## Analytical forms of deuteron wave function for potentials Nijmegen group and density distribution

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In the short range the structure of a deuteron is visually described by means of density distribution  $\rho_d^{M_d}(r',\theta)$  [1] for a projections  $M_d=0;\pm 1$  of full angular momentum

$$\begin{cases} \rho_d^0 = \frac{4}{\pi} \left[ C_0(2r') - 2C_2(2r')P_2(\cos\theta) \right]; \\ \rho_d^{\pm 1} = \frac{4}{\pi} \left[ C_0(2r') + C_2(2r')P_2(\cos\theta) \right]; \end{cases}$$

where  $C_0 = R_0^2 + R_2^2$ ;  $C_2 = \sqrt{2}R_0R_2 - \frac{1}{2}R_2^2$  - components of density distribution;  $R_0 = u/r$ ;  $R_2 = w/r$  - the radial functions for S- and D- states;  $P_2$  - Legendre's polynom; r' - distance from the center of masses;  $\theta$  - polar angle to r'; r = 2r' - between partial distance.

Values of density distribution  $\rho_d^{M_d}$  and transition density  $\rho_{tr}^{\pm 1}$  [1] are calculated using the earlier obtained coefficients of DWFs [2, 3] analytical forms in coordinate representation for a nucleon-nucleon potentials Nijmegen group (Nijm1, Nijm2, Nijm93).

Results of similar calculations of density distribution and transition density for Argonne v18 potentials are quoted in paper [4]. Depending on a choice of approximation for DWF the calculated values  $\rho_d^{M_d}$  and  $\rho_{tr}^{\pm 1}$  differ only in the area at 0-0.3 fm. In fact, this indicates, which of the approximations applied is the "best" for beginning of coordinates, despite the absence of redundant knots of the radial DWF.

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