ELECTRIC DIPOLE PHOTON ABSORPTION IN 32S†

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Summary

The particle—hole model has been applied to the calculation of E1 photon absorption by 32 S. A plausible unperturbed spectrum was obtained from investigation of the results of the 32 S(d,p) 33 S and 32 S(p,2p) 31 P reactions. The results of the calculation are compared with experiment, and suggestions are made with reference to the shortcomings of the calculation.

Introduction

The particle-hole model, proposed by Brown, Castillejo, and Evans (1961) enjoyed a large measure of success in accounting for the features of photodisintegration cross sections of doubly-magic ¹⁶O and ⁴⁰Ca. This success has been extended to the closed subshell nuclei ¹²C and ²⁸Si by Vinh-Mau and Brown (1962) and Bolen and Eisenberg (1964) respectively. The existence of a considerable body of data on ³²S from direct or inverse photodisintegration reactions has motivated the present calculation of 1⁻ states in ³²S.

CALCULATION AND RESULTS

 32 S is a closed subshell nucleus, the $2s_{1/2}$ subshell being completed there. The energies of the single-particle states in the unperturbed 32 S spectrum are extracted in the usual way from data on 33 S and 31 P. The known levels of 33 S, which are strongly excited in the 32 S(d,p) reaction and which have the correct spin and parity, have been taken as the single-particle states of 33 S (McFarlane and French 1960). Thus, the ground state of 33 S is the $1d_{3/2}$ state, and states at $2 \cdot 94$, $3 \cdot 22$, and $5 \cdot 71$ MeV are taken to be the $1f_{7/2}$, $2p_{3/2}$, and $2p_{1/2}$ single-particle states respectively. These states have reduced widths in the stripping reaction, which support these assignments. The $1f_{5/2}$ state is estimated to be about 6 MeV above the $1f_{7/2}$ state; the approximate value of the separation being taken from data on 40 Ca.

The 31 P states are more difficult to locate, and it must be done from two studies of the 32 S(p,2p) reaction. First, Pugh and Riley (1961) report 31 P states corresponding to the ejection of an s-wave proton from 32 S, and a non-s-wave proton, which are separated by 5 MeV. These correspond to the ground state of 31 P(2s_{1/2}) and a state which is presumably the $1d_{5/2}$ single-particle state at 5 MeV excitation. Gottschalk (1962) has also given spectra from this (p,2p) reaction and shows that there is an asymmetric peak which could well be due to the two states just mentioned. Also, one can read into his spectra the presence of two smaller "bumps" which should be

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due to the $1p_{1/2}$ and $1p_{3/2}$ states in ³¹P. These are necessary for the calculation, but, as the results show, play a very small part in it. It must be emphasized that large uncertainties should be attached to the energies of these two states. The whole scheme of particle and hole states is shown in Figure 1.

The residual interaction was taken to be of the form

$$V = V_0(1-\eta+\eta\sigma_1\cdot\sigma_2)\delta(r_1-r_2),$$

where the parameters are taken to be $\eta=0\cdot135$ and $V_0a^3/4\pi=-10\cdot22$ MeV, a being the harmonic oscillator range parameter. The value of $V_0a^3/4\pi$ was fixed by interpolating between the successfully used values of this parameter for ⁴⁰Ca and ²⁸Si, with the proper A-dependence.

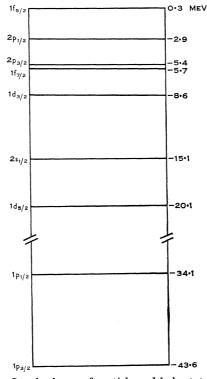


Fig. 1.—Level scheme of particle and hole states in 32S.

The diagonalization of the secular matrices based on the unperturbed spectrum shown in Figure 1, together with the residual interaction indicated, yields the results shown in Table 1 for 1⁻ states. The calculations were performed under the assumption that isospin is a good quantum number, and results for T=0 and T=1 states are presented. This assumption is not expected to be a good one for states in 32 S at such high excitation, and thus considerable isospin mixing is expected to take place.

The lowest T=0 state, given at 1.76 MeV and having a very coherent nature, is simply the spurious state of the shell model due to translational motion of the

nuclear centre of mass. The inclusion of ground state correlations in the calculations will place its energy much closer to zero, where it should be, and increase its coherence. These properties are exactly those expected of the $T=\mathbf{0}$, $\mathbf{1}^-$ spurious state.

Table 1 $\begin{array}{c} {\rm Table\ 1} \\ {\rm energy\ eigenvalues\ and\ eigenvectors\ for\ 1^-\ states\ in\ ^{32}S} \\ E\ {\rm denotes\ the\ transition\ energy.}\ The\ numbers\ in\ the\ second\ row\ are\ unperturbed\ energies\ in\ MeV.} \\ {\rm All\ eigenvectors\ are\ unnormalized.} \end{array}$

		1	l		l	1	
$E \ ({ m MeV})$	$\begin{vmatrix} 2s_{1/2}^{}^{-1}2p_{3/2} \\ 9\cdot 7 \end{vmatrix}$	$\begin{vmatrix} 1d_{5/2}^{-1}1f_{7/2} \\ 14\cdot 4 \end{vmatrix}$	$\begin{array}{c c} 1d_{5/2}^{-1}2p_{3/2} \\ 14\cdot 7 \end{array}$	$\begin{bmatrix} 2s_{1/2}^{-1}2p_{1/2} \\ 12 \cdot 2 \end{bmatrix}$	$\begin{array}{c c} 1d_{5/2}^{-1}1f_{5/2} \\ 20\cdot 4 \end{array}$	$\begin{vmatrix} 1p_{1/2}^{-1}1d_{3/2} \\ 25 \cdot 5 \end{vmatrix}$	$\begin{vmatrix} 1p_{3/2}^{-1}1d_{3/2} \\ 35 \cdot 0 \end{vmatrix}$
			<u>'</u>	T = 1			
11.89	1.000	-0.155	0.016	-0.188	0.020	-0.010	0.007
$14 \cdot 07$	$0 \cdot 152$	-0.234	0.066	1.000	-0.091	-0.024	-0.026
$16 \cdot 77$	-0.061	-0.177	1.000	-0.105	-0.096	0.096	0.023
$19 \cdot 07$	0.183	1.000	$0 \cdot 254$	0.204	0.220	-0.248	0.009
$22 \cdot 26$	-0.057	-0.263	0.064	0.028	1.000	-0.133	-0.115
$28 \cdot 29$	0.053	0.199	-0.029	0.072	0.159	1.000	-0.255
$38 \cdot 02$	0.004	0.010	-0.023	0.048	0.148	$0 \cdot 225$	1.000
			<u>′</u>	T = 0			
1.76	0.900	1.000	0.051	0.755	0.136	0.321	0.105
$8 \cdot 34$	1.000	-0.702	-0.197	-0.177	-0.073	-0.130	-0.038
11.81	0.148	-0.005	1.000	-0.209	0.016	-0.068	-0.014
$13 \cdot 63$	-0.178	-0.507	0.219	1.000	-0.142	-0.227	-0.062
$20 \cdot 86$	-0.026	-0.154	-0.001	0.058	1.000	0.001	-0.026
$23 \cdot 59$	-0.059	-0.302	0.085	0.041	-0.050	1.000	0.050
$35 \cdot 06$	-0.018	-0.072	0.012	0.005	0.014	-0.073	1.000

Calculat	ion I	Calculat	ion II	Calculation III		
$V_0 \alpha^3 / 4\pi = -$	$10 \cdot 22 \text{ MeV}$	$V_0 a^3/4\pi = -$	$-9.5~\mathrm{MeV}$	$V_0 a^3 / 4\pi = -10 \cdot 22 \text{ MeV}^*$		
$E \hspace{1cm} M ^2$		· · E	$ M ^{2}$	E	$ M ^2$	
(MeV)	(%)	(MeV)	(%)	(MeV)	(%)	
11.89	1	11.75	1	11.89	1	
$14 \cdot 07$	0	13.94	0	$14 \cdot 07$	0	
$16\cdot 77$	0	16.63	0	16.78	0	
$19 \cdot 07$	76	18.78	78	19.10	77	
$22 \cdot 26$	5	22 · 12	4	23.12	$3\frac{1}{2}$	
$28 \cdot 29$	17	28.09	16	28.07	16	
38.02	1	37 · 76	1	$35 \cdot 34$	$2\frac{1}{2}$	

^{*} Spectrum changed.

The eigenvectors obtained from the diagonalization were also used in the calculation of the dipole strengths of the T=1 states, and these are shown in Table 2 (calculation I). Also shown in Table 2 are the dipole strengths which result from calculations using two variations in the basic assumptions. In calculation II, the

TABLE 3

COMPARISON OF PRESENT CALCULATION WITH EXPERIMENTAL RESULTS

	-	COMPARISO	N OF PRESENT C.	COMPARISON OF PRESENT CALCULATION WITH EXPERIMENTAL RESULTS	KIMENTAL KESU	LIS		
Author:	Thompson and Taylor*	Firk (1964)	Mutsuro <i>et al.</i> (1963)	Bolen and Whitehead (1963)	Kimura <i>et al.</i> (1963)	Gemmell and Jones (1962)	Present Calculation	ent
Reaction studied:	$^{32}\mathrm{S}(\gamma,\mathrm{n})$	$^{32}\mathrm{S}(\gamma,\mathrm{n})$	$^{32}\mathrm{S}(\gamma,\mathrm{n})$	$^{82}\mathrm{S}(\gamma,\mathrm{n})$	$^{81}\mathrm{P}(\mathrm{p},\gamma_0)$	$^{31}\mathrm{P}(\mathrm{p},\gamma_0)$	T = 0	T = 1
						3	9	11.89
						13.5 14.3	13.62	14.06
						15.0		
	15.6		15.7		15.7	15.6		
	15.9			16.0	16.1	16.0		
	(16.7)		16.8		16.9	16.8		16.77
	17.4	17.4		17.5	17.8	17.6		
	17.6	17.7						
		17.9						
	18.05	18.1	18.0	18.25	18.2	18.1		
	18.4	18.4						
	18.8	18.8	18.7		18.8	18.9		
	19.1	19.0	19.3		19.1			19.07
		19.4						
	19.6	$\begin{array}{c} 19.8 \\ 20.2 \end{array}$	19.7	19.75	19.75			
*	(20.5)	$20 \cdot 6$	20.5		20.5		20.86	
	(21.5)			21.25				
				22.75				$22\cdot 26$
				24.0	•		$23 \cdot 59$	
							-	$28 \cdot 28$
							35.06	
								38.02

* Personal communication, 1964.

residual interaction strength was set at $V_0\alpha^3/4\pi = -9\cdot 5$ instead of $-10\cdot 22$ MeV. In calculation III, the assumed positions of the $1p_{3/2}$ and $1f_{5\ 2}$ states were shifted upward by 3 and 1 MeV respectively, with $V_0\alpha^3/4\pi$ set equal to $-10\cdot 22$ MeV. It is to be noted that the relative dipole strengths, taken as $|M|^2$ where $M = \int \psi_f^* z \psi_i \, d\tau$, are quite insensitive to these parameter variations. Also, the eigenvectors showed little change due to the two variations. However, it must be emphasized that the wave functions given by a calculation such as this are not considered to be reliable, so that the only firm prediction which can be made is that most of the dipole strength is concentrated in states at 19 and 28 MeV.

COMPARISON WITH EXPERIMENT

The states predicted by this calculation are to be compared with states found in the $^{31}\text{P}(p,\gamma_0)^{32}\text{S}$ reaction (Gemmell and Jones 1962; Kimura *et al.* 1963), in the measurement of the neutron spectra from $^{32}\text{S}(\gamma,n)^{31}\text{S}$ by time-of-flight methods (Firk 1964; Mutsuro *et al.* 1963), and in direct measurements of the $^{32}\text{S}(\gamma,n)^{31}\text{S}$ cross section by yield-curve techniques (Bolen and Whitehead 1963; Thompson and Taylor 1964, personal communication).

The comparison of the calculation with these experimental results is shown in Table 3. It should be noted that only the strong peaks obtained in Firk's measurement are quoted. This is done so that the experimental results may be compared with approximately the same resolution. This approximation in no way removes the necessity for understanding the finer structure.

There exist at present no experimental data above an excitation of 21 MeV, so that the prediction of much cross section at 28 MeV cannot be tested as yet. It is of interest, however, that the possibility exists for the use of the photodisintegration reaction to determine the energies of the two 1p states in 32 S.

DISCUSSION

It is apparent that there are many more states in the experimental data than are predicted by the present calculation. The extra peaks may be due to M1 or E2 excitations which were not included in the calculation, or they could be due to a deficiency of the particle—hole model.

This model takes the ground state of the initial nucleus to be the physical vacuum, and does so irrespective of any properties of this ground state. For example, the ³²S and ²⁸Si nuclei are expected to be softer against deformation than the doubly-magic ¹⁶O and ⁴⁰Ca nuclei, but the simple model is not able to build in this fact. Perhaps it would show in terms of ground state correlations being much more important for ³²S and ²⁸Si, if these correlations should be included in the calculation.

The other limiting feature of the model is that it assumes pure j-j coupling, and thus assumes that all of the single-particle strength is concentrated in one level. This is clearly an idealization, and an alternative much nearer reality was pointed out by Lane, Thomas, and Wigner (1955) to cope with a different situation. If one takes the view that the single-particle nature is concentrated in a given region of excitation, but is spread among many levels with some sort of peaked distribution

about the mean, then in nuclei where the level densities are relatively low there will be more strongly absorbing 1^- states than the simple particle–hole model predicts. This may well be the case in 32 S, where, for example, McFarlane and French (1960) state that the $1f_{7/2}$ state is apparently distributed among three or more levels.

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