 - Excludes internal potential





Basic Problem-Time Evolution

- Chebyshev Polynomials
 - Time Evolution

- Extending into a set of polynomials
- Recurrence Relation

$$\psi(r,t) = e^{-i\frac{Ht}{\hbar}}\psi(r,0).$$

$$\psi_0(t) = \psi(t)$$

$$\psi_1(t) = H\psi_0(t)$$

$$\exp^{-iH\Delta t} = J_0(\tau)\sum_{n=1}^{\infty} J_n(\tau)T_n(H)$$

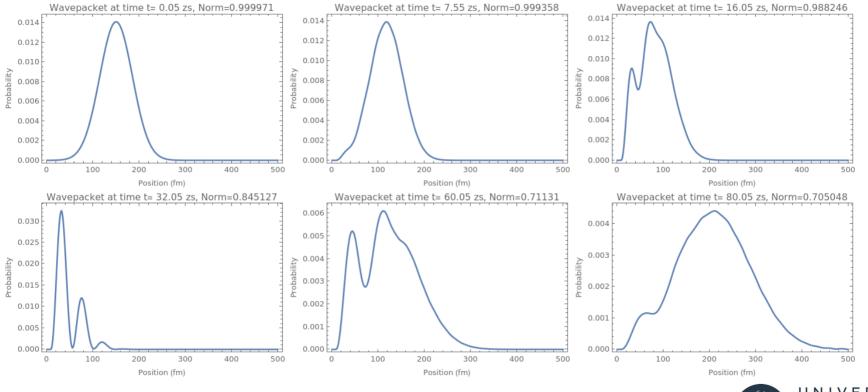
$$\psi_n(t) = 2H\psi_{n-1}(t) - \psi_{n-2}(t)$$



Basic Problem-Absorption

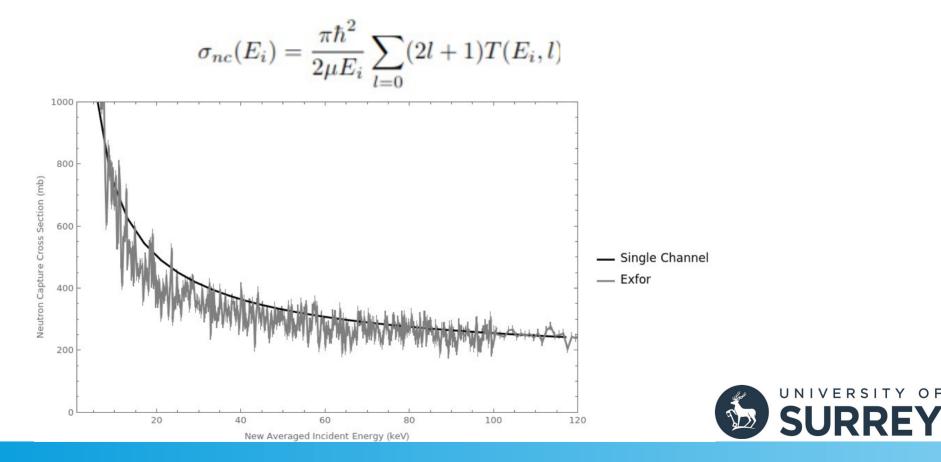
- Absorption Potential
- Modified Chebyshev Polynomials
- Woods Saxon Shaped Potential

Results-Wavepacket Propagation



SURREY

Results-Neutron Capture Cross Sections



Coupled Channels-Theory

Shifted Woods-Saxon

$$V_{CC} = -\frac{U_{WS}}{1 + \exp((r - r_{WS} - \hat{r}_{cc})/a_{WS})}$$

- Generate and Diagonalise the Coupling Matrix • $\langle I_n | \hat{r}_{cc} | I_{n'} \rangle = r_{coup} \beta_2 F(2, I_n, I'_n)$
- Generate a full coupled channels Hamiltonian

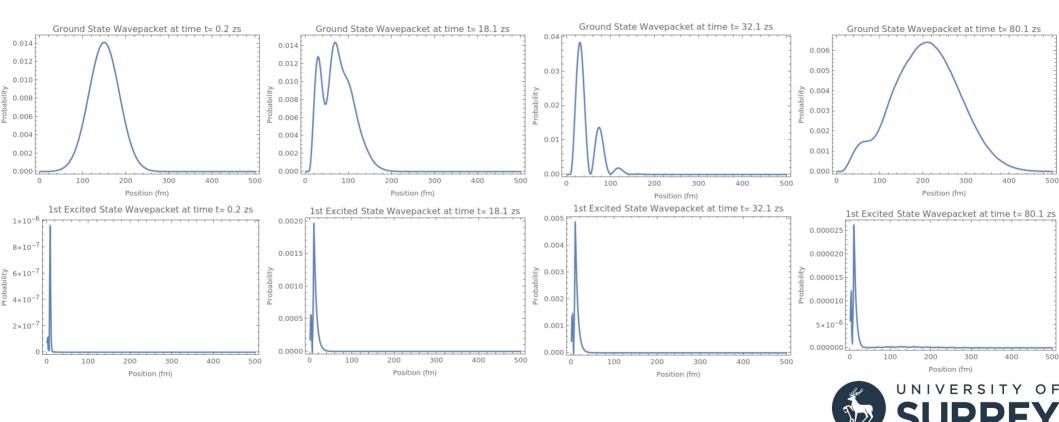
$$H = \begin{pmatrix} 0 - 0 & 0 - 2 \\ 2 - 0 & 2 - 2 \end{pmatrix} \begin{pmatrix} |\psi_1\rangle \\ |\psi_2\rangle \end{pmatrix}$$

Diagonal Entries
 Off-Diagonal Entries

$$(In - In) = K + V_C + V_{WS} + V_{cc} + \epsilon_n I. \qquad (I_n - I_{n'}) = V_{nn'}$$

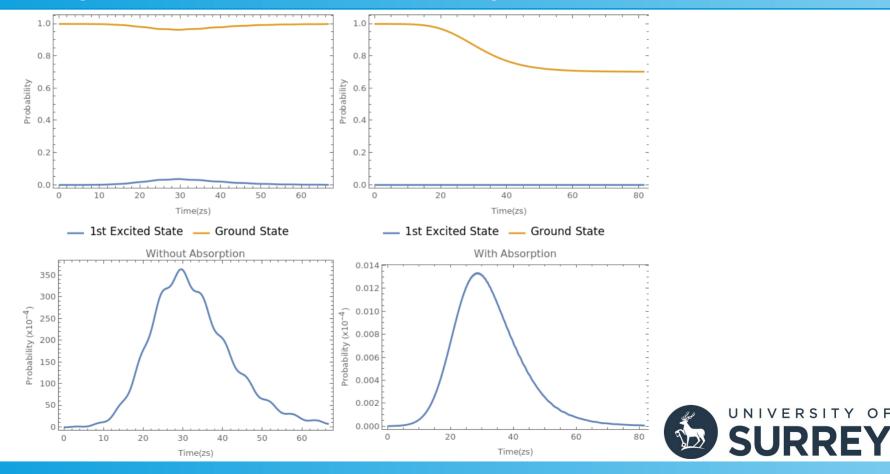


Coupled Channels-Wave packet Propagation

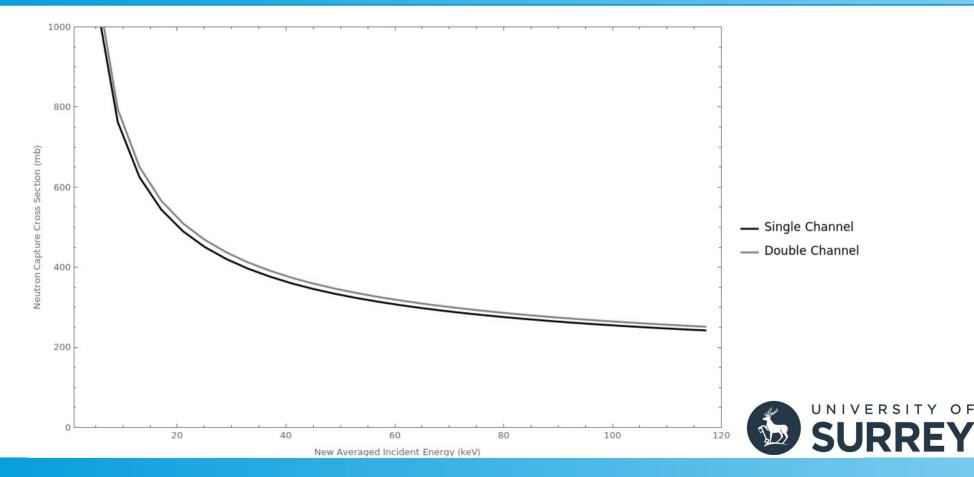




Coupled Channels-Absorption effects

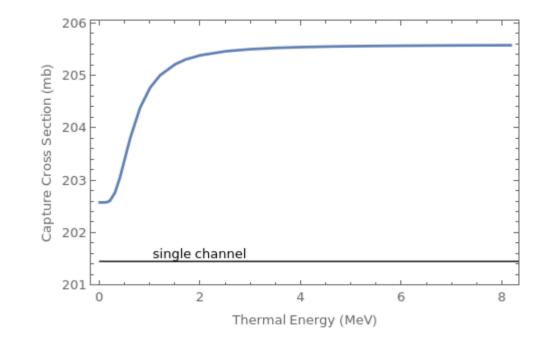


Coupled Channels-Cross Section Comparisons

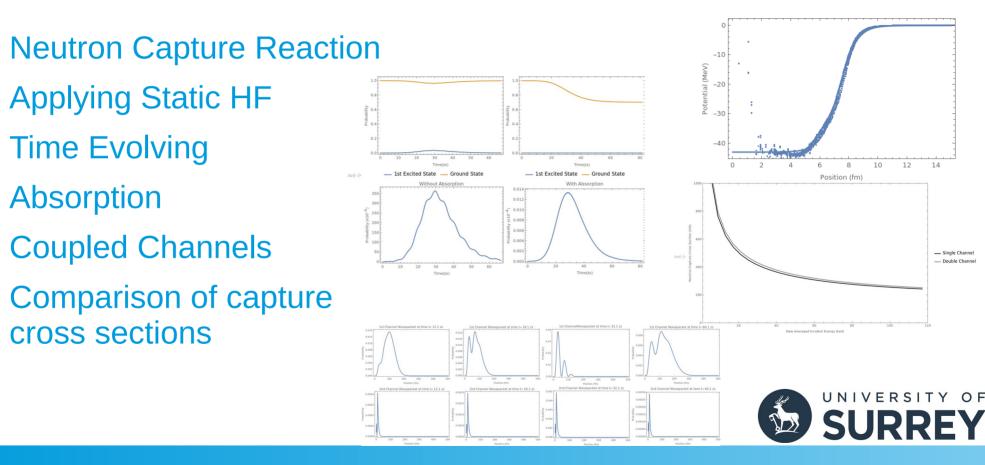


Future Steps

- Thermal Effects
- Reaction Rates
- Comparison to FRESCO,CCFULL, etc.
- 187-Os, 186-Os, etc.

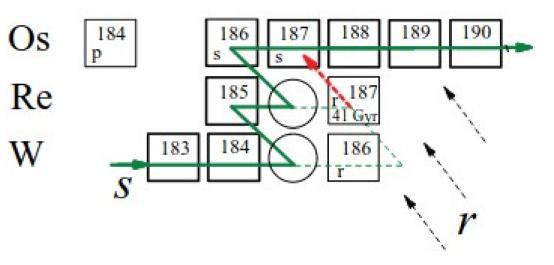


Recap



Introduction

- The reaction
 - Os-Re clock
- Progression
 - Basic Problem
 - Single Channel
 - Coupled Channels



Mosconi, M., et al. "Neutron physics of the Re/Os clock. I. Measurement of the (n,gamma) cross sections of Os 186,187,188 at the CERN n_TOF facility." *Physical Review C-Nuclear Physics* 82.1 (2010):015802





Thermal Effects-Theory I

- Liouville Equation
 - In the case with a pure state
 - Our initial non-thermalised state

- Lindblad Equation
 - Gives environmental effects
 - Thermalisation of the system

$$i\hbar \frac{\delta \rho(x, x', t)}{\delta t} = [\hat{H}, \hat{\rho}]_{x, x', t}$$
$$\hat{\rho} = |\psi^*(x, t)\rangle \langle \psi(x', t)|$$

$$i\hbar\frac{\delta\rho(x,x',t)}{\delta t} = [\hat{H},\hat{\rho}]_{x,x',t} + \hat{L}[\ldots].$$



Coupled Channels-Theory

Shifted Woods-Saxon

$$V_{CC} = -\frac{U_{WS}}{1 + \exp((r - r_{WS} - \hat{r}_{cc})/a_{WS})}$$

Generate the Coupling Matrix

$$\langle I_n | \hat{r}_{cc} | I_{n'} \rangle = r_{coup} \beta_2 F(2, I_n, I'_n)$$

$$V_{N,nn'} = \langle I_n | V_N(r, \hat{r}_{CC}) | I_{n'} \rangle - V_N(r, 0) \delta_{nn'}$$
$$= \sum_{\alpha} \langle I_n | \alpha \rangle V_N(r, r_{cc,\alpha}) - V_N(r, 0) \delta_{nn'}$$





Form Factor

- Form factor
- Generates using a Jsymbol

$$F(I, I_n, I_{n'}) = \sqrt{\frac{(2I+1)(2I_n+1)(2I_{n'}+1)}{4\pi}} \begin{pmatrix} I_n & I & I_{n'} \\ 0 & 0 & 0 \end{pmatrix}^2$$

• Similar to a Glebsch-Gordon coefficient.

