

Final Project

TALENT course no. 5 (2017)

Group 1

Introduction

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NushellX - Spectroscopic Factors

- Calculate the spectroscopic factors from the ground state of ^{23}O to all states in ^{22}O in the full $1s0d$ model space. Use the sum rule to obtain the orbital occupations in ^{23}O for $0d_{5/2}$, $1s_{1/2}$ and $0d_{3/2}$. Compare these to those given in the so-called xxx.occ file.

NushellX - Spectroscopic Factors (1)

*.ans file

23O

22O

1 = 1 nucleon transfer

from GS of 23O (1=GS) to 22O (-1=to all)

```
1 -----
2 lpe,  0          ! option (lpe or lan), neig (zero=10)
3 sd              ! model space (*.sp) name (a8)
4 n              ! any restrictions (y/n)
5 usdb           ! interaction (*.int) name (a8)
6 8              ! number of protons
7 23             ! number of nucleons
8 0.5, 6.5, 1.0, ! min J, max J, del J
9 0              ! parity (0 for +) (1 for -) (2 for both)
10 -----
11 lpe,  0          ! option (lpe or lan), neig (zero=10)
12 8              ! number of protons
13 22             ! number of nucleons
14 0.0, 7.0, 1.0, ! min J, max J, del J
15 0              ! parity (0 for +) (1 for -) (2 for both)
16 -----
17 den            ! option
18 1              ! 1, 2 or t
19 bb0600         ! initial file name (default from lpe above)
20 -1             ! max number (-1 for all)
21 bb0701         ! final file name (default = from lpe above)
22 1              ! max number (-1 for all)
23 0.0, 4.0, 1.0, ! min, max J, del J for bb0600
24 0.5, 0.5, 1.0, ! min, max J, del J for bb0701
25 n              ! restrict coupling for operator
26 -----
27 st             ! option
```

NushellX - Spectroscopic Factors (1)

Sum rules

Look at *.occ

N	NJ	Ex	2J	P	0d3	0d5	1s1	0d3	0d5	1s1
1	1	0.000	1	1	0.00	0.00	0.00	0.17	5.81	1.02

Compare with *.lsf

According to sum rules (particle removal) we should get occupation number (40.13):

$$\sum_{f-} S_{i,f,k} = \langle n_k \rangle_i$$

NushellX - Spectroscopic Factors (1)

Eg. $1s_{1/2}$

Sum rule (1.02)

0.9389+

0.0801=

1.019

```
! model space = sd
! interaction = usdb
```

(Ai	Tzi)	(Af	Tzf)	(type n,l,2j)	Ji	Jf	ni	nf	C^2S	Ei	Ef	C^2S*(Ef-Ei)	Exi	Exf
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	1	1	0.7823	-34.498	-37.079	-2.019	0.000	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	2	1	0.1505	-29.736	-37.079	-1.105	4.762	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	3	1	0.0038	-24.849	-37.079	-0.046	9.649	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	4	1	0.0001	-22.361	-37.079	-0.001	12.137	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	5	1	0.0014	-19.986	-37.079	-0.024	14.512	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	6	1	0.0000	-16.492	-37.079	-0.000	18.006	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	7	1	0.0008	-16.341	-37.079	-0.017	18.157	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	8	1	0.0000	-14.791	-37.079	-0.000	19.707	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	9	1	0.0000	-13.558	-37.079	-0.000	20.940	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	0.0+	0.5+	10	1	0.0000	-11.207	-37.079	-0.000	23.291	0.000
sum									0.9389			-3.213		

(Ai	Tzi)	(Af	Tzf)	(type n,l,2j)	Ji	Jf	ni	nf	C^2S	Ei	Ef	C^2S*(Ef-Ei)	Exi	Exf
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	1	1	0.0030	-24.340	-37.079	-0.038	10.158	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	2	1	0.0528	-23.900	-37.079	-0.696	10.598	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	3	1	0.0218	-23.345	-37.079	-0.299	11.153	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	4	1	0.0010	-21.741	-37.079	-0.015	12.757	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	5	1	0.0001	-19.890	-37.079	-0.002	14.608	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	6	1	0.0002	-18.360	-37.079	-0.004	16.138	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	7	1	0.0000	-17.881	-37.079	-0.000	16.617	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	8	1	0.0003	-17.261	-37.079	-0.006	17.237	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	9	1	0.0006	-15.029	-37.079	-0.013	19.469	0.000
(22	3.0)	(23	3.5)	(n 2 0 1)	1.0+	0.5+	10	1	0.0003	-14.893	-37.079	-0.007	19.605	0.000
sum									0.0801			-4.293		

NushellX - Spectroscopic Factors (1)

Sum over all orbitals = Valence neutrons

(22 3.0)	(23 3.5)	(n 1 2 5)	3.0+	0.5+	9	1	0.0005	-18.384	-37.079	-0.009	16.114	0.000
(22 3.0)	(23 3.5)	(n 1 2 5)	3.0+	0.5+	10	1	0.0008	-18.102	-37.079	-0.015	16.396	0.000
						sum	3.3424			-47.421		
						total sum	6.9875					

(22 3.0)	(23 3.5)	(n 1 2 5)	3.0+	0.5+	27	1	0.0000	-5.344	-37.079	-0.000	29.154	0.000
(22 3.0)	(23 3.5)	(n 1 2 5)	3.0+	0.5+	28	1	0.0000	-2.963	-37.079	-0.000	31.535	0.000
(22 3.0)	(23 3.5)	(n 1 2 5)	3.0+	0.5+	29	1	0.0000	2.495	-37.079	-0.000	36.993	0.000
						sum	3.3478			-47.683		
						total sum	6.9997					

NushellX - Spectroscopic Factors (2)

- Calculate the spectroscopic factors from the ground state of ^{23}O to all states in ^{24}O in the full $1s0d$ model space. Use the sum rule to obtain the number of holes in those three orbits in ^{23}O . Compare these to those given in the xxx.occ file.

Similar procedure but different sum rule (**particle addition**), eq. 40.14:

$$\sum_{f+} \frac{(2J_f + 1)}{(2J_i + 1)} S_{i,f,k} = (2j + 1) - \langle n_k \rangle_i$$

NushellX - Exploring gamma decay

Example: ^{59}Ni

Thank you