neglected. One finds that

$$\frac{E_{\omega_3}}{E_{\omega_2}} = \frac{3\pi n_0 e^{2l}}{2E_G q_3} \left(\frac{\omega_3}{\omega_2}\right) \frac{1 + 8E_F / 5E_G}{(1 + 4E_F / E_G)^{5/2}} \left(\frac{eE_{\omega_1}}{m * \omega_1 c}\right)^2.$$
(11)

Here E_{ω_1} , E_{ω_2} , and E_{ω_3} are the amplitudes of the three beams, which have been assumed collinear and of parallel polarization. To facilitate comparison with PSF, it is also convenient to rewrite this expression in terms of a nonlinear susceptibility and the beam powers. The formula which results is the same as the equation in their Table I, with $\chi^{(3)}$ defined as

$$\chi^{(3)} = \left[\frac{n_0 e^4}{4(m^*)^2 E_G \omega_1^2 \omega_2 \omega_3}\right] \left[\frac{1+8E_F / 5E_G}{(1+4E_F / E_G)^{5/2}}\right].$$
(12)

As shown in PSF, numerical values calculated from this formula are in reasonable agreement with experiment. The formula also gives the correct ratio of susceptibilities for InAs and InSb. This ratio is more accurately known experimentally than are the absolute values of $\chi^{(3)}$. Finally, Eq. (12) accounts for the measured variation of P_{ω_3} with electron density in InAs, for densities above $10^{16}/\text{cc.}$ This power is proportional to the square of $\chi^{(3)}$ and varies less rapidly than n_0^2 because of the second term in Eq. (12), which reduces the nonlinearity at higher doping levels. Crudely speaking, this reduction can be thought of as arising from an increase in carrier mass with doping.

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⁴L. Spitzer, <u>Physics of Fully Ionized Gases</u> (Interscience Publishers, Inc., New York, 1962).

⁵E. O. Kane, J. Phys. Chem. Solids <u>1</u>, 249 (1957). ⁶The use of a single-band Hamiltonian to describe the response of an electron to a perturbation is only correct when the perturbing frequencies are small compared with E_G/\hbar . This criterion is not particularly well satisfied in our case. Thus one should expect corrections of order $(\hbar \omega_1/E_G)^2 \lesssim \frac{1}{4}$ to succeeding formulas.

⁷J. A. Giordmaine, private communication.

ISOSPIN MIXING IN DEUTERON REACTIONS

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The reaction $C^{12}(d, \alpha)B^{10}$ has been employed recently¹ to study the isospin impurity of N¹⁴ as a function of excitation energy, since the deuteron, the alpha particle, and the C¹² nucleus are in T=0 states. It has been suggested² that since the deuteron is distorted by the electric field of the target, some admixture of T=1 states will be produced and will contribute to the observed isospin impurity. A crude upper limit of 10% was obtained² for the similar reaction Ca⁴⁰(d, α)K³⁸ by assuming that the neutron and proton are completely uncoupled in the nuclear Coulomb field.

An adiabatic approximation has been applied by the author³ to evaluate the polarization potential acting on a deuteron at any distance from a fixed point charge. In this Letter we will show that this previous work³ implies a value for the isospin impurity carried into the reaction by the distorted deuteron, and we will present numerical results.

Expanded in relative partial waves, the deuteron wave function has the form³

$$\psi(\mathbf{\vec{r}},\mathbf{\vec{x}}) = \varphi(\mathbf{r}) \left[1 + \sum_{L=0}^{\infty} F_L(x,r) P_L(\cos\theta) \right], \quad (1)$$

where $\vec{\mathbf{x}}$ is the center-of-mass coordinate of the deuteron measured from the target nucleus, the relative coordinate $\vec{\mathbf{r}} = \vec{\mathbf{r}}_n - \vec{\mathbf{r}}_p$, the deuteron ground state is approximated by

$$\varphi(r) = (\gamma/2\pi)^{1/2} e^{-\gamma r}/r, \qquad (2)$$

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and $F_L(x, r)P_L(\cos)$ is the multipole distortion of order L induced in the deuteron while it is outside a nucleus of charge Ze, and is orthogonal to $\psi^2(r)$.

Since no spin flip is produced by the Coulomb field, a simple relation between isospin and internal orbital angular momentum of the neutron-proton system holds: Even L corresponds to T=0 and odd L corresponds to T=1. It is thus easily seen that the ratio R of T=1 to T=0 intensity is given by⁴

$$R = \frac{2\gamma \sum_{\text{odd } L} (2L+1)^{-1} \int_{0}^{\infty} dr \, e^{-2\gamma r} F_{L}^{2}(x,r)}{1+2\gamma \sum_{\text{even } L} (2L+1)^{-1} \int_{0}^{\infty} dr \, e^{-2\gamma r} F_{L}^{2}(x,r)}.$$
(3)

An excellent approximation⁵ to R is obtained by keeping only the leading (L=1) term, and using the results of Ref. 3 we find the following analytic expression for R:

$$R = Z^{2} \alpha^{2} \{ (8t^{4})^{-1} + \text{Ei}(-2t) [\frac{4}{3}t^{2} - 2t + 2 + t^{-1}(e^{-2t} - 1)] + e^{-2t} [\frac{1}{3}(2t - 4 + 5/t) + t^{-4}(-\frac{3}{2} + 2t - 5t^{2}/2)] + e^{-4t}t^{-4} [11/8 + \frac{1}{2}t + \frac{1}{2}t^{2}] \},$$
(4)

where $t = 2\gamma x$ and $\alpha = Me^2/\gamma\hbar^2$. (Numerically, the inverse deuteron radius $\gamma = 2.316 \times 10^{12} \text{ cm}^{-1}$, and $\alpha = 0.150$.) The universal function G(x) $\equiv Z^{-2}R$ is plotted in Fig. 1 against the radial variable x.

To make use of the curve in Fig. 1, one determines x_0 , the effective radius of the target. Then $Z^2G(x_0) \equiv R$ gives the isospin impurity in the deuteron wave function just before the com-



FIG. 1. The universal function G(x).

pound nucleus is formed. Assuming that $x_0 = r_0 A^{1/3}$ and $T_z = 0$ (i.e., $Z = \frac{1}{2}A$), we have plotted R in Fig. 2 as a function of the mass number A, for three values of r_0 (1.2, 1.4, and 1.6 F).

For the case reported in Ref. 1, we obtain R = 0.67, 0.55, and 0.44% for these three choices of r_0 , compared with the measured isospin impurity of 1-2% at 11 MeV and about 10% at 9 MeV. At these energies, the adiabatic approximation is expected to overestimate the effect.



FIG. 2. Isospin impurity percentage $R = Z^2 G$ versus mass number A, for three choices of r_0 , where the interaction radius $x_0 = r_0 A^{1/3}$.

There is some evidence,⁶ however, which supports the use of a larger asymptotic normalization constant for the deuteron wave function than that of Eq. (2), and an increase of R by as much as a factor of 1.5 might result.

The numerical evaluation of G(x) from Eq. (4) was done by Edward Monasterski using the IBM 7094 at the Goddard Laboratory for Theoretical Studies.

¹L. Meyer-Schützmeister, D. von Ehrenstein, and R. G. Allas, Phys. Rev. <u>147</u>, 743 (1966).

²Y. Hashimoto and W. P. Alford, Phys. Rev. <u>116</u>, 981 (1959). [These authors acknowledge a private communication from J. B. French.]

³R. J. Drachman, Phys. Rev. <u>132</u>, 374 (1963). [See also the erratum, Phys. Rev. <u>139</u>, AB4 (1965), which does not affect the present application.]

⁴The original form of the denominator should be

$$2\gamma \int_{0}^{\infty} dr \, e^{-2\gamma \, r} [1 + F_{0}]^{2} + 2\gamma \sum_{\text{even } L > 0} (2L + 1)^{-1} \int_{0}^{\infty} dr \, e^{-2\gamma \, r} F_{L}^{2}$$

Now, evaluating the first integral we have

$$2\gamma \int_{0}^{\infty} dr \, e^{-2\gamma \, r} [1 + 2F_{0} + F_{0}^{2}]$$

= 1 + 4\gamma \int_{0}^{\infty} dr \, e^{-2\gamma \, r} F_{0} + 2\gamma \int_{0}^{\infty} dr \, e^{-2\gamma \, r} F_{0}^{2}.

The second term vanishes because of the orthogonality of the ground-state function $|0\rangle$ and the first-order perturbation correction

$$F\varphi = \sum_{n \neq 0} \frac{|n\rangle \langle n|V|0\rangle}{E_0 - E_n}.$$

⁵An estimate of the error incurred by neglecting L > 1 terms can be obtained by using the approximate form $F_L \alpha (r/2x)^{L+1}/(L+1)$ for r < 2x, and $F_L = 0$ for r > 2x (see Ref. 3, Sec. IV). Then it is easily shown, for example, that the L = 3 term is about 5% of the L = 1 contribution at x = 2, and 3% at x = 4.

⁶C. F. Clement, Phys. Rev. <u>128</u>, 2724 (1962).

SU(3) ASSIGNMENT AND COUPLING OF $N*(1688)*\dagger$

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A comparison of the two reactions

$$\gamma + p \rightarrow \eta + p$$

and

$$\gamma + p \rightarrow \pi^0 + p$$

in the region of the third nucleon isobar $N^{***}(1688)$ has produced evidence that (1) $N^{***}(1688)$ is a member of a unitary octet, and (2) the D/Fratio, which relates the couplings $N^{***N\pi}$ and $N^{***N\eta}$, is similar to that which relates $NN\pi$ and $NN\eta$. One will note that (1) is in accord with the usual hypothesis that the N^{***} is the first Regge recurrence of the nucleon. Point (2) is not predicted by any existing theory but may have some simple dynamical origin or follow from some higher symmetry.

The above conclusions are arrived at as follows: An $I = \frac{1}{2}$ object decaying into πN (like N^{***}) must, in SU(3), belong to 8, 10*, or 27. Of these possible assignments, we may immediately rule out 10* since $\gamma + p \rightarrow 10^*$ is forbidden by SU(3)¹ whereas N^{***} is strongly observed in pion photoproduction. To decide between <u>8</u> and <u>27</u>, we compare the couplings $\gamma_N * * *_{N\pi^0}$ and $\gamma_N * * *_{N\eta}$; according to SU(3), the ratios are

$$\gamma_{N^{***N\eta}} / \gamma_{N^{***N\pi}} = 3 \text{ if } N^{***} \in \underline{27},$$
$$= \frac{1}{3} (3 - 4\alpha)^2 \text{ if } N^{***} \in \underline{8}, \quad (1)$$

where $\alpha/(1-\alpha)$ is the D/F ratio. It is worth noting at this point that the ratio of the couplings can be small only if N^{***} belongs to an octet with $\alpha \sim \frac{3}{4}$.

To compare with experiment, we write²

$$R = \frac{\Gamma(N^{***} - p + \eta)}{\Gamma(N^{***} - p + \pi^0)} = \frac{f(q_{\eta})}{f(q_{\pi})} \frac{\gamma_{\eta NN^{***}}}{\gamma_{\pi^0 NN^{***}}},$$
 (2)

where the f's are kinematic factors. The ratio R can, of course, be determined by comparing the height of the N^{***} bump in the cross sections for $\gamma + p \rightarrow p + \eta$ and $\gamma + p \rightarrow p + \pi^0$. The kinematic factor, which is taken to be

$$f(q) = q \left(\frac{q^2}{q^2 + X^2}\right)^L,$$
 (3)

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