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Comparison between Dirac Equation and Its Equivalent Schrödinger Equation for Inelastic Scattering

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Dirac equation has been used since some years to study proton elastic scattering at intermediate energies¹⁾ (above 200MeV). This equation is :

$$\begin{bmatrix} \vec{\alpha}, \vec{p} + \beta \left\{ m + U_{s}(r) \right\} + U_{v}(\vec{r}) + V_{c}(\vec{r}) \end{bmatrix} \Psi(\vec{r}) = E\Psi(\vec{r})$$
(1)

where V (r) is the coulomb potential, U (r) and U (r), two complex potentials of which the real and the imaginary parts are approximated by a Woods-Saxon form-factor. For a given target and a given laboratory energy, the center of mass momentum k is fixed; we choose the reduced mass for m to define E with $E^2 = m^2 + k^2$. Using the decomposition into large and small components $\Psi(\vec{r}) = [F(\vec{r}), iG(\vec{r})]$, we get a system of two linear equations. The small components $G(\vec{r})$ can be eliminated. Using $F(\vec{r}) = D(\vec{r})^{1/2}f(\vec{r})$ where $D(\vec{r}) = E+m+U_s(\vec{r})-U_v(\vec{r})-V_c(\vec{r})$, we get the Schrödinger equation

$$\begin{cases} \Delta - v_1(\vec{r}) - i\vec{o}\vec{\nabla}v_2(\vec{r})_{\mathbf{A}}\vec{\nabla} - k^2 \end{cases} \mathbf{f}(\vec{r}) = 0 \tag{2}$$

$$v_1(\vec{r}) = \frac{3}{4} \left(\frac{\vec{\nabla}D}{D}\right)^2 - \frac{\Delta D}{2D} + \frac{E}{m} \left\{ U_v(\vec{r}) + v_c(\vec{r}) \right\} + U_s(r) + \frac{1}{2m} \left\{ U_s^2(\vec{r}) - (U_v(\vec{r}) + v_c(\vec{r}))^2 \right\}$$

$$v_c(\vec{r}) = Ln\{D(\vec{r})\} \tag{3}$$

If we consider the potentials $U_{\vec{r}}(\vec{r})$ and $U_{\vec{v}}(\vec{r})$ as describing the target and its excited states (vibrational or rotational models) and the wave functions $\Psi(\vec{r})$ as describing the incident proton and the target, we can derive a set of coupled equations from eq.(1). Their solution is described in another communication for a more general equation. Furthermore, if we neglect the variation of E with the channels in $D(\vec{r})$ and the mixing of the channels due to the factor $D(\vec{r})^{1/2}$ in the definition of $f(\vec{r})$, we get from eq.(2) a set of coupled equations which differs from the conventional ones only by the radial dependence of the potentials.



Fig. 1 : Elastic cross-section



Fig. 2 : Elastic polarization



Fig. 3 : Inelastic analysing power

We have done calculations for the inelastic scattering of 497.5MeV protons on ⁴⁰Ca to the 3⁻ excited state at 3.7MeV, using the first set of parameters published in the analysis of elastic scattering by Clark et al². Results obtained by using eq.(1) are quite the same as those obtained by using eq.(2). Effects of coupled channels are very small and DWBA³) is enough for the inelastic scattering. However higher order effects show up in the elastic scattering as can be seen on Fig. 1 and 2. To show these effects without interference with other problems, we present here results obtained without charge (neutron scattering) between 30° and 60°. Below 30°, difference are too weak, above 60°, cross sections are too small. The conclusion is that eq.(2) slightly overestimates the coupling(the total cross-section for the 3⁻ is 2.527 mb with eq.(1) and 2.617 with eq.(2)). The largest difference at "experimental" angles is found in the elastic polarization. For the inelastic cross-section, the difference could not be seen on a figure. These two computations took almost the same time.

This comparison with a potential found in the literature does not prove that results are always similar. Indeed, in the study of inelastic scattering of 225MeV protons on 9 Be, we obtained a good fit to the analysing power with eq.(1) with a potential which gives worse results with eq.(2). We have still to see if slightly different parameters would give a good fit with eq.(2) instead of eq.(1).

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