Comparison of sub-Coulomb (d, p) stripping to analog resonance results for the $N = 82$ isotones*

G. A. Norton, H. J. Hausman, and J. F. Morgan Department of Physics, The Ohio State University, Columbus, Ohio 43210 (Received 6 August 1973)

Reduced normalizations have been extracted from proton elastic scattering data via isobaric-analog resonance for analogs of low-lying parent states in ^{139}Ba , ^{141}Ce , ^{143}Nd , and 145 Sm, by the use of three different analog resonance theories. These reduced normalizations were compared to those obtained from sub-Coulomb (d, p) stripping to the low-lying parent states of the above nuclei. This comparison shows that the R-matrix theory gives the best agreement to the (d, p) results.

NUCLEAR REACTIONS 138 Ba, 140 Ce, 142 Nd, 144 Sm (d,p) comparison to (p,p₀) results. Calculated Λ for (p, p_0) IAR results.

I. INTRODUCTION

In the study of sub-Coulomb (d, p) stripping reactions a quantity, known as the reduced normalization Λ_{ij} , has been shown to be insensitive to the optical-model parameters used in the distortedwave Born-approximation (DWBA) analysis.^{1,2} $:$ to stor $\frac{1}{2}$ This same quantity can be extracted from (p, p_0) isobaric analog resonance (IAR) data by the use of three different IAR theories.³ By comparison of sub-Coulomb (d, p) stripping and the (p, p_0) IAR results one may be able to choose among the three IAR theories. In the past this comparison has been attempted by comparing spectroscopic factors $S_{\bm{i}}$, derived from $(p,p_{\scriptscriptstyle 0})$ scattering via IAR with those found from (d, p) stripping to the low-lying parent states.

In the DWBA analysis of the (d, p) data, S_{ij} is strongly dependent on the optical-model parameters used. In many cases, the analysis of the (d, d) elastic scattering data leads to several equally good families of parameters, which when applied to the (d, p) reaction yield spectroscopic factors which may differ by as much as 50% (e.g. Ref. 4). The dependence of S_{ij} on the deutero and proton potential parameters can be strongly reduced by performing the (d, p) experiments at energies in which both entrance and exit channels are below the Coulomb barrier. However, the dependence on the bound-state neutron potential parameters does not decrease appreciably below the Coulomb barrier. Thus, S_{ij} still cannot be determined uniquely. However, the reduced normalization, which is closely related to S_{ij} , can be determined uniquely for both the sub-Coulomb (d, p) stripping reactions and the (p, p_0) IAR experiments.

 Λ_{ij} is essentially the square of the ratio of the

transferred neutron's asymptotic wave function to a spherical Hankel function, and is related to S_{ij} by

$$
k^3 \Lambda_{ij} = N_{ij}^2 S_{ij}
$$
,

with

$$
k = (2\mu |B_n|)^{1/2}/\hbar ,
$$

where $N_{i,j}$ is the ratio of the DWBA neutron boundstate wave function to a spherical Hankel function evaluated outside the nuclear radius, μ is the reduced mass in the exit channel, and B_n is the binding energy of the last neutron. Rapaport and Kerman¹ have shown that Λ_{IJ} is nearly independe of the geometrical parameters used to describe the neutron bound-state potential for sub-Coulomb (d, b) stripping. This has also been shown by Kent, Morgan, and Seyler⁵ and Norton et al.²

Clarkson, Von Brentano, and Harney' have extracted reduced normalizations from the (p, p_0) reactions in the context of three IAR theories: the R -matrix approach of Thompson, Adams, and Robson' (TAR) and two shell-model methods, that of Mekjian and McDonald' (MM) and that of Zaidi and Darmodjo⁸ and Harney⁹ (ZDH). These theories have been used to yield the single-particle proton width Γ_{\bullet}^{sp} of the analog state. Clarkson et al. have defined a term called the reduced single-particle proton width G_{ρ} , which is related to Γ_{ρ}^{sp} by

$$
\Gamma_{p}^{\mathbf{sp}}=\frac{N_{lj}^{2}}{k^{3}}G_{p}.
$$

For (p, p_0) reactions, the spectroscopic factor is given by

$$
S_{ij} = \Gamma_p / \Gamma_p^{sp}.
$$

Then, the experimentally measured proton partial width Γ_{p} is related to the reduced normalization

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by $\Gamma_{p} = \Lambda_{ij} G_{p}$. Thus, to determine the value of Λ_{IJ} from (p, p_0) IAR experiments, G_p is calculate from one of the three theories and compared to the experimental value of Γ_{p} . However, the value of G_{ρ} is strongly dependent on the IAR theory used, and therefore the value of the reduced normalization, which can be determined uniquely from sub-Coulomb (d, p) to the parent state, is dependent on the analog resonance theory used in its extraction from (p, p_0) scattering data.

Harney and Weidenmüller¹⁰ (HW) have compared these theories on a theoretical level and have shown that there are fundamental differences which can produce spectroscopic factors that may differ by as much as 50% . It is now possible to compare these theories on an experimental level by use of the reduced normalization.

In this paper reduced normalizations are calculated for each of the three theories mentioned above, using analog resonance parameters found
in the available literature for the $N=82$ isotones in the available literature for the $N = 82$ isotones;
¹³⁸Ba, ¹⁴⁰Ce, ¹⁴²Nd, and ¹⁴⁴Sm. These reduced normalizations are compared to those reported by Rapaport and Kerman¹ for sub-Coulomb (d, p) stripping to parent states in ^{139}Ba , and by Norton stripping to parent states in \mathcal{A} , and by Not ton
et al.² for sub-Coulomb stripping to parent states in 141 Ce, 143 Nd, and 145 Sm. This comparison is a continuation of that started by Morgan, Seyler, and Kent¹¹ near the $N=50$ region. The opticalmodel parameter dependence of Λ_{IJ} from the various IAR theories is also discussed. The preliminary results of the comparisons, for both the $N=50$ and the $N=82$ region, have been previously reported.¹² reported.

II. OPTICAL-MODEL PARAMETERS FOR (p, p_0) IAR IAR ANALYSIS

Reduced normalizations were calculated from each of the three theories by use of code BETTINA¹³ with the optical-model parameters taken from with the optical-model parameters taken from
Wiedner et al.¹⁴ The proton partial widths were Wiedner *et al*.¹⁴ The proton partial widths wer obtained from Williams *et al*.¹⁵ for ¹³⁸Ba, Marobtained from Williams *et al*.¹⁵ for ¹³⁸Ba, Mar-
quardt *et al*.¹⁶ for ¹⁴⁰Ce, Grosse *et al*.¹⁷ for ¹⁴²Nd
and Fiarman *et al*.¹⁸ for ¹⁴⁴Sm, and are shown in and Fiarman et al.¹⁸ for 144 Sm, and are shown in Table I.

The optical-model potential used in code BETTINA ls

$$
U(r) = V(1 + e^{x})^{-1} + V_c(r, r_{0c}) + 4i W \frac{e^{x'}}{(1 + e^{x'})^2}
$$

$$
+ \alpha(l) V_{\infty} \left[\frac{-1.998}{r} \frac{d}{dr} (1 + e^{x})^{-1} \right],
$$

where

$$
x = \frac{\gamma - \gamma_0 A^{1/3}}{a}, \quad x' = \frac{\gamma - \gamma_0' A^{1/3}}{a}
$$

and

$$
\alpha(l) = \begin{cases} l \text{ if } j = l + \frac{1}{2}, \\ -l - 1 \text{ if } j = l - \frac{1}{2}. \end{cases}
$$

				(p,p_0) IAR			
Parent	E_{x}	l_j	Γ_{ϕ} (keV)	Λ_{MM}	Λ_{TAR}	Λ_{ZDH}	Λ_{dp} ^a
139 Ba	0.00	$f_{1/2}$	17.2	26	21	15	19.6°
	0.63	$p_{3/2}$	26.0	78	88	42	107
	1.08	$\boldsymbol{p}_{1/2}$	22.5	58	58	28	66
	1.42	$f_{5/2}$	9.5	1.41	1.20	0.75	\cdots
141 Ce	0.00	$f_{7/2}$	12.3	50	39	26	33
	0.67	$p_{3/2}$	21.8	149	145	86	168
	1,14	$\rlap/p_{1/2}$	20.4	124	103	51	104
	1,50	$f_{5/2}$	7.3	3.4	2.6	1.6	2.6
	1,78	$f_{5/2}$	(7.2)	3.2	2.3	1.5	2.1
	2.41	$(p_{1/2})$.	.	\ddotsc	\cdots	36
		$(p_{3/2})$	\cdots	.	.	\cdots	19
143 Nd	0.00	$f_{7/2}$	10.5	111	76	52	49
	0.74	$p_{3/2}$	23.5	344	272	139	250
	1,31	$P_{1/2}$	22.7	365	210	115	165
	1,56	$f_{5/2}$	6.0	6.26	4.32	2.63	5.08
	1.92	$f_{5/2}$	\ddotsc	.	.	.	1.68
145 Sm	0.00	$f_{1/2}$	8	212	123	84	77
	0.89	$p_{3/2}$	27	944	517	282	442
	1,61	$p_{1/2}$	30	789	350	196	367

TABLE I. Final results.

^a Taken from Norton *et al.* (Ref. 2) and from Rapaport and Kerman (Ref. 1) for ^{139}Ba .

The factor of 1.998 is the square of the Compton wavelength of the π meson in femtometers and V_c is the Coulomb potential of a uniformly charged sphere of radius $r_{0c}A^{1/3}$.

In order to calculate $\Gamma^{\mathfrak{sp}}_b$, code BETTINA assumes that the last neutron in the parent nucleus is in a single-particle state. Therefore, since the lowlying states of the $N = 83$ isotones are not pure single-particle states, the binding energy required to determine V_n is not the binding energy of the last neutron, but a single-particle binding energy, E_{Bn} , given by

$$
E_{Bn} = Q_{\rm g.s.}(d, p) - \epsilon_{\rm sp} + 2.224 \text{ MeV}
$$

and

$$
\epsilon_{sp} = \frac{\sum S_{ij} E_{ij}}{\sum S_{ij}}
$$

where the sum is over states of the same spin and parity, and E_{ij} is the excitation energy of the state whose spectroscopic factor is S_{ij} . Estimates of $\epsilon_{\rm sp}$ are available in the literature¹⁹⁻²² from (d, p) work done above the Coulomb barrier.

As mentioned earlier, the reduced normaliza-

FIG. 1. Results of the variation of the neutron well radius for the ¹⁴⁰Ce(p, p₀) reaction. r_n^0 was varied from 0.9 to 1.4 fm while all other parameters were held constant. Although each theory exhibits a different r_x^0 dependence, in every case, Λ_{ij} is less dependent on r_n^0 than is S_{ij} .

tions obtained from sub-Coulomb (d, p) stripping reactions are nearly independent of the opticalmodel parameters used to describe the neutron bound-state well, as opposed to the strong dependence exhibited by the spectroscopic factors obtained from the sub-Coulomb (d, p) reactions. For the proton elastic scattering data, Λ_{ij} is again more insensitive to the neutron parameter than is S_{ij} , as shown in Fig. 1. Here, as in parameter variations to follow, only one parameter was varied; in this case the neutron bound-state well radius, while all others were held constant.

Of the three theories, the R -matrix approach of TAR is the least dependent on the neutron radius for S_{ij} . For this theory, S_{ij} decreases by 80%, while a 90% decrease is noted in the MM and ZDH theories. However, for Λ_{ij} the TAR theory exhibits the greatest neutron radius dependence, since Λ_{1i} increases almost 50%, while the MM value increases by 20% and Λ_{ij} from ZDH decreases by I2%.

In the proton channel the potential parameters were varied individually to see hom the various theories differed as to parameter dependence. Figure 2 shows the dependence of S_{ij} and Λ_{ij} on

FIG. 2. Results of the variation of the absorption potential depth for the 140 Ce(p,p₀) reaction. W was varied from 0 to 25 MeV in code &ETTINA, while all other parameters were held constant.

the surface absorption potential W which was varied from 0 to 25 MeV. At $W = 0$, MM and ZDH give identical values for S_{ij} and Λ_{ij} with TAR and ZDH theories showing identical W dependence, which is due to the fact that TAR and ZDH have nearly the same expressions for the resonance mixing phase same expressions for the resonance mixing phase
 ϕ_c , as noted in Harney and Weidenmüller.¹⁰ In the variation of the real proton potential (Fig. 2), strong peaking was noted in the values of S_{ij} and Λ_{1} , at V_{\bullet} ~ 50 MeV for the MM theory only. An identical curve is obtained when the real potential radius r_{00} is varied, with a peak occurring at r_{00} =1.15 fm for the MM theory. Also the spin-orbit potential depth $V_{\rm so}$ was varied from 0 to 10 MeV, which produced changes of 5 to 10% in S_{ij} and Λ_{ij} for all three theories, with the MM theory showing the greatest dependence.

These parameter variations illustrate some of the substantive differences that exist among the

FIG. 3. Results of the variation of the real proton potential depth for the ¹⁴⁰Ce(p,p₀) reaction. V_p was varied from 30 to 70 MeV in code BETTINA, while all other parameters were held constant.

three theories. These differences are explored in detail by Harney and Weidenmüller.¹⁰

III. COMPARISON OF A_{ij} FROM IAR AND (d,p) RESULTS

In obtaining Λ_{1} , from the analog resonance data, the largest uncertainties are from the proton partial widths and the center-of-gravity excitation energy ϵ_{sp} . Assuming that ϵ_{sp} is known to within 150 keV, an average uncertainty of $\pm 10\%$ can be assigned to Λ_{1} , due to this parameter. This uncertainty, combined with the 6 to 12% experimental errors associated with Γ_{ϕ} , leads to an estimate of $\pm 20\%$ error in Λ_{14} . The values of the reduced normalizations are listed in Table I, along with those obtained from sub-Coulomb (d, p) stripping reac- $\frac{\text{total}}{\text{cross}}$ if $\frac{\text{total}}{\text{cross}}$ and Rapaport and Ra Kerman.¹ These reduced normalizations also have uncertainties of $\pm 20\%$.

The reduced normalizations with their associated uncertainties versus excitation energy are shown in Fig. 4. The cross-hatched area represents Λ_{ij} obtained from the (d, p) analysis. Note that in each case the error bars of at least one of the analog resonance theories overlap with the (d, p) results. From this figure and Table I it can be seen that the R-matrix approach of TAR comes closest to the (d, p) reduced normalization a total of 11 times, with MM and TAR producing the same values of Λ_{ij} for the $p_{1/2}$ state of ¹³⁹Ba.

In order to determine which of the analog resonance theories has the best agreement to the (d, p) reduced normalizations, a "goodness-of-fit" parameter¹¹ I is defined similarly to χ^2 . This parameter is given by

$$
I = \sum_{\text{states}} \frac{(\Lambda_{dp} - \Lambda_{pp})^2}{(\Delta \Lambda_{dp})^2 + (\Delta \Lambda_{pp})^2} ,
$$

where the sum is over states of the same spin and parity and Δ is equal to 0.2, representing the 20% uncertainty in Λ_{dp} and Λ_{pp} . Note that if I is equal to 1 for a given state the difference between Λ_{ab} and $\Lambda_{\rho\theta}$ is 1.4 Δ , since there are two equal terms

TABLE II. Goodness-of-fit parameter for the $N = 82$ isotones.

	No. of				Isotones
State	states	мм	TA R	ZDH	included
$2f_{1/2}$	4	18.6	5.3	1.7	Ba. Ce. Nd. Sm
$3p_{3/2}$	4	8.4	1.1	18.8	Ba. Ce. Nd. Sm
$3p_{1/2}$	4	12.7	0.9	18.2	Ba. Ce. Nd. Sm
$2f_{5/2}$	3	3.5	0.4	8.7	Ce. Nd
Total	15	43.1	7.8	47.5	
Total/States		2.9	0.5	3.2	

FIG. 4. The reduced normalization versus excitation energy for states in the $N = 83$ isotones. The cross-hatch marks denote the value of the reduced normalizations obtained from the sub-Coulomb (d, p) stripping reactions.

in the error denominator. In the usual case of only one error term a χ^2 of 1 implies that the quantities agree to within one Δ . In this case, however, I must be less than 0.707 for the Λ 's to agree within one Δ .

Table II shows the comparison of the three theories to the experimental (d, p) results. For. each state except the $2f_{7/2}$ state, TAR does give the best agreement for the 139 Ba and 141 Ce isotones; only in 143 Nd and 145 Sm does the ZDH approach come closer to $\Lambda_{d\rho}$. It should be noted that the $2f_{7/2}$ ground states of ¹⁴³Nd and ¹⁴⁵Sm are the states closest to the top of the Coulomb barrier, within 8% and 5% respectively, for the proton channel, thereby causing the reduced normalizations for these states to have a greater dependence on the outgoing channel's potential parameters. Also, the TAR approach yields the lowest total value of I , and is the only theory that produces an average value for I per state that is less than 0.707.

IV. CONCLUSION

The sub-Coulomb (d, p) stripping comparisons to the proton elastic analog resonance reactions for the $N=82$ isotones indicate rather strongly that the theory of Thompson, Adams, and Robson gives the best agreement to the experimental results in this mass region. This supports the work done by this mass region. This supports the work done by Morgan, Seyler, and Kent $et\ al.^{11}$ near the N =50 region. The reason why the TAR R -matrix approach yields the best over-all agreement to the sub-Coulomb (d, p) stripping results is not clear. In their extensive comparison of these three theories, HW pointed out that true substantive differences exist among the theories. In particular both of the shell-model approaches utilize statistical assumptions in the construction of the analog states. These assumptions ignore second-order effects in the imaginary optical potential W , an assumption which HW show is violated even for very small values of W . Also, HW show that application of an R -matrix theory to analog resonances appears to violate the R-matrix assumptions of no internal mixing and no external polarizing potential. It may be that the better results obtained with the TAR method is an indication of which of these violations has a stronger effect on the calculation. It should be pointed out that the R-matrix approach does have an adjustable parameter, the channel radius, outside of which

there exists no nuclear potential for the channel in question. In every case this parameter was set at the first maximum in the single-particle proton width outside the nuclear radius. Since this was done automatically by the code BETTINA for every state in this comparison, it is not clear

that this extra parameter significantly influenced the results.

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