Neutron Transfer Reactions for Deformed Nuclei

Sturmian Basis Approach

to the $P(J^{\pi}, E)$ distribution in ¹⁵⁶Gd* from ¹⁵⁷Gd(³He,⁴He)¹⁵⁶Gd*







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Outline of the Talk

- Introduction and Motivation
- Surrogate Reactions
- Strongly Deformed Nuclei
 - Single particle states
 - ¹⁵⁶Gd* spectrum (Bohr-Mottelson rotor plus particle model)
 - Sturmian basis approach for reaction cross sections
 - J^{π} distribution estimates for ${}^{157}Gd+{}^{3}He \rightarrow {}^{4}He+{}^{156}Gd*$
- Outlook

Origin of the Elements

s-process branch point nuclei are unstable



•All elements starting with carbon (C) are products of stellar evolution.

•Most light to intermediate nuclei are products of various nuclear burning-cycles.

- •Heavier nuclei along the valley of stability are produced via neutron capture reactions.
- •Elements away from the valley of stability are produced in explosive events.

The Surrogate Method

deducing reaction cross sections on unstable nuclei from reactions on stable nuclei



If two reactions proceed via formation of the same intermediate equilibrated compound system B^{*} then the cross section for the desired reaction $a+A \rightarrow B^* \rightarrow C+c$, which might involve an *unstable* target A, may be deduced using theoretical modeling and experimental data for the *surrogate* reaction $d+D \rightarrow b+B^* \rightarrow C+c$ for a *stable target* D.

Reactions with Deformed Nuclei Computational Strategy and Methods

• Main assumptions:

- Direct particle transfer to/from a *deformed single particle state* (sps).
- Axially deformed Woods-Saxon potential defines the deformed sps.
- Reaction cross section can be computed with the available reaction codes by using superposition of *Sturmian spherical single particle states*.

• Computational steps where coding or scripting is completed:

- WSBETA code to obtain *energies and deformed states* (Ψ_E) for a given *deformed potential*, *big deformations careful numerical treatment*.
- **DWUCK**^{*} code generates the set of *spherical WS potentials* and *non-orthogonal Sturmian basis wave functions* ϕ_{Enli} for the desired particle binding energy *E*.
- **WSBETA** code uses DWUCK's *spherical WS potentials* to compute the *Sturmian basis wave functions* in the same basis where the *deformed state* Ψ_E is given.
- **CalculateCij** code computes the overlaps $C_{nlj} = \langle \phi_{Enlj} | \Psi_E \rangle$ using WSBETA wave functions.
- **CHUCK**^{*} CC-code uses $\Psi_{E}(r) = \sum C_{nli} \Phi_{Enli}$ to compute the reaction cross sections.

Auxiliary codes

- AddStates code uses C_{nlj} and $\Phi_{Enlj}(r)$ to verify that the Sturmian basis is sufficiently big and to construct $\Psi_{E}(r)$ as superposition of Sturmian basis states.
- Various python scripts and Mathematica notebooks to prepare inputs, run the codes, collect, organize and visualize the results....
- * Essential help from P. D. Kunz, Dept. of Physics & Astrophysics, U. of Colorado, Boulder, CO

Need for Careful Numerical Treatment

$$V(r) = \frac{V_0}{1 + \exp\left(\frac{r-R}{a}\right)}$$

 $R(\theta, \phi) = R_0 \left(1 + \beta_2 Y_{20} \left(\theta, \phi \right) + \beta_4 Y_{40} \left(\theta, \phi \right) \right)$

N=4 HO Neutron Orbits

11

9

9

• 3

- Q

1h_{9/2} & 2f_{7/2} Valence Neutron Orbits

Solid lines - exact numerical results **Dashed lines - linear approximation**



Neutron Single Particle States

n	Energy	2Mj	Parity	n	Energy	2Mj	Parity	n	Energy	2Mj	Parity
1	-39.706	1	1	20	-18.202	5	1	40	-7.932	1	-1
2	-36.096	1	-1	21	-18.096	5	-1	41	-7.867	1	1
3	-34.408	3	-1	22	-17.785	3	-1	42	-7.607	3	-1
4	-33.782	1	-1	23	-16.681	1	-1	43	-7.404	3	1
5	-31.401	1	1	24	-16.586	1	1	44	-7.058	1	1
6	-30.202	3	1	25	-16.579	7	1	45	-6.842	11	-1
7	-29.193	1	1	26	-14.649	1	1	46	-6.603	3	1
8	-28.322	5	1	27	-14.645	3	1	47	-6.163	3	-1
9	-26.503	3	1	28	-14.435	9	1	48	-5.765	5	1
10	-25.983	1	-1	29	-13.723	1	-1	49	-5.367	5	-1
11	-25.804	1	1	30	-13.179	3	-1	50	-4.815	1	-1
12	-25.100	3	-1	31	-12.396	3	1	51	-4.583	7	1
13	-23.653	5	-1	32	-12.198	5	-1	52	-4.102	5	-1
14	-23.370	1	-1	33	-12.073	5	1	53	-3.076	9	1
15	-21.625	7	-1	34	-11.033	1	1	54	-2.665	7	-1
16	-21.116	3	-1	35	-10.829	7	-1	55	-2.512	1	1
17	-20.678	1	-1	36	-9.785	5	1	56	-2.483	1	-1
18	-20.056	1	1	37	-9.308	1	-1	57	-1.901	3	-1
19	-19.379	3	1	38	-9.072	9	-1	58	-1.772	7	-1
				39	-8.791	7	1	59	-1.224	11	1
								60	-1.103	3	1
	157	C	1					61	-0.627	1	1
	. • ·	J						62	-0.521	3	-1
								63	-0.171	1	-1
								64	0.120	1	-1

Unpaired valence neutron at level 47. A=157, N=93, Z=64 ¹⁵⁷Gd: $S_n = 6.36 \text{ MeV}$ E_n=0 ... 1 MeV n+155Gd= =¹⁵⁶Gd+E_{ext} ¹⁵⁷Gd-n= E_{ext}=8.5 ... 9.5 MeV ¹⁵⁶Gd: $S_n = 8.536 \text{ MeV}$ **BE**_n ~ -14.8987 MeV single particle levels relevant for the surrogate reaction

Excited states in ¹⁵⁶Gd model: neutron hole in the ¹⁵⁷Gd core

$$|^{156}Gd, \Omega = |K \pm \nu| > = \psi^{\dagger}_{\pm \nu}|^{157}Gd, K = 3/2^{-} > 0$$

$$E(J^{\pi}; \Omega = |K \pm \nu|) = \epsilon_0 - \epsilon_{\nu} + \frac{\hbar^2}{2\mathcal{I}}(J(J+1) + \delta_{\pm})$$

J_n^{π}	E_{exp} [MeV]	E_{th} [MeV]	J_n^{π}	E_{exp} [MeV]	E_{th} [MeV]
0^+_{gs}	0	0.004	3_{1}^{+}	1.248	1.255
2^{+}_{1}	0.089	0.085	4_{3}^{+}	1.355	1.363
4_{1}^{+}	0.288	0.275	5_{1}^{+}	1.507	1.499
$ 6_1^+ $	0.585	0.574	6^{+}_{3}	1.644	1.662
81	0.965	0.982	7_{1}^{+}	1.85	1.852

TABLE III: Comparison of the experimentally-observed lowenergy states in ¹⁵⁶Gd [14] and excitation energies calculated considering a neutron hole $\nu = 3/2^-$ at $\epsilon_{47} = -6.361$ coupled to the core system ¹⁵⁷Gd with $K = 3/2^-$. The first three columns represent the $\Omega = 0^+$ ground state band; the next three columns represent the $\Omega = 3^+$ band. The theoretical ground state is not exactly zero because ϵ_{47} is not exactly equal the neutron separation energy S_n in ¹⁵⁷Gd.
$$\begin{split} H_C &= -\frac{\hbar^2}{2\mathcal{I}}(I_+ j_- + I_- j_+).\\ \Delta E(\Omega_{\pm} &= |K \pm \nu|) = \pm c\frac{\hbar^2}{2\mathcal{T}}K\nu. \end{split}$$

Coriolis coupling



Excited States in ¹⁵⁶Gd

 $E_{v}^{J,K=\Omega_{v}\pm\Omega_{0}} = E_{F} - BE_{v} + \frac{\hbar^{2}}{2\wp_{1,2}}J(J+1) \pm \frac{\Omega_{v}\Omega_{0}}{\wp_{3}}$





Model space convergence in Sturmian basis



s.p.s. in Sturmian basis



Individual x-sections



Total cross section $\sigma(E)$ smearing function $\rho(E,\Gamma)$ dependence

$$\sigma(E) = \sum_{v} \rho_{v}(E)\sigma_{v}, \ \rho_{v}(E) = \frac{1}{2\pi} \frac{4\Gamma}{4(E - E_{v})^{2} - \Gamma^{2}}, \ \Gamma = a + bE$$



Total cross section σ(E) model space N-dependence

$$\sigma(E) = \sum_{v} \rho_{v}(E)\sigma_{v}, \ \rho_{v}(E) = \frac{1}{2\pi} \frac{4\Gamma}{4(E - E_{v})^{2} - \Gamma^{2}}, \ \Gamma = a + bE$$



P(J^π, E) Distribution



P(J^π) Distributions *E*= 8, 8.5, 9, 9.5 MeV

Ρ

Р





Probability distrinution at E=9 MeV





Outlook

Sturmian basis:

- Guarantees the correctness of the wave-function tail;
- Provides fast convergence!
- Consistency checks:
 - Calculate basis overlap matrix for DWUCK w.f.
 - Compare the basis overlap matrices
- Need better understanding of:
 - the norm issue for few of the states;
 - E-deviations for large model spaces;
 - Smearing function for comparison to experiments!
 - When are the Pairing and Coriolis mixing important?
- Compare to other methods, codes, and experiments ...
- Study the x-sections for the desired reaction n+¹⁵⁵Gd!