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## SENSITIVITY OF PROTON POLARIZATION IN d(e,e'p)n REACTION 3.18 TO NEUTRON ELECTRIC FORM FACTOR

## M.P. Rekalo, G.I. Gakh, A.P. Rekalo

## Institute of Physics and Technology, the Ukrainian Academy of Sciences, Kharkov 310108, USSR

We carried out the theoretical analysis of the proton polarization from the d(e,e'p)n reaction in order to investigate its sensitivity to the neutron electric form factor  $\mathtt{G}_{\mathrm{En}}$  . In the general case, all components of the proton polarization vector  $\overrightarrow{P}$  are nonzero. They are defined by five structure functions 1)

$$\begin{aligned} & GP_{x} = \lambda_{e} N \left[ \sqrt{1 - x^{2}} P_{x}^{(0)} + \cos \varphi \sqrt{2x(1 - x)} P_{x}^{(1)} \right], \\ & GP_{y} = \lambda_{e} N \sin \varphi \sqrt{2x(1 - x)} P_{y}^{(1)}, \\ & GP_{z} = \lambda_{e} N \left[ \sqrt{1 - x^{2}} P_{z}^{(0)} + \cos \varphi \sqrt{2x(1 - x)} P_{z}^{(1)} \right], \end{aligned}$$
(1)

where  $\bigcirc$  is the differential cross section for the d(e,e'p)n process with unpolarized particles,  $\lambda_e = \pm 1$  depending on the mutual orientation of the initial electron spin and momentum,  $\mathcal{N}$  is the kinematical factor,  $\mathcal Z$  is the parameter describing the virtual photon linear polarization, arphi is the angle between the electron scattering plane and the  $\int_{c}^{*} d \rightarrow np$  reaction plane (xz),  $\mathcal{D}_{c}^{(0)}$  and  $\mathcal{D}_{c}^{(1)}$  are the structure functions, the z axis is directed along the virtual photon 3-momentum.

The structure functions  $P_i^{(o,1)}(\theta_{\rho}, E_{n\rho}, \kappa^2)$  are calculated here in the framework of the relativistic impulse approximation (Fig.1)(  $\partial_{
ho}$  is the proton scattering angle in the  $\chi^* d \rightarrow np$  reaction c.m.s.,  $E_{np}$  is the n-p c.m.s. relative energy). The dnp-vertex form factors are taken from 2) using the Paris deuteron wave function 3). The dipole fit and the scaling law relations are used for the proton electromagnetic form factors and for the neutron magnetic form factor.

The sensitivity of some components of the  $\overline{\mathcal{P}}$  to the  $\mathtt{G}_{\mathrm{En}}$  is demonstrated by using the following three parametrizations 4 for  $G_{En}$ :

(1)-  $G_{En} = M_n \mathcal{C} G_{Ep} / (1 - 5.6 \mathcal{C})$ , (2)-  $G_{En} = 0$  and (3)-  $G_{En} = \mathcal{C} G_{Mn}$ , where  $M_n = -1.913$ ,  $\mathcal{C} = K^2 / 4 m^2$ , m is the nucleon mass. It is seen from Fig.2 that at  $\Theta_p < 90^\circ$  the  $R_{X,Z}^{(1)} = P_{X,Z}^{(1)} / (H_{11} + H_{22})$  weakly depend on  $\Theta_p$  ( $H_{11}$ ,  $H_{22}$  are the transverse structure functions for the d(e,e'p)n cross section in the unpolarized case). However, in the region of  $\partial_{
ho}pprox$  90° (where the interference between the neutron

and proton diagrams is essential) the sign of the  $R_{\chi}^{(1)}$  changes. It is natural that at  $\partial_{\rho}$  > 90° the main contribution must be from the neutron diagram. Low  $R_{\chi}^{(1)}$  polarization values near  $\Theta_{\rho} \approx 180^{\circ}$  are determined by the smallness of the  $G_{Mn}G_{En}$  product. The high energy neutron polarization (or the low energy proton polarization in the lab.system) appears to be sensitive to the  $G_{En}$  value just at  $\theta_{\rho} \approx 180^{\circ}$ . The angle  $\theta_{\rho}^{\prime o}$  at which the sign of the  $R_{\chi}^{(1)}$  changes depends on the  $G_{En}$  form factor parametrization. This may, in principle, be used for determining  $G_{En}$ .

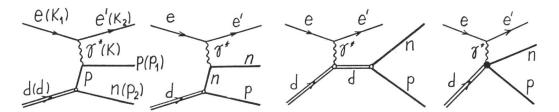


Fig.1. Feynman diagrams

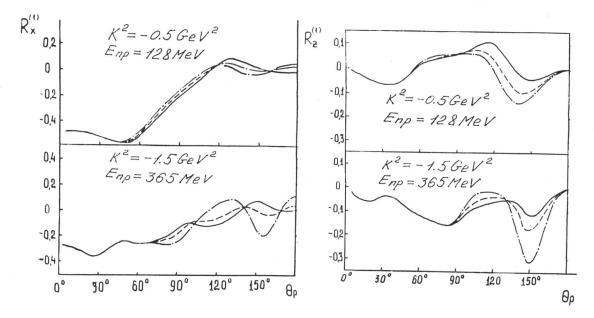


Fig.2. The dependence of the  $R_{\chi}^{(1)}$ ,  $R_{Z}^{(1)}$  on the choice of  $G_{En}$  parametrization. The dashed, solid and dash-dotted curves correspond to parametrizations 1, 2, 3 (see the text) of  $G_{En}$ , respectively. References

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