#### The Emission of Negative Electrons from Lithium and Fluorine Bombarded with Deuterons

In the last issue of *The Physical Review* we published the energy spectrum of beta-rays from boron bombarded with deuterons, which we were able to obtain by bombarding a target inside the cloud chamber. Continuing with the same experimental set-up we have obtained the beta-ray spectra from two other substances, which we suppose are  $Li^8$  and  $F^{20}$ , and which also have short halflives. In both cases an automatic switching device was used which shut off the bombarding ion beam before the chamber expansion, so that only delayed effects were observed.

### LITHIUM BOMBARDED WITH DEUTERONS

The points in Fig. 1 show the distribution in energy of 1646 electrons obtained from 1000 cloud chamber photographs, taken about  $\frac{1}{2}$  second after bombardment, with  $\frac{1}{2}$  microampere deuteron current at 0.8 M.E.V. The tracks were curved in a magnetic field of 1500 gauss to determine their energy. 0.3 M.E.V. has been added to all points to compensate for the stopping power of the foil surrounding the target, so that they give directly the energy of the



FIG. 1. Energy distribution of negative electrons emitted from a lithium chloride target after bombardment with deuterons.



FIG. 2. Energy distribution of negative electrons emitted from a calcium fluoride target after bombardment with deuterons.

electrons. The reaction is probably

### $Li^7 + H^2 \rightarrow Li^8 + H^1 \rightarrow Be^8 + e^- + H^1 \rightarrow He^4 + He^4 + e^- + H^1.$

If the first part of the reaction is correct, the second disintegration, in which  $Be^8$  splits into two alpha-particles probably follows, since several pieces of information indicate that  $Be^8$  is unstable. There is no assurance, however, that the two disintegrations do not take place in the reverse order, that is,

$$Li^7 + H^2 \rightarrow Li^8 + H^1 \rightarrow He^4 + H^4 + H^1 \rightarrow He^4 + He^4 + e^- + H^1.$$

The protons emitted during bombardment may be expected to have an energy in the neighborhood of 4 M.E.V.

To determine the half-life of the active constituent, we adjusted the timing device so that the ion beam was shut off at  $\frac{1}{4}$ ,  $\frac{1}{2}$ ,  $\frac{3}{4}$  and 1 second before the chamber expansion. 50 photographs were taken at each of these settings, and the average numbers of tracks per photograph were found to be 7.08, 4.84, 3.70 and 2.45, respectively. These, plotted on a log scale lie quite closely on a straight line, and indicate a half-life of  $0.5 \pm 0.1$  second.

#### FLUORINE BOMBARDED WITH DEUTERONS

The points in Fig. 2 show the energy distribution of 1363 electron tracks obtained from 800 cloud chamber photographs taken with  $\frac{1}{2}$  microampere deuteron current, at 0.8 M.E.V., and 1500 gauss magnetic field. These points have also been corrected for the stopping power of the foil. The electrons probably arise from the reaction

$$F^{19}+H^2 \rightarrow F^{20}+H^1 \rightarrow Ne^{20}+e^-+H^1$$

We have measured the half-life of the radio-fluorine by means of an ionization chamber and found it to be  $12\pm 2$  seconds.

#### COMPARISON WITH THEORY

The energy distributions of the electrons in Figs. 1 and 2 do not fit at all with those derived from Fermi's theory in its original form, according to which we should expect the maxima of the curves, for such high energy spectra, to lie at points only slightly less than half the upper limit of energy. It should be kept in mind, however, that the experimentally obtained energy distributions are somewhat influenced by instrumental conditions, which tend, in general, to make unreliable the relative numbers of low energy tracks measured, but we feel that this distortion is important only below 1 or 1.5 M.E.V. It is not probable, therefore, that this can account for the large difference between the observed positions of the maxima and those predicted by Fermi's theory. Professor Oppenheimer has pointed out to us that the data are in very much better agreement with a formula of the form:

# $N(E) = \text{const.} E^2(E_0 - E)^4.$

Such a formula would follow if the coupling energy between the neutron-proton and the electron-neutrino field were taken proportional to the first derivative of the neutrino function, and this possibility has been advanced by Uhlenbeck and Konopinski on empirical grounds. The curves in Figs. 1 and 2 have been drawn according to this formula, fitting the data as well as possible, which necessitated taking as upper limits  $E_0$ , 10.5 and 5.25 M.E.V., respectively. It is seen that these fit the experimental points well within the probable statistical fluctuations, although, from the standpoint of the data alone, one might be inclined to place the upper limits slightly lower than 10.5 and 5.25 in consideration of the possibility of a little straggling. The lifetimes of these substances, on the other hand, lie too near those to be expected on the basis of Fermi's original formulation to permit the inclusion of higher derivatives in the interaction energy.

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## Interaction of Nuclear Particles

Heisenberg,1 in an important set of papers, attributed nuclear binding entirely to proton-neutron attractions. Supposing these to be of a classical nature it is easy to show (if the proton-neutron potential energy is a simple monotonic function of the inter-particle distance) that the nuclear binding energy is

$$W \sim -Z(A - Z). \tag{1}$$

This is essentially what is obtained if the Coulomb repulsion between protons is neglected. Inclusion of Coulomb forces would, however, still give a binding energy varying quadratically with the number of nuclear particles. If this were right we could not understand natural radioactivity and the ending of the periodic system of elements. The failure of a classical interaction law is not apparent for the lightest nuclei and Wigner<sup>2</sup> was able to show that the mass defects of the deuteron and the  $\alpha$ -particle were compatible with classical forces. From the experimental law that mass defects are roughly proportional to the number of nuclear particles we see that our theory must give instead of Eq. (1)

$$W \sim -A$$
. (2)

Now Z(A-Z) is just the number of interactions of Z protons with A - Z neutrons. Heisenberg, and later Majorana,3 saw that the simplest way of obtaining something like Eq. (2) instead of Eq. (1) was the introduction of interactions of such a nature that a given proton or neutron feels only those neutrons or protons in the same quantum state. They introduced such an interaction and obtained instead of Eq. (2)

$$W \sim -Z.$$
 (3)

Heisenberg's proposed interaction gave saturation with one-proton-one-neutron, but the proposal of Majorana allowed two-protons-two-neutrons to interact in an effective way. The latter is much more satisfactory.

These were steps in the right direction but, for heavy nuclei, Eqs. (2) and (3) are not equivalent and experiment favors Eq. (2). To improve the theory we must depart from the original Heisenberg hypothesis that only neutronproton bonds are important. With neutron-neutron and proton-proton bonds of the type suggested by Majorana for the proton-neutron bond, it is possible to develop a theory giving binding energies  $\sim -A$  and nuclear radii  $\sim A^{\frac{1}{3}}$ . These bonds are taken to be the same without regard to the kind of interacting particles. To make this statement more definite we will give, for comparison, expressions for the nuclear potential energy in the theories discussed in this letter.

(Classical)	$V=-\int\rho_{\pi}(\mathbf{r}_{1})I(r_{12})\rho_{\nu}(\mathbf{r}_{2})d\tau_{1}d\tau_{2},$
(Majorana)	$V = -\int \rho_{\pi}(\mathbf{r}_{1},  \mathbf{r}_{2}) I(r_{12}) \rho_{\nu}(\mathbf{r}_{2},  \mathbf{r}_{1}) d\tau_{1} d\tau_{2},$
(Proposed)	$V = -\int \rho(\mathbf{r}_{1},  \mathbf{r}_{2}) I(r_{12}) \rho(\mathbf{r}_{2},  \mathbf{r}_{1}) d\tau_{1} d\tau_{2} + S.$

In these expressions  $\rho_{\pi}$  and  $\rho_{\nu}$  are the density functions for protons and neutrons, respectively. Where one argument is written, these are just the usual functions but if two arguments are indicated the density matrices of Dirac are to be used. In the last equation  $\rho = \rho_{\pi} + \rho_{\nu}$  and S is such a function that this expression does not include the interaction of any particle with itself.

The assumption that I(r) has the form of a simple potential hole of depth  $\alpha mc^2$  and radius  $\beta \epsilon^2/mc^2$  has been made. To get agreement with experimental binding energies of nuclei it was found that the parameters  $\alpha$  and  $\beta$  must be

$$\alpha \sim 40, \qquad \beta \sim 1.00.$$

The calculations on which these results rest are not accurate enough to allow one to decide whether or not all the bonds are of the same strength. It is probable that they are not, so the above numbers must be a sort of average.

Independent arguments for the existence of strong neutron-neutron and proton-proton binding can be found in the number theory properties of stable nuclei. For Zodd and A > 2Z the number of neutrons A - Z is always even. The most simple and unforced explanation of this seems to be that strong attractive interactions exist between paired neutrons. In order that the balance between protons and neutrons, within the nucleus, be preserved, similar forces must exist between paired protons.

White4 has just published preliminary results on protonproton scattering which indicate just such large deviations from the Coulomb law at distances of the order  $10^{-13}$  cm. It is to be hoped that such experiments and furthermore accurate theoretical work will teach us a great deal about these fundamental interactions.

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<sup>1</sup> W. Heisenberg, Zeits. f. Physik 77, 1 (1932); 78, 156 (1933); 80, <sup>1</sup> W. REISCHOULS, 2011
<sup>2</sup> E. Wigner, Phys. Rev. 43, 252 (1933).
<sup>3</sup> E. Majorana, Zeits. f. Physik 82, 137 (1933).
<sup>4</sup> M. G. White, Phys. Rev. 47, 573 (1935).