

ON THE CALCULATION OF D_A IN THE PHASE-SHIFT APPROXIMATION

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The accuracy of the phase-shift approximation in calculating the well-depth of a A -particle in nuclear matter, D_A , is investigated. A model case of a simple A - N potential is considered. The value of D_A calculated in the phase-shift approximation is about 10 MeV higher than the value obtained by the complete K -matrix method. This indicates that the phase-shift approximation is too rough for application in the D_A problem.

1. Introduction

The well-depth D_A of a A -particle in nuclear matter is of importance and interest because it gives some informations about the A -nucleon interaction.

The most accurate method of calculating D_A is probably the K -matrix method based on the Brueckner theory [1], [2]. Recently, an approximation to the K -matrix method, namely, the phase-shift approximation (PSA) has been applied by Bhaduri and Law in calculating D_A [3]. However, in the case of pure nuclear matter the phase-shift approximation is known to fail [4]. The purpose of the present paper is to investigate the applicability of the PSA in the D_A problem.

To check the accuracy of the PSA we consider a model case, in which we assume the AN interaction to be represented by one of the simple spin-independent potentials considered by Downs and Ware [5]:

$$V_{AN} = \begin{cases} \infty & \text{for } r \leq r_c \\ -V_0 \exp[-3.5412(r-r_c)/b] & \text{for } r > r_c \end{cases} \quad (1)$$

with $r_c = 0.4$ fm, $b = 1.1$ fm, $V_0 = 330.9$ MeV.

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In this case we calculate D_A in the PSA. On the other hand, the "exact" result for D_A , *i. e.*, the result obtained with the K -matrix method, is known for V_{AN} , Eq. (1) [1].

The results obtained indicate a large difference (of about 25%) between the PSA and the "exact", result for D_A . This seems to indicate that the PSA is too rough to be applied in the D_A problem. The calculational method is presented in Section 2, and it follows in essence the procedure applied by Bhaduri and Law [3]. The results obtained are presented and discussed in Section 3.

2. Calculation of D_A in the PSA

In nuclear matter, according to the present state of the Brueckner theory, the potential energy of a A -particle is

$$-D_A = \frac{4}{(2\pi)^3} \frac{1}{\beta^3} \int_0^{\beta k_F} \langle \bar{\mathbf{k}} | t^N | \bar{\mathbf{k}} \rangle d_3 k, \quad (2)$$

where $\beta = m_A / (m_A + m_N)$ (m_A and m_N are the masses of the A -particle and the nucleon), k_F is the Fermi momentum of the nucleons, \mathbf{k} is the relative momentum, and t^N is the AN reaction matrix inside nuclear matter.

Expressing t^N in terms of t^F (the reaction matrix for free AN scattering) one gets:

$$\begin{aligned} t^N &= t^F + t^F \left(\frac{P}{e_0} - \frac{Q'}{e_N} \right) t^N \\ &= t^F + t^F \left(\frac{P}{e_0} - \frac{Q'}{e_N} \right) t^F + \text{higher order terms,} \end{aligned} \quad (3)$$

where P is the principal value operator, Q' is the exclusion principle operator, e_0 and e_N are the energy denominators, which will be specified later. Combining Eqs (2) and (3) one gets $D_A^{(1)}$, the first-order term of D_A in the PSA,

$$D_A^{(1)} = - \frac{4}{(2\pi)^3} \frac{1}{\beta^3} \int_0^{\beta k_F} \langle \mathbf{k} | t^F | \mathbf{k} \rangle d_3 k \quad (4)$$

and the second-order term,

$$D_A^{(2)} = - \frac{4}{(2\pi)^3} \frac{1}{\beta^3} \int_0^{\beta k_F} \left\langle \mathbf{k} \left| t^F \left(\frac{P}{e_0} - \frac{Q'}{e_N} \right) t^F \right| \mathbf{k} \right\rangle d_3 k. \quad (5)$$

To evaluate $D_A^{(1)}$ one has to know the diagonal matrix elements of the reaction matrix t^F . These matrix elements can be expressed in terms of δ_l^i , the l -partial wave phase-shifts generated by the AN potential,

$$\langle \mathbf{k} | t^F | \mathbf{k} \rangle = -4\pi \frac{\hbar^2}{2\mu_{AN}} \frac{1}{k} \sum_l (2l+1) \delta_l^i, \quad (6)$$

where μ_{AN} is the AN reduced mass.

Inserting (6) into Eq. (4) one obtains $D_{A,l}^{(1)}$, the l -partial wave contribution to the first-order term of D_A ,

$$D_{A,l}^{(1)} = \frac{8}{\pi} \frac{1}{\beta^3} \frac{\hbar^2}{2\mu_{AN}} \int_0^{\beta k_F} (2l+1) \delta_l k dk. \quad (7)$$

In our calculation we restrict ourselves to the first three phase-shifts, $l = 0, 1, 2$. In our model case the phase-shifts are obtained by the numerical solution of the Schrödinger equation with V_{AN} , Eq. (1). The values of $\delta_0(k)$, $\delta_1(k)$, $\delta_2(k)$ thus obtained are shown in Fig. 1. Using these values of δ_l and taking $k_F = 1.366 \text{ fm}^{-1}$ we have performed numerically the integrations in Eq. (7). The resulting values of $D_{A,l}^{(1)}$ are presented in Table I.

The calculation of $D_A^{(2)}$ is more complicated. Let us denote by \mathbf{q} the momentum transfer of the interacting lambda and nucleon whose momenta are initially \mathbf{k}_A and \mathbf{k}_1 . Then, the

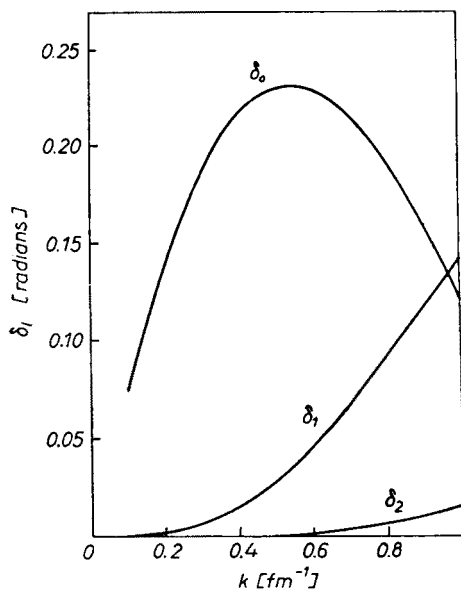


Fig. 1. The phase-shifts (in radians) generated by the V_{AN} potential, Eq. (1), as functions of the λN relative momentum, k

TABLE I

The results of D_A (in MeV) obtained in the phase-shift approximation and the exact results of Ref. [1]

l	First-order term	Second-order term	Exact [1]
0	34.7	-4.4	21.2
1	16.9	—	14.0
2	1.2	—	1.1
Total	52.8	~ -7.1	36.3

energy denominator e_0 is defined by

$$e_0 = \frac{\hbar^2(\mathbf{k}_1 + \mathbf{q})^2}{2m_N} + \frac{\hbar^2(\mathbf{k}_A - \mathbf{q})^2}{2m_A} - \frac{\hbar^2 k_1^2}{2m_N} - \frac{\hbar^2 k_A^2}{2m_A} = \frac{\hbar^2}{2\mu_{AN}} (q^2 + 2\mathbf{k}\mathbf{q}). \quad (8)$$

To determine the energy denominator e_N we assume that the nucleon and the A -particle are moving initially in a potential well U_N and U_A respectively and they are free after scattering, *i. e.*, we assume the single nucleon and lambda potentials to be equal to zero in the intermediate states. Then¹

$$e_N = \frac{\hbar^2}{2\mu_{AN}} (q^2 + 2\mathbf{k}\mathbf{q}) - U_N - U_A. \quad (9)$$

We approximate U_N by a quadratic function

$$U_N = Ak_1^2 + C, \quad (10)$$

with the constants A and C fixed by the two requirements [1]:

$$\bar{\varepsilon}_0 + \frac{1}{2}\bar{U}_N = \frac{1}{2}(\bar{\varepsilon}_0 + \bar{\varepsilon}_N) = \varepsilon_{\text{vol}}, \quad (11)$$

$$\varepsilon_N(k_F) = \varepsilon_{\text{vol}}, \quad (12)$$

where ε_0 denotes the nucleon kinetic energy, $\varepsilon_N = \varepsilon_0 + U_N$, the bars denote average values over the Fermi sea, and $\varepsilon_{\text{vol}} = -15.8$ MeV is the energy per nucleon in nuclear matter. Solving Eqs (11) and (12) we obtain

$$A = 63.407 \text{ MeV fm}^2, \quad C = -113.33 \text{ MeV}.$$

Taking $U_A = -36.3$ MeV (which is the ‘‘exact’’ value obtained in Ref. [1]) we may write Eq. (9) in a modified form

$$e_N = \frac{\hbar^2}{2\mu_{AN}} (q^2 + 2\mathbf{k}\mathbf{q} + \Delta - \nu k^2), \quad (13)$$

with $\Delta = 3.917 \text{ fm}^{-2}$, $\nu = 5.6254$.

Now, the second-order contribution to D_A is given by:

$$D_A^{(2)} = -\frac{4}{(2\pi)^3} \frac{1}{\beta^3} \frac{1}{(2\pi)^3} \int_0^{\beta k_F} d_3 k \int_0^\infty d_3 q |\langle \mathbf{k} | t^F | \mathbf{k} + \mathbf{q} \rangle|^2 \left(\frac{P}{e_0(\mathbf{k}, \mathbf{q})} - \frac{Q'}{e_N(\mathbf{k}, \mathbf{q})} \right). \quad (14)$$

The main problem is to evaluate off-the-momentum-shell matrix elements $\langle \mathbf{k} | t^F | \mathbf{k} + \mathbf{q} \rangle$. Assuming that t^F is local and energy-independent, *i. e.*,

$$\langle \mathbf{k} | t^F | \mathbf{k} + \mathbf{q} \rangle \approx \int_0^\infty t^F(r) e^{i\mathbf{q}\mathbf{r}} d_3 r \equiv t^F(\mathbf{q}), \quad (15)$$

¹ The gap in the single lambda particle spectrum has been neglected in Ref. [3].

and replacing $t^F(q)$ by $t_0^F(q)$ (the s -wave contribution is predominant) one arrives at the following formula:

$$t_0^F(q) = -\frac{16\pi}{q} \frac{\hbar^2}{2\mu_{\Lambda N}} \frac{d}{dq} \left\{ \int_0^\infty dk k \delta_0(k) [\delta(2k-q) + \delta(2k+q)] \right\}. \quad (16)$$

Now, the total ΛN cross-section σ_T may be expressed in terms of δ_l , *i. e.*,

$$\sigma_T = \frac{4\pi}{k^2} \sum_l (2l+1) \sin^2 \delta_l \equiv \sum_l \sigma_l, \quad (17)$$

where σ_l is the l -partial wave cross-section. Hence $t_0^F(q)$ may also be written as a function of σ_0 , the s -wave cross-section,

$$t_0^F(q) = -\frac{2\pi}{q} \frac{\hbar^2}{2\mu_{\Lambda N}} \frac{d}{dq} \left[q^2 \sqrt{\frac{\sigma_0(q/2)}{4\pi}} \right]. \quad (18)$$

Changing the parameter Δ to

$$\delta = \frac{\Delta}{(\beta k_F)^2},$$

and performing part of the integrations in Eq. (14) we obtain:

$$D_\Lambda^{(2)} = -\frac{8}{(2\pi)^4} \frac{2\mu_{\Lambda N}}{\hbar^2} \frac{(\beta k_F)^4}{\beta^3} \int_0^\infty dq |t^F(\beta k_F q)|^2 \eta(q), \quad (19)$$

with

$$\begin{aligned} \eta(q) &= q[f_<(q) - f_>(q)] \text{ for } \beta q \leq 2, \\ &= qf_>(q) \text{ for } \beta q > 2, \end{aligned} \quad (20)$$

where

$$f(q) = \frac{1}{16} \left[(4-q^2) \ln \frac{2+q}{|2-q|} + 4q \right], \quad (21)$$

$$f_<(q) = \frac{1}{2} \int_{1-\beta q}^1 k dk \ln \left[\frac{q^2 + 2kq - \nu k^2 + \delta}{q^2(1-\beta) - \left(\nu + \frac{1}{\beta}\right) k^2 + \left(\delta + \frac{1}{\beta}\right)} \right], \quad (22)$$

$$f_>(q) = \frac{1}{2} \int_0^1 k dk \ln \left[1 + \frac{4kq}{q^2 - 2kq - \nu k^2 + \delta} \right]. \quad (23)$$

The cross-sections σ_0 and σ_T calculated with the phase-shifts obtained previously are shown in Figs 2 and 3. These cross-sections are reproduced by the following forms

$$\sigma_0 = 7.31 \exp(-3.5 k^{1.8}), \quad (24)$$

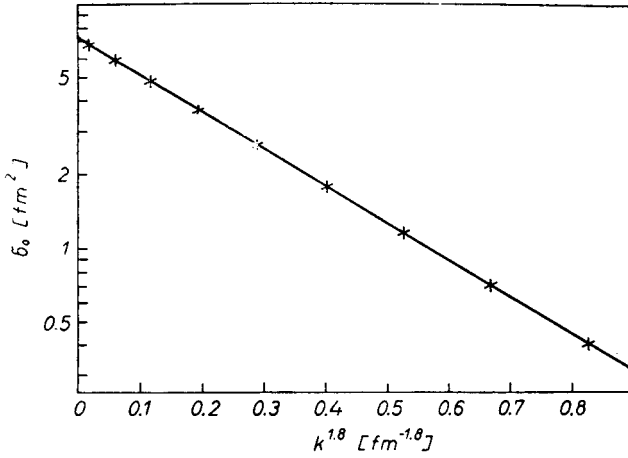


Fig. 2. The s -wave cross-section given by Eq. (24). Crosses represent values calculated with the δ_0 phase-shift

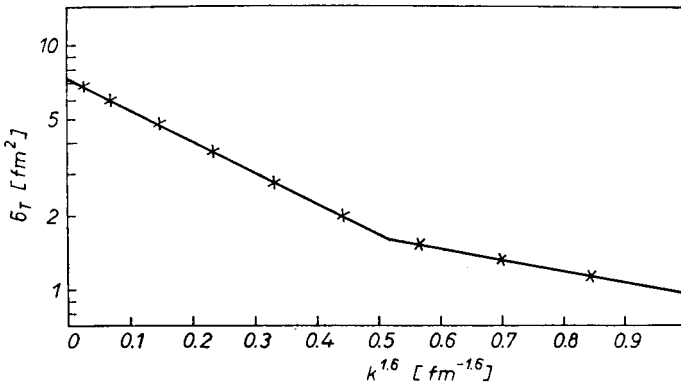


Fig. 3. The total cross-section given by Eq. (25a, b). Crosses represent values calculated with the phase-shifts δ_0 , δ_1 , δ_2 , Eq. (17)

and

$$\sigma_T = \begin{cases} 7.5 \exp(-3k^{1.6}) & \text{for } k \leq 0.66 \text{ fm}^{-1}, \\ 2.67 \exp(-k^{1.6}) & \text{for } k > 0.66 \text{ fm}^{-1}. \end{cases} \quad (25a)$$

$$(25b)$$

Inserting expression (24) into Eq. (18) we have

$$t^F(q) \sim 2\sqrt{7.31} [1 - 0.2643 q^{1.8}] \exp(-0.2937 q^{1.8}). \quad (26)$$

Inserting (25a) into Eq. (18) we get

$$t^F(q) \sim 2\sqrt{7.5} [1 - 0.2455 q^{1.6}] \exp(-0.3069 q^{1.6}), \quad (27a)$$

and inserting (25b) into Eq. (18) we obtain

$$t^F(q) \sim 2\sqrt{2.67} [1 - 0.0818 q^{1.6}] \exp(-0.1023 q^{1.6}). \quad (27b)$$

Introducing expressions (26) and (27a, b) into Eq. (19) and performing the q -integration numerically we find $D_{A,0}^{(2)}$ (the s -wave contribution to the second-order term of D_A) and $D_{A,T}^{(2)}$ (obtained by replacing σ_0 by σ_T in Eq. (18)).

3. Discussion

The results of our calculation are shown in Table I, which also contains the “exact” results of Ref. [1]. Our total first-order contribution, $D_A^{(1)} = 52.8$ MeV, *i. e.* it is 16.5 MeV higher than the “exact” result, $D_A = 36.3$ MeV.

Strictly speaking, the second-order contribution, $D_A^{(2)}$, goes beyond the PSA. Only by introducing drastic approximations (locality and energy independence of t^F) is it possible to express the second-order contribution through the AN phase-shifts. Actually, Bhaduri and Law [3] express the second-order contribution through the AN total cross-section and obtain in this way, in our notation, $D_{A,T}^{(2)}$. In the spirit of the approach of Ref. [3], the value of $D_{A,T}^{(2)} = -7.1$ MeV is supposed to approximate the second-order contribution. This reduces the PSA value of the A potential depth from 52.8 MeV to 45.7 MeV, which is still higher than the “exact” value of D_A by 9.4 MeV. If, on the other hand, we assume that $D_A^{(2)}$ may be approximated by $D_{A,0}^{(2)} = -4.4$ MeV we are lead to a difference of 12.1 MeV between the “exact” and approximate value of D_A . In any case we see that the PSA procedure of Bhaduri and Law [3] applied to our model case leads to a value of D_A which is about 10 MeV higher than the “exact” result. Now, the difference between the D_A values calculated with the contemporary, phenomenological AN potentials and the D_A values estimated empirically is probably less than 10 MeV [1]. For this reason it seems to us that the PSA is too rough to be applied in the D_A problem. On the other hand, as seen from Table I, the PSA may be useful for estimating the higher partial wave contribution to D_A .

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