

SYSTEMATIC STUDY ON THE ELASTIC SCATTERING OF 65 MEV POLARIZED PROTONS

By

Harutaka SAKAGUCHI

Department of Physics, Kyoto University, Kyoto 606, Japan

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ABSTRACT

Elastic scattering of 65 MeV polarized protons from 25 nuclei (^{16}O - ^{209}Bi) has been measured. Systematics of the optical potential, which reproduces both the analyzing power and the differential cross section data remarkably, was pursued. The volume integral of the real central part of the optical potential (J_R) shows a behavior similar to the binding energy curve for the target mass number. The mean square radius of the real central part of the optical potential is found to obey the relation $\langle r^2 \rangle_{\text{pot}} = (0.937 \pm 0.012) A^{2/3} + (6.42 \pm 0.21) \text{ fm}^2$. By comparing with the systematics of the charge distributions obtained from electron scattering data, it is found that the effective two-body interaction range between an incident proton and a nucleon in the target has a target mass number dependence given by $\langle r^2 \rangle_{\text{int}} = (0.132 \pm 0.013) A^{2/3} + (4.24 \pm 0.24) \text{ fm}^2$. Assuming this relation, root mean square radii of the point nucleon distributions are obtained. The dependences of the J_R -value and the $\langle r^2 \rangle_{\text{pot}}$ -value on the mass number and energy obtained here are compared critically with recent microscopic optical potential calculations.

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1. Introduction

Recent progress in nuclear matter theories has made it possible to understand the nuclear optical potential microscopically in terms of a two-body nucleon-nucleon interaction. Jeukenne, Lejeune and Mahaux (JLM)^{1)~4)} at Liege and Brieva and Rook^{5)~9)} at Oxford have calculated the nuclear optical potential microscopically. In order to check critically these global optical potential theories and to extract new aspects in many body problems, it is necessary to measure accurately proton elastic scattering over a wide range of target nuclei and over a wide range of energies relative to the Fermi energy. In applying nuclear matter theory to scattering problems there are many difficulties to overcome by using suitable approximations. The most ambiguous process among them is the transformation procedure from infinite nuclear matter to finite nuclei such as the process using the local density approximation (LDA). As was already pointed out by Wong¹⁰⁾ and Negele,¹¹⁾ the LDA is not accurate in the nuclear surface region. On the other hand, the nuclear surface is the region most sensitively explored by nuclear scattering, and surface effects will be exemplified by their mass number dependence. In order to clarify the role of the nuclear surface and to check the approximation used in the theory, it is important to determine accurately the mass number dependence of the optical potential. For the LDA, energy dependence of the t-matrix in nuclear matter is reflected directly in the potential depth. But in the folding potential of Brieva and Rook the situation is not so straightforward as in the LDA. For checking the validity of their approximation, the energy dependence of the optical potential will be a useful guide.

On the experimental side, the study of the optical potential has progressed after Becchetti and Greenless's¹²⁾ work. If we plot the data available on elastic scattering in a two dimensional plane of incident proton energy versus target mass number, we notice that data are concentrated at energies below 35 MeV and for targets near magic nuclei. In addition, systematic experimental studies on the optical potential using a polarized proton beam are still scarce. Recently Fabrici et al. reported^{13)~14)} measurements of the elastic scattering at several proton energies between 20 MeV and 45 MeV. Although their systematic analysis utilized polarization data only partially, it has clarified the energy dependence in that region. In the intermediate energy region a group at Indiana Univ. is investigating^{15)~17)} the systematics of the optical potential in a relativistic framework. In this paper we report systematic measurements on elastic scattering of polarized protons and an analysis of data over a wide range of target mass numbers, from ¹⁶O to ²⁰⁹Bi, at an incident proton energy of 65 MeV. Partial results have been published.^{18)~20)} The use of high-purity germanium (HP-Ge) detectors has made possible rapid data acquisition at this energy. The 65 MeV data are thought to be valuable not only because they fill a gap in experimental data but also because of the simplicity of the reaction mechanism; they are relatively free from the giant resonance

effects and the multi-step processes observed in the lower energy region.

2. Experimental method

The polarized proton beam (of 10 keV energy) from the atomic beam type polarized ion source²¹⁾ (PIS) is injected axially into the Research Center for Nuclear Physics (RCNP) Osaka Univ. AVF cyclotron.²²⁾ A beam buncher in the injection system intensifies the beam current by about a factor of 3. The extracted beam of 65 MeV was energy analyzed and transported to the polarization experiment area.²³⁾ As shown in Fig. 1, the beam was first focused on a target to a beam spot size of about $1 \times 2 \text{ mm}^2$. After passing through the target, the polarized beam was again focused on a polarimeter target foil and then collected by a Faraday cup located downstream of the polarimeter. The beam current was monitored by a current digitizer whose output pulses were routed by a spin controller²⁴⁾, depending on the beam polarization direction.

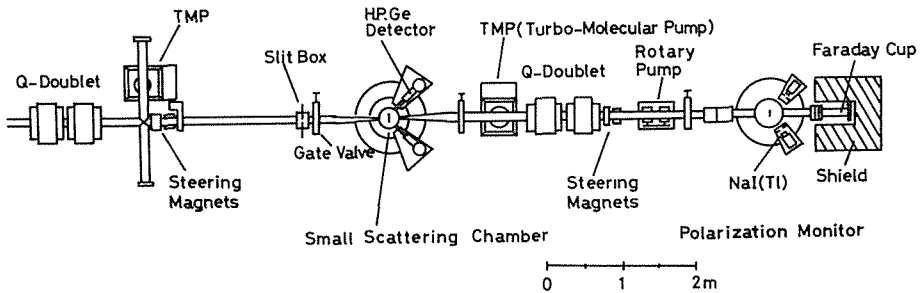


Fig. 1. A layout of the polarization experiment area.

Scattered protons were detected by 1.5 cm thick HP-Ge detectors, which were located at symmetric scattering angle to the left and right of the beam on the goniometers outside the scattering chamber. In the angle region of rapidly changing angular distributions, the detectors were placed at a distance of more than 30 cm from the scattering chamber (of 26.6 cm diameter) to obtain better angular resolution. A vacuum bag was inserted between the scattering chamber and the HP-Ge cryostat to reduce the energy loss and range straggling of the scattered particles in air. For solid targets, the acceptance solid angles of the detector were 0.1453 msr at forward angles ($\theta < 32.5^\circ$) and 0.6902 msr at backward angles ($\theta \geq 32.5^\circ$). For gas targets the vacuum bag was inserted between the front and back slits of a double-slit collimator. The gas target G-factors of the slit system were 1.09×10^{-4} at the forward angles and 2.15×10^{-4} at the backward angles. The beam direction was determined by two methods. One was the conventional kinematical crossover method. The other was to search for the scattering angles on both sides of the beam direction at which the analyzing power of $p + {}^{208}\text{Pb}$

elastic scattering changes sign rapidly. The uncertainty of the scattering angles was estimated to be less than 0.05 degrees.

The degree of beam polarization changes depending on the vacuum in the ionizer region of the PIS, the out-gassing history of the ionizer, and on other ionizer conditions. So the beam polarization was monitored continuously by a polarimeter located downstream of the scattering chamber. Scattered protons from the polarimeter target foil (a stacked 3 mg/cm² thick polyethylene foil) were detected by NaI(Tl) scintillators placed at the $\theta_{\text{LAB}} = 47.5$ degrees. The left-right asymmetry of the elastic scattering from ¹²C was used to deduce the beam polarization. The polarimeter target foil was changed after an appropriate beam irradiation in order to reduce the effect of contaminant buildup. The beam polarization error due to the contaminant peak was estimated to be less than 0.2%. As the analyzing power of the ¹²C polarimeter at $\theta_{\text{LAB}} = 47.5^\circ$, we adopted a value of 0.975 ± 0.011 , which was obtained in the double scattering experiment at $E_p = 65$ MeV by Kato et al.²⁵⁾ The direction of the proton spin was reversed after every 200 pC of integrated beam charge by reversing the solenoid magnetic field direction at the ionizer of the PIS. In the later stage of the experiment the spin direction was reversed every 1 s by alternating the atomic beam rf transition mode between the weak field and strong field transition. Signals from a micro-processor triggered the spin controller which controlled the rf transition mode, the scalers, and the data storage locations in the memory of the pulse height

Table 1. Targets

Nucleus	Form	Thickness (mg/cm ²)	Enrichment
¹⁶ O	gas (O ₂)	1-2 atm	natural (99.8%)
²⁰ Ne	gas (Ne)	1-2 atm	natural (90.51%)
²⁴ Mg	metal foil	3.18	99.94%
²⁸ Si	metal foil	2.98	natural (92.21%)
⁴⁰ Ar	gas (Ar)	1-2 atm	natural (99.60%)
⁴⁰ Ca	metal foil	2.19	natural (96.97%)
⁴⁴ Ca	metal foil	2.78	98.55%
⁴⁸ Ca	CaCO ₃ +mylar	1.34	97.27%
⁴⁶ Ti	metal foil	0.52	81.20%
⁴⁸ Ti	metal foil	0.99	99.25%
⁵⁰ Ti	metal foil	0.82	83.2%
⁵⁴ Fe	metal foil	1.57	96.81%
⁵⁶ Fe	metal foil	1.02	99.93%
⁵⁹ Co	metal foil	2.09	99.83%
⁵⁸ Ni	metal foil	2.04	99.83%
⁶⁰ Ni	metal foil	2.04	99.79%
⁶² Ni	metal foil	1.71	96.48%
⁶⁴ Ni	metal foil	3.55	96.48%
⁸⁹ Y	metal foil	1.376	natural (100%)
⁹⁰ Zr	metal foil	2.672	97.65%
⁹⁸ Mo	metal foil	0.900	97.01%
¹⁰⁰ Mo	metal foil	0.372	97.27%
¹⁴⁴ Sm	metal foil	1.71	96.33%
²⁰⁸ Pb	metal foil	15.1	99.14%
²⁰⁹ Bi	metal foil	4.23	natural (100%)

analyzer, where energy spectra were stored in different memory locations depending on the polarization direction. The average beam intensity during the measurements was about 30 nA and the beam polarization was about 60–70%. The overall energy resolution detected by the HP-Ge system was 180keV–250 keV FWHM, including the beam energy spread and the range straggling due to window foils of the scattering chamber, the vacuum bag, and the HP-Ge cryostat. Table 1 lists the forms, thicknesses and enrichments of the target foils used. The thicknesses of the solid target foils were measured by dividing the total weight by the area, and for gas targets the gas pressure was monitored using a strain gauge sensor.

3. Data reduction

Analyzing powers were measured by the left and right detectors located at the same scattering angle. We denote by $L\uparrow$ the number of particles detected by the left detector in the spin up mode. $L\downarrow$, $R\uparrow$, and $R\downarrow$ are defined in an analogous way. The analyzing power $A_y(\theta)$ and its statistical error $\delta A_y(\theta)$ are calculated as follows

$$\begin{aligned}
 A_y(\theta) &= \frac{1}{P_{\text{Beam}}} \left(\frac{r-1}{r+1} \right), \quad r \equiv \left(\frac{L\uparrow \cdot R\downarrow}{L\downarrow \cdot R\uparrow} \right)^{1/2} \\
 P_{\text{Beam}} &= \frac{1}{A_y(^{12}\text{C})} \left(\frac{r'-1}{r'+1} \right), \quad r' \equiv \left(\frac{L\uparrow_{\text{pol}} \cdot R\downarrow_{\text{pol}}}{L\downarrow_{\text{pol}} \cdot R\uparrow_{\text{pol}}} \right)^{1/2} \\
 \delta A_y(\theta) &= A_y(\theta) \left[\left(\frac{\delta P_{\text{Beam}}}{P_{\text{Beam}}} \right)^2 + \left(\frac{\delta A}{A} \right)^2 \right]^{1/2} \\
 &= A_y(\theta) \left[\left(\frac{\delta P_{\text{Beam}}}{P_{\text{Beam}}} \right)^2 + \left(\frac{r}{r^2-1} \right)^2 \left(\frac{1}{R\uparrow} + \frac{1}{R\downarrow} + \frac{1}{L\uparrow} + \frac{1}{L\downarrow} \right) \right]^{1/2} \\
 \delta P_{\text{Beam}} &= P_{\text{Beam}}(\theta) \left[\left(\frac{\delta A_y(^{12}\text{C})}{A_y(^{12}\text{C})} \right)^2 \right. \\
 &\quad \left. + \left(\frac{r'}{r'^2-1} \right)^2 \left(\frac{1}{R\uparrow_{\text{pol}}} + \frac{1}{R\downarrow_{\text{pol}}} + \frac{1}{L\uparrow_{\text{pol}}} + \frac{1}{L\downarrow_{\text{pol}}} \right) \right]^{1/2} \\
 A_y(^{12}\text{C}) &= 0.975 \quad \delta A_y(^{12}\text{C}) = 0.011
 \end{aligned}$$

where $L\uparrow_{\text{pol}}$ denotes carbon elastic peak sum for the left polarimeter detector in the spin up mode, and $L\downarrow_{\text{pol}}$, $R\uparrow_{\text{pol}}$, and $R\downarrow_{\text{pol}}$ are similarly defined. $A_y(^{12}\text{C})$ and $\delta A_y(^{12}\text{C})$ are the carbon analyzing power and its uncertainty at the scattering angle $\theta_{\text{LAB}} = 47.5^\circ$ and $E_p = 65$ MeV. Differential cross section data were corrected by the detector efficiency due to the nuclear reaction in the HP-Ge itself. Energy dependent detector efficiency $\varepsilon(E_p)$ in the energy region of 45 MeV to 65 MeV is obtained from the Makino's data²⁶⁾ according to the relation

$$\varepsilon(E) = 1.0267 - 0.0011333E$$

Using this formula, the difference of the elastic peak detecting efficiency between $\theta_{\text{LAB}}=0^\circ$ and $\theta_{\text{LAB}}=80^\circ$ was 0.7% for ^{16}O and 0.3% for ^{40}Ca . Thus the correction due to Makino's data affected the relative angular distributions of the differential cross sections negligibly. Measured cross section and analyzing power data are plotted in Fig. 2. Numerical values of the data are listed in the Appendix. Error bars shown in the figure and the Appendix are only the statistical ones. In the analyzing power data, uncertainties of the ^{12}C -polarimeter analyzing power are included. We notice a systematic shift of the diffraction pattern as the target mass number increases. In particular, a sharp rise near 30° in the analysing power data shifts forwards as the target mass number increases.

4. Optical potential fitting

The optical potential fitting to the measured data was performed using the automatic search code MAGALI of Raynal.²⁷⁾

The following optical potential was as used;

$$\begin{aligned}
 U(r) = & V_{\text{Coul}}(r) - V_r f(r; r_R, a_R) - iW_v f(r; r_{vv}, a_{vv}) \\
 & + 4a_{ws} \mathbf{W}_s \cdot \mathbf{i} \frac{d}{dr} f(r; r_{ws}, a_{ws}) \\
 & + V_{ls} \left(\frac{\hbar}{m_\pi c} \right)^2 \left(\frac{1}{r} \frac{d}{dr} f(r; r_{ls}, a_{ls}) \right) (\boldsymbol{\sigma} \cdot \mathbf{L})
 \end{aligned}$$

where

$$\begin{aligned}
 f(r; r_0, a_0) &= (1 + \exp((r - r_0 A^{1/3})/a_0))^{-1} \\
 V_{\text{Coul}}(r) &= \begin{cases} \frac{Ze^2}{2r_c A^{1/3}} \left(3 - \frac{r^2}{r_c^2 A^{2/3}} \right) & r \leq r_c A^{1/3} \\ \frac{Ze^2}{r} & r \geq r_c A^{1/3} \end{cases}
 \end{aligned}$$

A search for best-fit values of the optical potential parameters was started using the gas-target data. The probability of becoming trapped in a false local minimum during the search for a χ^2 -minimum was thought to be small, since uncertainties in the cross section due to uncertainties in the target thickness measurement are small for the gas targets. For ^{40}Ar , we started with the Becchetti and Greelees parameter values. The initial parameter values for ^{16}O , ^{20}Ne were obtained from the best-fit ^{40}Ar parameter set. For other targets, potential parameters of the neighbouring target which had already been fitted were adopted as a starting set. Also, a renormalization of the calculation was introduced because of the target thickness uncertainty. In Fig. 2 measured differential cross sections and analyzing powers are shown together with the best-fit optical potential calculations. The optical potential parameters and the associated χ^2 -values obtained are listed in Table 2.

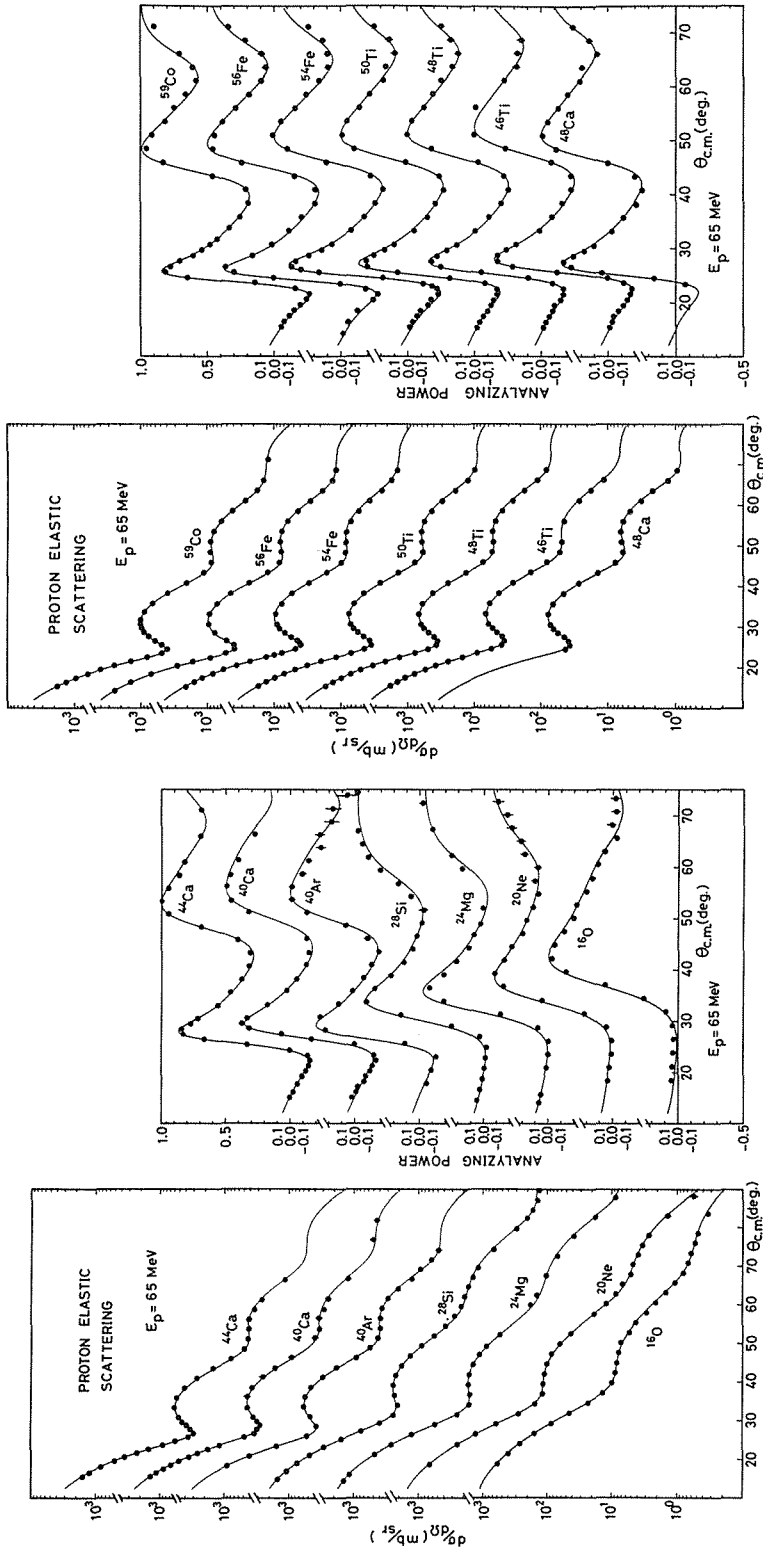


Fig. 2. (b) Measured cross sections and analyzing powers. The solid curves are optical potential calculations. The parameters used are listed in Table 2.

Fig. 2. (a) Measured cross sections and analyzing powers. The solid curves are optical potential calculations. The parameters used are listed in Table 2. In fitting the analyzing power data, the data used were extended to $\theta_{cm}=90^\circ$ for the light nuclei although only the data ($\theta_{cm} \leq 75^\circ$) are shown in the figure.

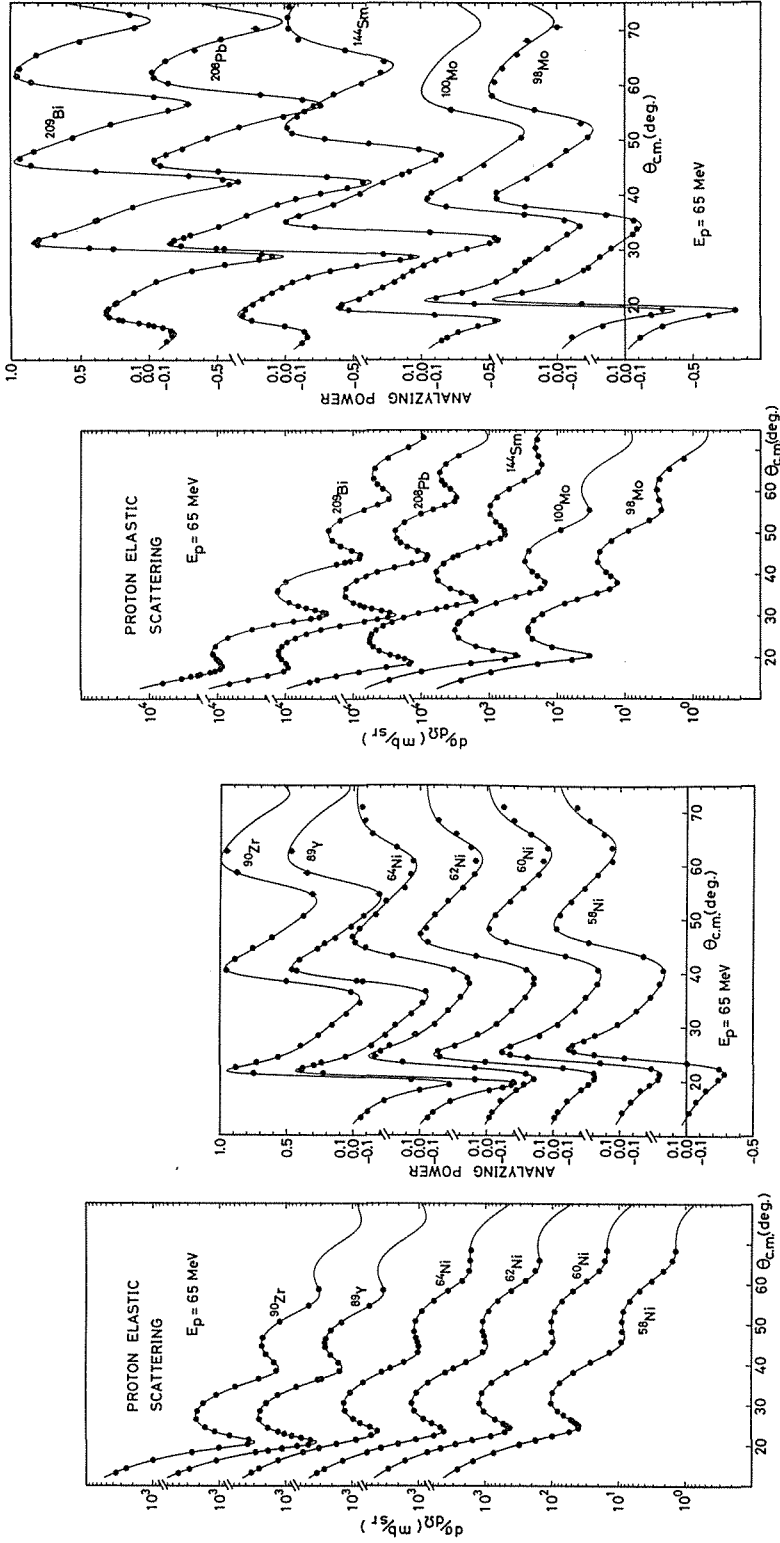


Fig. 2. (d) Measured cross sections and analyzing powers. The solid curves are optical potential calculations. The parameters used are listed in Table 2.

Fig. 2. (c) Measured cross sections and analyzing powers. The solid curves are optical potential calculations. The parameters used are listed in Table 2.

Table 2. Best fit optical potential parameters Coulomb radius is fixed to $r_c = 1.25$ fm.

Nucleus	V_R	r_R	a_R	W_0	r_{wp}	a_{wp}	W_1	r_{w1}	a_{w1}	V_{LS}	r_{LS}	a_{LS}	χ^2/F	Renorm. Factor
¹⁶ O	27.172	1.297	0.6556	12.847	0.2762	1.198	3.290	1.350	0.375	5.793	1.057	0.5807	1.50	1.04
²⁰ Ne	31.162	1.212	0.7439	7.709	0.9552	0.9484	2.658	1.320	0.5367	5.402	1.015	0.6860	1.25	1.05
²⁴ Mg	27.242	1.251	0.6864	10.525	0.8908	1.041	1.809	1.395	0.3850	5.269	1.022	0.6364	2.29	1.23
²⁸ Si	33.912	1.176	0.7248	9.917	0.8388	0.3660	5.348	1.241	0.5656	5.970	1.007	0.6181	1.68	1.05
⁴⁰ Ar	34.249	1.208	0.7266	9.063	1.137	0.7498	2.897	1.357	0.4928	5.606	1.061	0.6722	1.64	1.07
⁴⁸ Ca	33.285	1.232	0.7080	8.131	1.164	0.7993	2.737	1.344	0.4413	5.796	1.088	0.6317	0.806	0.807
⁴⁸ Ca	31.474	1.246	0.7007	9.703	1.120	0.9099	2.553	1.329	0.4476	5.449	1.094	0.6554	1.16	0.956
⁴⁸ Ca	32.907	1.230	0.6808	10.228	1.187	0.7824	2.515	1.344	0.3975	5.680	1.091	0.6475	0.792	1.03
⁴⁶ Ti	31.691	1.241	0.7082	8.034	1.182	0.8090	2.376	1.347	0.4949	5.285	1.094	0.6554	0.772	1.03
⁴⁶ Ti	31.425	1.242	0.7041	8.244	1.191	0.8346	2.189	1.349	0.4199	5.209	1.094	0.6554	0.608	0.955
⁵⁰ Ti	32.610	1.226	0.6898	9.521	1.162	0.8216	2.506	1.346	0.3941	5.248	1.083	0.6276	0.550	0.982
⁵⁴ Fe	33.195	1.212	0.6907	10.410	1.047	0.7670	2.922	1.330	0.4803	5.608	1.071	0.6155	0.550	1.11
⁵⁶ Fe	34.870	1.197	0.7138	11.354	1.004	0.8162	3.333	1.316	0.5307	5.581	1.051	0.6285	0.678	1.01
⁵⁸ Co	33.634	1.212	0.7177	9.606	1.133	0.8240	2.231	1.346	0.4654	5.559	1.071	0.6554	0.610	0.966
⁵⁸ Ni	35.466	1.173	0.7248	12.499	0.9729	0.6262	3.618	1.319	0.5639	5.907	1.027	0.6554	0.511	1.08
⁶⁰ Ni	34.812	1.195	0.7171	11.662	1.001	0.8083	3.282	1.317	0.5345	5.781	1.049	0.6554	0.439	1.11
⁶² Ni	33.831	1.218	0.7175	11.356	1.052	0.8815	3.195	1.320	0.4562	5.757	1.074	0.6554	0.638	1.04
⁶⁴ Ni	33.858	1.209	0.7234	11.518	1.019	0.7744	3.864	1.310	0.5422	5.822	1.063	0.6554	0.696	1.02
⁸⁹ Y	35.463	1.232	0.7084	8.545	1.274	0.7226	2.806	1.301	0.3847	5.506	1.147	0.6300	0.918	1.06
⁹⁰ Zr	34.861	1.233	0.7129	9.230	1.223	0.7070	2.621	1.352	0.4180	5.395	1.138	0.6527	0.498	0.936
⁹² Mo	34.460	1.240	0.7407	8.585	1.016	0.8210	5.163	1.261	0.6046	5.035	1.147	0.6603	0.729	0.951
¹⁰⁰ Mo	34.933	1.244	0.7311	7.058	1.072	0.7878	5.901	1.248	0.6005	5.267	1.156	0.6339	0.283	1.08
¹⁴⁴ Sm	36.162	1.230	0.7288	10.439	1.016	0.8210	5.676	1.248	0.6005	5.397	1.151	0.6224	0.584	1.10
²⁰⁹ Pb	39.105	1.223	0.7461	9.816	1.142	0.7775	5.758	1.253	0.5872	5.838	1.175	0.6103	0.652	1.13
²⁰⁹ Bi	37.399	1.229	0.7347	12.160	1.016	0.8210	6.236	1.252	0.5998	5.447	1.155	0.6330	0.696	1.09

Uncertainties in the experimental data are mainly due to inhomogenities of the target foil thickness and to the scattering angle errors (less than 0.1°) rather than to counting statistics. Therefore the uncertainties used during the parameter search were

$$\delta\left(\frac{d\sigma}{d\Omega}\right) = \text{Max}\left(0.03 \times \left(\frac{d\sigma}{d\Omega}\right), \left(\frac{d\sigma}{d\Omega}\right)_{\text{statistical}}\right)$$

and

$$\delta A(\theta) = \text{Max}(0.03, \delta A(\theta)_{\text{statistical}})$$

in order to avoid trapping in an unphysical local χ^2 -minimum. As for the data points in Fig. 2, the error bars for the cross sections include only the statistical errors, while the error bars for analyzing powers include the uncertainty of the ^{12}C -polarimeter analyzing power in addition.

5. Systematics of the mean square radius of the real central part of the optical potential and the effective interaction range

From the real central part of the optical potential obtained in this analysis, the mean square radius of the potential $\langle r^2 \rangle_{\text{pot}}$, was calculated and is plotted as a function of $A^{2/3}$ (A denotes the target mass number) in Fig. 3. The $\langle r^2 \rangle_{\text{pot}}$ data are remarkably linear in $A^{2/3}$. A linear least-squares fit to the $\langle r^2 \rangle_{\text{pot}}$ data gives

$$\langle r^2 \rangle_{\text{pot}} = (0.937 \pm 0.012) A^{2/3} + 6.42 \pm 0.21 \text{ fm}^2. \quad (1)$$

The error bars in Fig. 3 indicate the uncertainties in the optical potential fitting and were calculated using the following procedure. First, all the

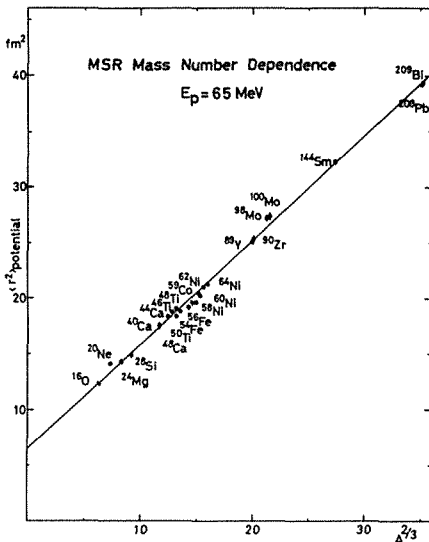


Fig. 3. Mean square radii of the real central part of the optical potentials are shown as a function of the target mass number A . The definition of the error bars in the figure is explained in the text. The numerical data with error bars are listed in Table 3. The solid line is obtained by least-square linear fitting to the data.

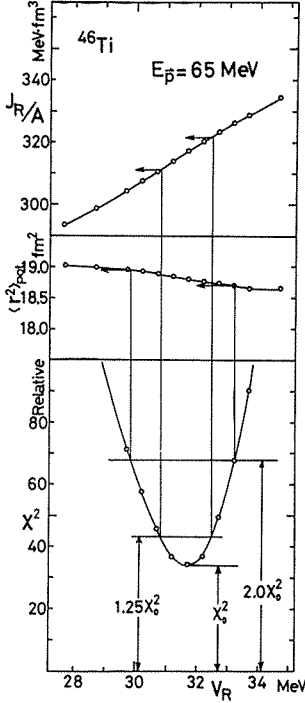


Fig. 4. An example of the procedure for deducing the error bar is shown. The volume integral per nucleon of the real central part of the optical potential, the mean square radius of the potential and the total χ^2 -value are plotted as a function of the potential depth V_R for ^{46}Ti . For $\langle r^2 \rangle_{\text{pot}}$, error bars are obtained from the $2\chi_0^2$ point and for J_R/A from $1.25\chi_0^2$ point.

parameters except V_r and r_c were searched for to obtain the $\chi_{\text{min}}^2(V_r)$ as a function of V_r . A $\chi_{\text{min}}^2(V_r)$ curve for ^{46}Ti case is shown in Fig. 4. The curve resembles a parabola. Then the error of $\langle r^2 \rangle_{\text{pot}}$ was calculated from the $\langle r^2 \rangle_{\text{pot},1}$ and $\langle r^2 \rangle_{\text{pot},2}$ values, which were obtained with the parameter sets at $\chi_{\text{min}}^2(V_r) = 2\chi_0^2$ (best fit). Thus, the uncertainty due to the well known $VR^n = \text{constant}$ parameter correlation is included in the error bars.

The linear relation between $\langle r^2 \rangle_{\text{pot}}$ and $A^{2/3}$ is understood by using a simple folding model. As was shown already by Greenlees, Pyle and Tang,²⁸⁾ the real central part of the optical potential can be written as

$$U(\mathbf{r}_0) = \int \rho(\mathbf{r}) V_{\text{int}}(|\mathbf{r} - \mathbf{r}_0|) d\mathbf{r}^3, \quad (2)$$

where $\rho(\mathbf{r})$ is the density distribution of point nucleons and $V_{\text{int}}(\mathbf{r})$ is an effective two body interaction. For a nucleus with a rotationally symmetric density distribution, the mean square radius of the potential deduced from Eq. (2) is

$$\langle r^2 \rangle_{\text{pot}} = \langle r^2 \rangle_{\text{matt}} + \langle r^2 \rangle_{\text{int}}, \quad (3)$$

where $\langle r^2 \rangle_{\text{pot}} = \int r^2 U(\mathbf{r}) d\mathbf{r}^3 / \int U(\mathbf{r}) d\mathbf{r}^3$, $\langle r^2 \rangle_{\text{matt}} = \int r^2 \rho(\mathbf{r}) d\mathbf{r}^3 / \int \rho(\mathbf{r}) d\mathbf{r}^3$ and $\langle r^2 \rangle_{\text{int}} = \int r^2 V_{\text{int}}(\mathbf{r}) d\mathbf{r}^3 / \int V_{\text{int}}(\mathbf{r}) d\mathbf{r}^3$. If we assume that $U(\mathbf{r})$ and $\rho(\mathbf{r})$ are spherically symmetric Fermi functions, $\langle r^2 \rangle_{\text{matt}}$ can be calculated in a good approximation^{29),30)} as

$$\langle r^2 \rangle_{\text{matt}} = \frac{3}{5} R_m^2 + \frac{7}{5} \pi^2 a_m^2, \quad (4)$$

where R_m and a_m are the half density radius and diffuseness of the point nucleon density distribution, respectively. Therefore $\langle r^2 \rangle_{\text{pot}}$ is expressed as

$$\langle r^2 \rangle_{\text{pot}} = \frac{3}{5} R_m^2 + \frac{7}{5} \pi^2 a_m^2 + \langle r^2 \rangle_{\text{int}}. \quad (5)$$

In order to obtain a relation between the half density radius and the mass number, we use the volume integral of the Fermi-type density distribution:

$$\int \rho(r) d\mathbf{r}^3 = \frac{4\pi}{3} R_m^3 \left(1 + \frac{\pi^2 a_m^2}{R_m^2} \right) \rho_0 = A \quad (6)$$

where

$$\rho(r) = \frac{\rho_0}{1 + \exp\left(\frac{r - R_m}{a_m}\right)}$$

Since a_m and ρ_0 are reasonably constant with A , the half density radius R_m is calculated as a function of A , as

$$R_m = r_m A^{1/3} \left\{ 1 - \frac{1}{3} \left(\frac{\pi^2 a_m^2}{r_m^2 A^{2/3}} \right) + \frac{1}{81} \left(\frac{\pi^2 a_m^2}{r_m^2 A^{2/3}} \right)^3 + \dots \right\} \quad (7)$$

where

$$r_m = \left(\frac{3}{4\pi\rho_0} \right)^{1/3}$$

Inserting (7) into (5), we obtain for $\langle r^2 \rangle_{\text{pot}}$,

$$\langle r^2 \rangle_{\text{pot}} = \langle r^2 \rangle_{\text{int}} + \frac{3}{5} r_m^2 A^{2/3} + \pi^2 a_m^2 + \frac{1}{15} \pi^2 a_m^2 \left(\frac{\pi^2 a_m^2}{r_m^2 A^{2/3}} \right) + \dots \quad (8)$$

The 4th term can be neglected, since the ratio of the 4th term to the 3rd term is less than 0.02 for nuclei considered ($A \geq 16$). We finally obtain an approximate relation

$$\langle r^2 \rangle_{\text{pot}} = \frac{3}{5} r_m^2 A^{2/3} + \pi^2 a_m^2 + \langle r^2 \rangle_{\text{int}}. \quad (9)$$

If we treat $\langle r^2 \rangle_{\text{int}}$ as a constant value as usual, $\langle r^2 \rangle_{\text{pot}}$ is linear in $A^{2/3}$ with the coefficient $\frac{3}{5} r_m^2$. In order to obtain the $\langle r^2 \rangle_{\text{int}}$ -value, we need to know the a_m -value, which is inferred from the charge distribution data. By comparing the two linear relations (1) and (9), we obtain the value of $r_m = (1.25 \pm 0.01)$ fm.

If we assume that the point proton density distribution is also of the Fermi-type (as we did for nuclear matter distribution), then $\langle r^2 \rangle_{\text{charge}}$ can be written in terms of the half-density radius R_p , the diffuseness a_p of the point proton distribution, and the mean square radius of the charge distribution of the proton itself, $\langle r^2 \rangle_{\text{proton}}$, as

$$\langle r^2 \rangle_{\text{charge}} = \frac{3}{5} R_p^2 + \frac{7}{5} \pi^2 a_p^2 + \langle r^2 \rangle_{\text{proton}} . \quad (10)$$

For a relation between R_p and the mass number A , we use the same relation as (7)

$$R_p = r_p A^{1/3} \left\{ 1 - \frac{1}{3} \left(\frac{\pi^2 a_p^2}{r_p^2 A^{2/3}} \right) + \frac{1}{81} \left(\frac{\pi^2 a_p^2}{r_p^2 A^{2/3}} \right)^3 + \dots \right\} \quad (11)$$

We finally obtain a linear relation for $\langle r^2 \rangle_{\text{charge}}$ with $A^{2/3}$.

$$\langle r^2 \rangle_{\text{charge}} = \frac{3}{5} r_p^2 A^{2/3} + \pi^2 a_p^2 + \langle r^2 \rangle_{\text{proton}} \quad (12)$$

The mean square radius of the charge distribution, $\langle r^2 \rangle_{\text{charge}}$, from electron scattering³¹⁾ is plotted in Fig. 5 as a function of $A^{2/3}$. By least-squares linear fitting, we obtained $\langle r^2 \rangle_{\text{charge}} = (0.799 \pm 0.006) A^{2/3} + (2.50 \pm 0.12) \text{ fm}^2$. By introducing the $\langle r^2 \rangle_{\text{proton}} = 0.64 \text{ fm}^2$ and comparing values from electron scattering with equation (12), we obtain $r_p = (1.154 \pm 0.004) \text{ fm}$ and $a_p = (0.434 \pm 0.014) \text{ fm}$. These values are slightly modified by taking into account the

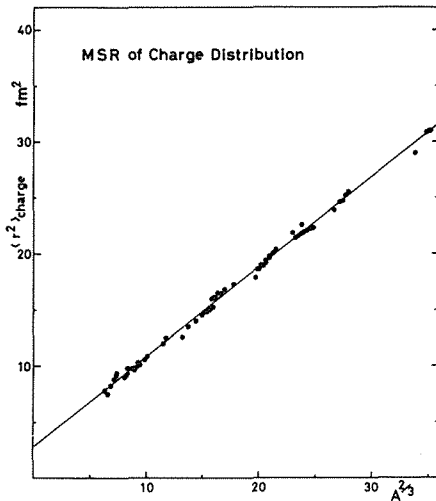


Fig. 5. Mean square radii of the charge distribution obtained from the electron scattering are plotted as a function of the mass number A .

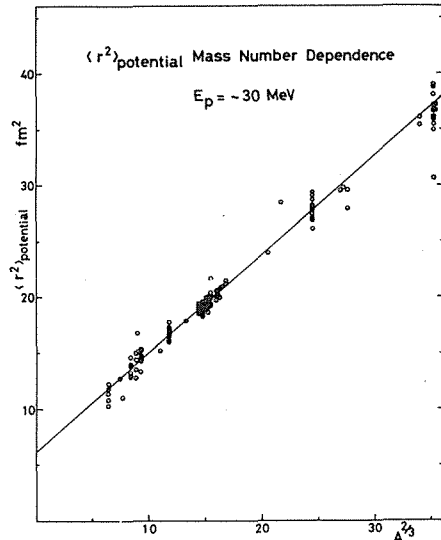


Fig. 6. Mean square radii of the real central part of the optical potentials at $E_p = 27\text{--}32 \text{ MeV}$ are plotted as a function of $A^{2/3}$.

neutron charge distribution.³¹⁾ Thus $r_p = (1.158 \pm 0.004)$ fm and $a_p = (0.470 \pm 0.012)$ fm are obtained. If the diffuseness of the point nucleon distribution is assumed to be equal to the diffuseness of the point proton distribution, we get the value of $\langle r^2 \rangle_{\text{int}} = (4.24 \pm 0.24)$ fm².

We notice that the value of r_m extracted from the present experiment is larger than that of r_p obtained from the electron scattering data. In order to show that such a difference is common to proton scattering, linear fits were made to the mean square potential radii from Perey's collection³²⁾ of optical potential results at 27-32 MeV and 47-52 MeV incident proton energies. These are shown in Fig. 6 and Fig. 7, respectively. The values obtained are $r_m = (1.21 \pm 0.03)$ fm and $\langle r^2 \rangle_{\text{int}} = (4.06 \pm 0.84)$ fm² at 30 MeV, and $r_m = (1.19 \pm 0.02)$ fm and $\langle r^2 \rangle_{\text{int}} = (4.98 \pm 0.74)$ fm² at 50 MeV. Although the 30 MeV data and the 50 MeV data consist of optical potential parameters by many authors and hence could contain many inconsistencies among the parameters obtained due to different fitting principles, the r_m and $\langle r^2 \rangle_{\text{int}}$ values from the 30 MeV and 50 MeV data are consistent with our values, and r_m is larger than the r_p value of 1.158 fm. From the analysis of the 800 MeV-1 GeV data, it is concluded that r_m is approximately equal to r_p .

One answer to this contradiction is to introduce a target mass number dependence into the effective interaction range $\langle r^2 \rangle_{\text{int}}$. As demonstrated in Fig. 3, the linear relation between $\langle r^2 \rangle_{\text{pot}}$ and $A^{2/3}$ is confirmed and acceptable. Therefore the mass number dependence to be introduced into $\langle r^2 \rangle_{\text{int}}$ must also be linear in $A^{2/3}$. A recent argument³⁴⁾ based on experimental data and the experimental results from the Los Alamos Meson Physics Facility (LAMPF),^{38)~43)} show that with a small correction from the charge distribution in the neutron itself, the difference between the root mean square

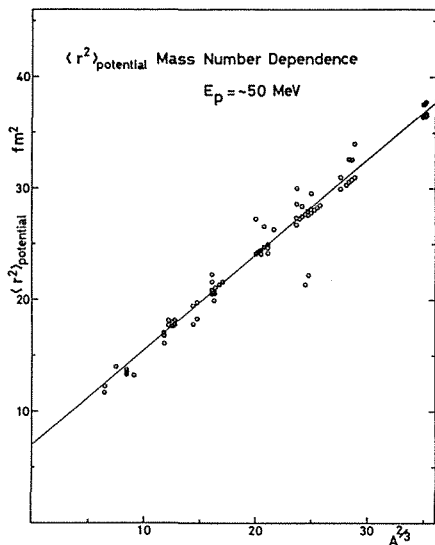


Fig. 7. Mean square radii of the real central part of the optical potentials at $E_p = 47-52$ MeV are plotted as a function of $A^{2/3}$.

radius of the point nucleon distribution and that of the point proton distribution is less than 0.1 fm, which is smaller than the error in our $\langle r^2 \rangle_{\text{int}}$ -value. Thus the average density distribution of point nucleons may be thought to be equal to that of point protons. Then the effective two-body interaction range obtained is

$$\langle r^2 \rangle_{\text{int}} = (0.132 \pm 0.013) A^{2/3} + (4.24 \pm 0.24) \text{ fm}^2 .$$

According to the recent theoretical work of Brieva⁹⁾ on the nucleon-nucleus optical potential using a realistic nucleon-nucleon interaction, the exchange term is repulsive in the nuclear center, whereas at the surface it is attractive. An exchange term of this type introduces a mass number dependence of the effective interaction range, if $\langle r^2 \rangle_{\text{int}}$ is defined as $\langle r^2 \rangle_{\text{int}} = \langle r^2 \rangle_{\text{pot}} - \langle r^2 \rangle_{\text{charge}} + \langle r^2 \rangle_{\text{proton}}$. Another source of the mass number dependence may come from a small difference between the point proton and the point neutron distributions, because the proton-neutron interaction is stronger than the proton-proton interaction, and this fact may enhance the effect from the density distribution difference. Thus the mass number dependence of $\langle r^2 \rangle_{\text{int}}$ is an empirical relation and reflects various many body effects. The true origin of this mass number dependence may be explained by an elaborate microscopic calculation. The possibility of a target dependence of the effective interaction range has been suggested already by B. Sinha.³⁰⁾ Our $\langle r^2 \rangle_{\text{int}}$ -value is larger than the GPT's value²⁸⁾ of $(2.25 \pm 0.6) \text{ fm}^2$. GPT's $\langle r^2 \rangle_{\text{int}}$ -value was obtained in the search for the χ^2 -minimum mainly of the cross section data because of the partial lack of polarization data at that time. By equally weighting the polarization and cross section data we were able to reduce the VR^n type ambiguity and have found a larger value for the mass number dependent $\langle r^2 \rangle_{\text{int}}$ -value. Bertsch et al. also obtained a large $\langle r^2 \rangle$ -value ($\sim 6 \text{ fm}^2$) for inelastic scattering by fitting the interaction to the matrix element of the scattering operator, t-matrix or G-matrix. Since their interaction is effective at the nuclear surface and is not density dependent, their $\langle r^2 \rangle_{\text{int}}$ does not have any mass number dependence. Our $\langle r^2 \rangle_{\text{int}}$ -values for medium weight nuclei are as large as the one obtained in the Bertsch's calculation.⁴⁴⁾

6. The mass number dependence of the effective two-body interaction range deduced from the optical potential for deuteron, helium-3 and alpha particles

The $\langle r^2 \rangle_{\text{pot}}$ -values for other light ion projectiles were calculated and plotted as a function of $A^{3/2}$ in Fig. 8 for 56 MeV deuteron optical potential data,⁴⁵⁾ in Fig. 9 for 109 MeV and 119 MeV ^3He ⁴⁶⁾ optical potentials and in Fig. 10 for 166 MeV alpha particles.³²⁾ In each case the solid lines are linear fits by the least square method and are expressed

$$\langle r^2 \rangle_{\text{pot}} = 0.859 A^{3/2} + 8.62 \text{ fm}^2 \quad \text{for } E_d = 56 \text{ MeV},$$

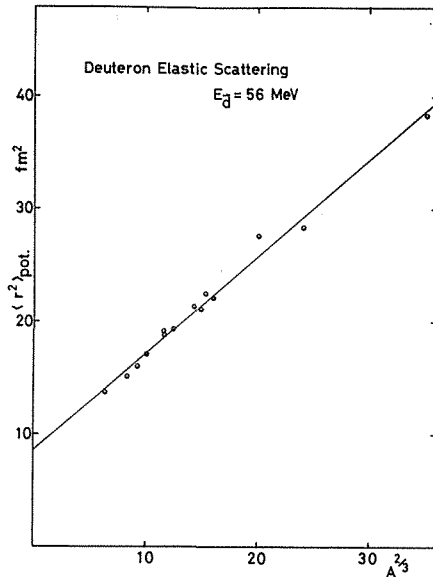


Fig. 8. Mean square radii of the real central part of the deuteron optical potential at $E_d=56$ MeV are plotted as a function of $A^{2/3}$. The solid line is a least linear fit to the data.

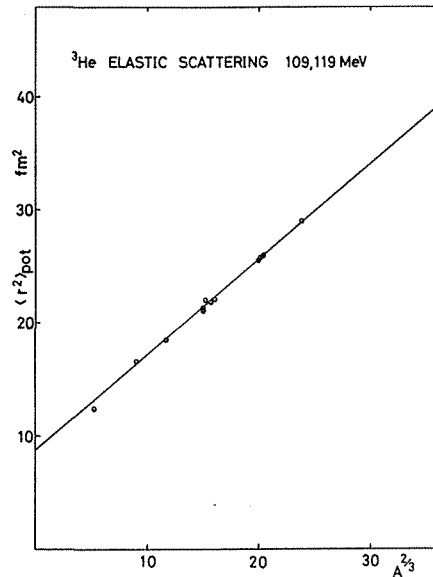


Fig. 9. Mean square radii of the real central part of the helium-3 optical potential at $E_{He}=109, 119$ MeV are plotted as a function of $A^{2/3}$. The solid line is a least linear fit to the data.

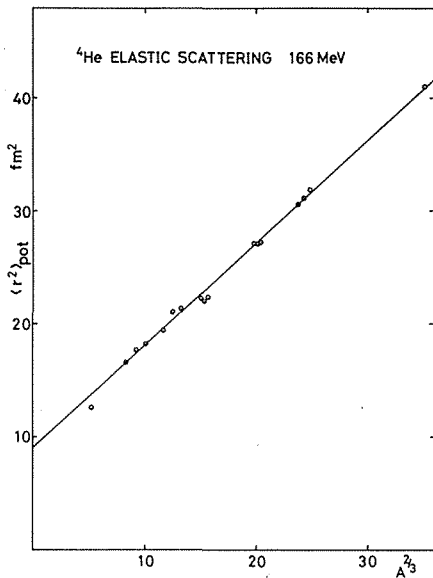


Fig. 10. Mean square radii of the real central part of the optical potential at $E_{He}=166$ MeV are plotted as a function of $A^{2/3}$. The solid line is a linear fit to the data.

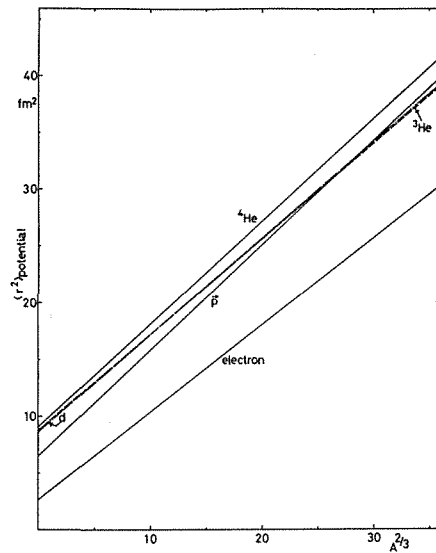


Fig. 11. Mass number dependences of the mean square radii of the real central part of the optical potential are compared with each other for $p, d, {}^3\text{He}$ and ${}^4\text{He}$ projectiles.

$$\langle r^2 \rangle_{\text{pot}} = 0.843A^{2/3} + 8.78 \text{ fm}^2 \quad \text{for } E_{\text{He}} = 109 \text{ MeV}, 119 \text{ MeV}$$

and

$$\langle r^2 \rangle_{\text{pot}} = 0.908A^{2/3} + 8.982 \text{ fm}^2 \quad \text{for } E_{\alpha} = 166 \text{ MeV},$$

respectively. If we combine all of these data, we notice in Fig. 11 that the slopes of the proton and the alpha lines are similar. The mean square radius of the real central part of the composite particle is described by the double folding model as;

$$\langle r^2 \rangle_{\text{pot}} = \frac{3}{5} r_m^2 A^{2/3} + \pi^2 a_m^2 + \langle r^2 \rangle_{\text{int}} + \langle r^2 \rangle_m^{\text{projectile}}$$

where

$$\langle r^2 \rangle_m^{\text{projectile}} = \langle r^2 \rangle_{\text{charge}}^{\text{projectile}} - \langle r^2 \rangle_{\text{charge}}^{\text{proton}} \quad \text{and}$$

$\langle r^2 \rangle_m^{\text{projectile}}$ means the mean square radius of the point nucleon distribution in the projectile. Thus the intersection in Fig. 11 is the sum of $\pi^2 a_m^2 + \langle r^2 \rangle_m^{\text{projectile}}$ and the mass number independent part of the $\langle r^2 \rangle_{\text{int}}$. The slope is the sum of $\frac{3}{5} r_m^2$ and the mass number dependent part of the $\langle r^2 \rangle_{\text{int}}$. Using the formula above and the electron scattering data for the $\langle r^2 \rangle_{\text{charge}}^{\text{projectile}}$ -values. The mean square radius of the effective two-body interaction $\langle r^2 \rangle_{\text{int}}$ is calculated as follows;

$$\begin{aligned} \langle r^2 \rangle_{\text{int}} &= (0.132A^{2/3} + 4.24) \text{ fm}^2 && \text{for proton} \\ \langle r^2 \rangle_{\text{int}} &= (0.087A^{2/3} + 2.45) \text{ fm}^2 && \text{for deuteron} \\ \langle r^2 \rangle_{\text{int}} &= (0.075A^{2/3} + 3.84) \text{ fm}^2 && \text{for helium-3} \\ \langle r^2 \rangle_{\text{int}} &= (0.140A^{2/3} + 4.89) \text{ fm}^2 && \text{for alpha} \end{aligned}$$

These values show that for the hard projectiles such as protons and alpha particles, mass number dependence of $\langle r^2 \rangle_{\text{int}}$ is large and that for the soft projectiles like deuterons and helium-3, the mass number dependences of the $\langle r^2 \rangle_{\text{int}}$ and $\langle r^2 \rangle_{\text{int}}$ -values themselves are small.

The origin of such target mass number dependence and the projectile dependence may be in the Pauli-principle. The a $A^{2/3} + b$ type mass number dependence is divergent on A and is not preferable. But it has by far the better χ^2 -value than the a $A^{-1/3} + b$ type or the a $A^{-2/3} + b$ type dependence. From the best of our knowledge we conclude that the mean square radius of the effective two-body interaction is a $A^{2/3} + b$ type. In that sense the formula of $\langle r^2 \rangle_{\text{int}}$ is an experimental one.

7. Comparison between experimentally obtained $\langle r^2 \rangle_{\text{pot}}$ -value and microscopic calculations

In the preceding section it was pointed out that there is a difference

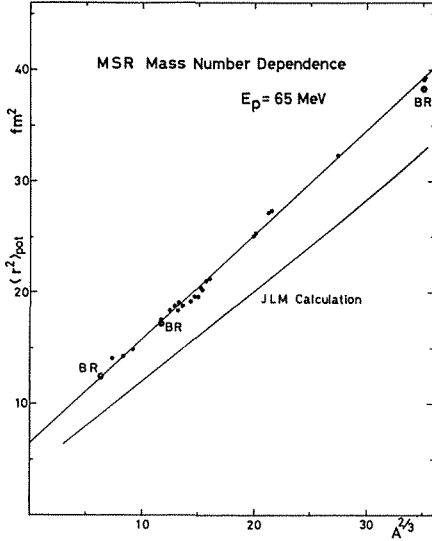


Fig. 12. Mean square radii deduced from the best fit optical potential are compared with the recent microscopic optical potential calculations. The \odot symbols denote Brieva-Rook values interpolated to $E_p=65$ MeV. The solid line marked JLM is obtained from the optical potential with the microscopically derived parameters of Jeukenne, Lejeune and Mahaux.

between $\langle r^2 \rangle_{\text{matt}}$ and $\langle r^2 \rangle_{\text{pot}}$ in the target mass number dependence. This difference was reduced to the mass number dependence of the effective two-body interaction. In Fig. 12 the calculated values based on the recent microscopic theories are shown with the experimental $\langle r^2 \rangle_{\text{pot}}$ -values. The line labeled JLM is the calculation using the JLM model. A detailed explanation of the JLM calculation will be given in Sec. 9. Brieva and Rook calculated the $\langle r^2 \rangle_{\text{pot}}$ -values of ^{16}O , ^{40}Ca and ^{208}Pb for several incident energies and their $\langle r^2 \rangle_{\text{pot}}$ -values varied smoothly with energy. The points labeled BR show the interpolation to 65 MeV proton energy of the BR calculations. The BR calculation reproduces remarkably our experimental values shown as the line labeled Kyoto data, although the BR value for ^{208}Pb is a little smaller than the observed value. Calculations in the JLM model differ greatly from the experimental values (by $2\sim 5$ fm 2) and cannot reproduce the slope of the mass number dependence. Lejeune and Hodgson⁴⁷⁾ pointed out that JLM calculation does not explain the $\langle r^2 \rangle_{\text{pot}}$ -values and angular distributions, and must be modified by introducing a phenomenological range parameter. Such a phenomenological parameter, however, will mask the validity of the theory to study the dynamics of the reaction. The main origin of the discrepancy may be in the LDA approximation used to transform the optical potential in nuclear matter to the optical potential in a finite nucleus. These two types of microscopic calculations suggest that the density dependence of the effective two-body interaction at the nuclear surface plays an essential role in explaining the A-dependence of the $\langle r^2 \rangle_{\text{int}}$ -value. It was already pointed out in nuclear matter theory that there is a density dependence⁴⁷⁾ in the nuclear matter effective interaction. As the density decreases, the effective two-body interaction increases due to the Pauli principle, and the depth of the optical potential well increases. But the A-dependence of $\langle r^2 \rangle_{\text{int}}$ is not explained by the density dependent JLM interaction using the

LDA alone, because the slope of the JLM calculation on the target mass number A is different from the experimentally observed one. (see Fig. 12) On the other hand, when obtaining the t -matrix of the r -representation by BR's calculation, the momentum sum up to the Fermi momentum gives another contribution to the density dependence⁴⁸⁾ of the interaction, in addition to the dependence coming explicitly from the Pauli principle. From our present knowledge of the BR calculation, we cannot discern the primary origin of the A -dependence of $\langle r^2 \rangle_{\text{int}}$.

8. Root mean square radius of the point nucleon distribution of the target nucleus

Applying the results obtained in the preceding section, we can extract the root mean square radius of the point nucleon distribution of the target nucleus, as follows

$$\langle r^2 \rangle_{\text{matt}}^{1/2} = (\langle r^2 \rangle_{\text{pot}} - \langle r^2 \rangle_{\text{int}})^{1/2}.$$

where

$$\langle r^2 \rangle_{\text{int}} = 4.24 \pm 0.24 + (0.132 \pm 0.013) A^{2/3} \text{ fm}^2$$

The calculated $\langle r^2 \rangle_{\text{matt}}^{1/2}$ -values from our elastic scattering data are listed in Table 3, together with the 800 MeV polarized proton elastic scattering results from LAMPF. The LAMPF $\langle r^2 \rangle_{\text{matt}}^{1/2}$ -values were calculated from the proton MSR values $\langle r^2 \rangle_p$ and the neutron MSR values $\langle r^2 \rangle_n$ of their data,^{31), 33)} using the relation

$$\langle r^2 \rangle_{\text{matt}}^{1/2} = \left(\frac{N}{A} \langle r^2 \rangle_n + \frac{Z}{A} \langle r^2 \rangle_p \right)^{1/2}$$

Although the LAMPF $\langle r^2 \rangle_n$ and $\langle r^2 \rangle_p$ values are model dependent, they are thought to be relatively free from the dynamical effects on the nucleon-nucleon interaction in the nucleus. We notice in Table 3 that our values agree remarkably well with the LAMPF results. This indicates the validity of our procedure for extracting the $\langle r^2 \rangle_{\text{int}}$ -value and the mass number dependence of the effective interaction range. We have thus obtained a new method to extract $\langle r^2 \rangle_{\text{matt}}^{1/2}$ value from polarized proton elastic scattering.

9. Volume Integral of the real central part of the optical potential

In the folding model the volume integral J_R of the real central part of the optical potential is calculated as

$$\begin{aligned} J_R &= \int V(\mathbf{r}_0) d\mathbf{r}_0^3 \\ &= \int d\mathbf{r}_0^3 \int d\mathbf{r}^3 \rho(r) V_{\text{int}}(|\mathbf{r} - \mathbf{r}_0|) \end{aligned}$$

Table 3. Volume integral per nucleon and mean square radius of the real central part of the optical potential are listed together with error bars. The effective interaction range and the nuclear matter radius are listed and compared with LAMPF-values. LAMPF-values are calculated by $\langle r^2 \rangle_{\text{matter}}^{1/2} = \left(\frac{N}{A} \langle r^2 \rangle_n + \frac{Z}{A} \langle r^2 \rangle_p \right)^{1/2}$.

Nuclei	$J_R A$ (MeV fm ³)	$\langle r^2 \rangle_{\text{pot.}}$ (fm ²)	$\langle r^2 \rangle_{\text{int.}}$ (fm ²)	$\langle r^2 \rangle_{\text{matter}}^{1/2}$ (fm)	
				Kyoto	LASL
¹⁶ O	346.8 +14.2 - 9.4	12.35 +0.06 -0.26	5.08	2.70	—
²⁰ Ne	349.5 + 9.64 -11.3	14.14 +0.02 -0.06	5.21	2.99	—
²⁴ Mg	303.0 +15.3 -14.5	14.32 +0.07 -0.23	5.34	3.00	—
²⁸ Si	325.3 + 7.8 -14.0	14.92 +0.06 -0.07	5.46	3.08	—
⁴⁰ Ar	330.4 + 7.4 -11.5	17.54 +0.05 -0.28	5.78	3.43	—
⁴⁰ Ca	333.6 + 5.2 - 8.6	17.58 +0.01 -0.18	5.78	3.43	3.39
⁴⁴ Ca	318.8 + 5.7 - 8.5	18.39 +0.03 -0.18	5.89	3.54	3.48
⁴⁸ Ca	315.0 + 6.9 - 7.5	18.39 +0.07 -0.09	5.98	3.52	3.47
⁴⁶ Ti	317.24 + 4.6 - 5.9	18.79 +0.16 -0.10	5.93	3.59	—
⁴⁸ Ti	313.1 + 4.9 - 4.0	19.08 +0.06 -0.10	5.98	3.62	—
⁵⁰ Ti	309.7 + 5.3 - 3.5	18.82 +0.08 -0.03	6.03	3.58	—
⁵⁴ Fe	303.1 + 5.9 - 5.1	19.18 +0.14 -0.24	6.13	3.61	3.57
⁵⁶ Fe	307.1 + 2.8 - 3.2	19.69 +0.07 -0.16	6.17	3.68	—
⁵⁹ Co	307.8 + 4.6 - 4.0	20.47 +0.09 -0.09	6.24	3.77	—
⁵⁸ Ni	300.2 + 2.6 - 2.0	19.63 +0.06 -0.20	6.22	3.66	3.67 or 3.70
⁶⁰ Ni	306.8 + 2.1 - 3.4	20.25 +0.07 -0.12	6.26	3.74	—
⁶² Ni	309.4 + 5.1 - 5.0	21.07 +0.21 -0.19	6.31	3.84	—
⁶⁴ Ni	315.1 + 3.3 - 3.2	21.27 +0.05 -0.12	6.35	3.86	3.86
⁸⁰ Y	323.4 + 2.7 - 2.4	25.09 +0.1 -0.35	6.87	4.27	—
⁹⁰ Zr	318.9 + 1.3 - 4.4	25.35 +0.19 -0.41	6.89	4.30	4.25
⁹⁸ Mo	321.0 + 2.7 - 5.8	27.20 +0.30 -0.13	7.05	4.49	—
¹⁰⁰ Mo	326.0 + 3.4 - 1.0	27.38 +0.16 -0.31	7.08	4.51	—
¹⁴⁴ Sm	317.4 + 3.4 - 2.2	32.18 +0.23 -0.39	7.87	4.93	—
²⁰⁸ Pb	330.7 + 2.3 - 7.0	39.18 +0.64 -0.64	8.87	5.51	5.55
²⁰⁹ Bi	319.7 + 3.3 - 3.6	39.36 +0.46 -0.71	8.89	5.52	—

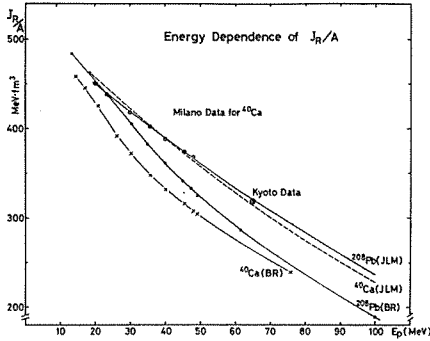


Fig. 13. Volume integral values per nucleon of the optical potentials are plotted as a function of incident proton energy. Open circles are Milan data for ^{40}Ca and double circle is a mean value over 25 targets at $E_p=65$ MeV. The solid dots and crosses are calculated microscopically by Brieva and Rook, connected by curves meant to guide the eyes. The solid line and the broken line labeled JLM are calculated values using procedure of Jeukenne, Lejeune and Mahaux.

$$= A \int V_{\text{int}}(a) da^3$$

The volume integral value J_R is proportional to the target mass number if the effective two-body interaction potential between the projectile and the target nucleon is independent of density and energy. A linear fit to J_R -values confirms the approximate validity of the above assumption at 65 MeV, and the volume integral is expressed as

$$J_R = (318 \pm 3) A \text{ MeV fm}^3$$

In order to show how the recent microscopic calculation explain the empirical volume integral values, J_R/A -values are plotted as a function of incident proton energy in Fig. 13. The double circle point at 65 MeV is the average value of the Kyoto data. The open circles show the recent measurement for ^{40}Ca by the Milan group.^{13),14)} We notice that the Kyoto data lie on a smooth extrapolation of lower energy Milan data. The solid and dashed curves labeled JLM indicate the volume integral of the microscopic optical potential calculated in the JLM model. According to the parameter table of the JLM calculation⁴⁾ the optical potential in infinite nuclear matter is obtained as a function of the matter density and the incident projectile energy. As for the point nucleon density distribution for the JLM calculation, we used Negele's density distribution obtained from the fitting to the electron scattering data as in JLM's work. The solid curve shows the calculation for ^{208}Pb and the dashed curve shows the ^{40}Ca case. The results of Brieva-Rook calculation are indicated by the points marked with a cross in the figure for ^{40}Ca and ^{208}Pb nuclei; the curves connecting these points are only meant to guide the eye. As is evident from the figure, the J_R/A value and its bombarding energy dependence are reproduced remarkably well by the JLM calculation. The local density approximation used in the JLM model seems to be effective for the J_R calculation. The energy dependence and the density dependence of the nuclear matter t -matrix are directly reflected in the JLM nuclear optical potential. On the other hand Brieva and Rook trans-

formed the nuclear matter t -matrix in momentum-representation to the t -matrix in the r -representation using a suitable approximation. They then calculated the optical potentials for finite nuclei by applying the folding approximation. The calculation of BR explains the empirical results at 30 MeV, but deviates from the observed values at higher energies. The origin of the discrepancy between the BR's calculation and the experimental data seems to be due to the approximation in their transformation to the r -representation.

In Fig. 14 our J_R/A values listed in Table 3 are plotted as a function of the target mass number. Error bars in the figure were defined similar as the error bars of $\langle r^2 \rangle_{\text{pot}}$ -values in Sec. 5 but in this case were deduced from the potential parameters at $\chi^2_{\text{min}}(V_r) = 1.25 \chi^2_0$ (best fit), so the error bars have no statistical meaning. Although observed J_R/A values scatter considerably, we notice the following global behavior around the average value of $J_R/A = 318$ MeV. As the mass number A increases, the J_R/A value decreases

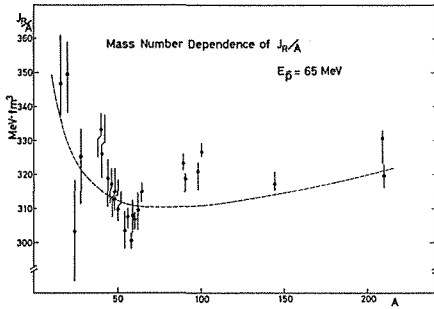


Fig. 14. Volume integral values per nucleon of the real central part of the optical potentials are plotted as a function of the target mass number A . The definition of the error bars is given in the text. The dashed curve is the JLM model calculation for $E_p = 65$ MeV.

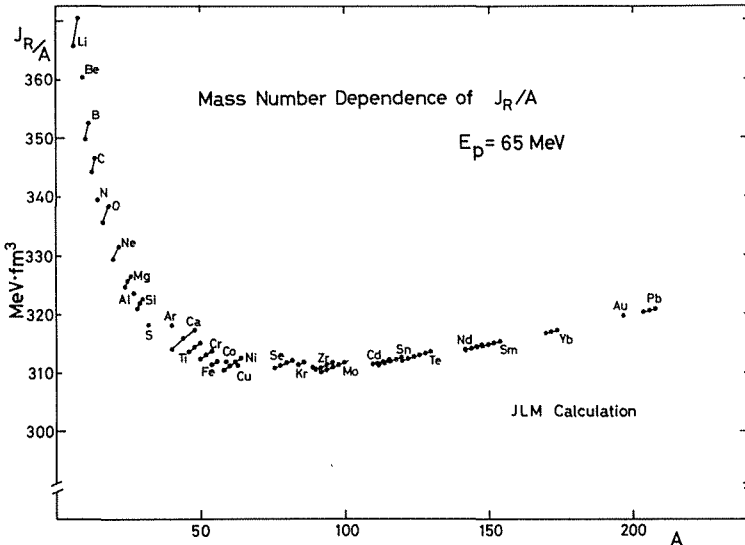


Fig. 15. Volume integral values per nucleon of the real central part of the optical potentials are plotted for each isotope, using Juekenne, Lejeune and Mahaux's model.

rapidly to the minimum in the Fe-Ni region and then increases gradually toward the Pb-Bi region. This global trend is remarkably reproduced by the JLM model calculation shown by the dashed curve in Fig. 14. According to the JLM model, the effective interaction is density dependent. In the lower density region, the effective interaction is stronger^{3),4)} due to the smaller Pauli blocking effect. The surface-to-volume ratio is large in light nuclei. As the target mass number increases, the surface-to-volume ratio decreases as $A^{-1/3}$ and the J_R/A -value decreases rapidly. The second gradual increase is explained in the JLM model by the isospin dependent interaction and by the velocity dependence of the effective interaction. (As the mass number increases, the velocity of the projectile inside the nucleus decreases due to the repulsive Coulomb potential.) This global trend in Fig. 14 is similar to the binding energy per nucleon curve, if we remind that the Coulomb potential is included in the binding energy calculation and the velocity dependent effect is included in the J_R/A calculation. The rapid change in J_R/A -values for lighter nuclei evident in Fig. 14 is possibly evidence of the density dependence of the effective interaction.

10. Anomalous isotope dependence of the real central part of the optical potential in the f - p shell region nuclei

C. M. Perey and F. C. Perey analyzed the elastic scattering of 11 MeV polarized protons on 20 target isotopes from ^{48}Ti to ^{76}Ge . For the depth of the real central part of the optical potential they obtained an anomalous linear relation of $V_R = V_0 + \alpha A$ without the isospin dependence. At 65 MeV incident proton energy a similar anomaly was reported in our previous work.²⁰⁾ Our previous conclusion was that there must be a linear mass number dependence of $V_{rT} = -1.8 + 0.72(A - 40)$ in the real isospin dependent part of the optical potential. In Fig. 16 the volume integral values per nucleon are plotted versus $(N - Z)/A$. The slope of the line connecting the same isotopes changes from negative to positive as the mass number of the isotope increases from ^{40}Ca to ^{64}Ni . In the case of the ordinary isospin dependence

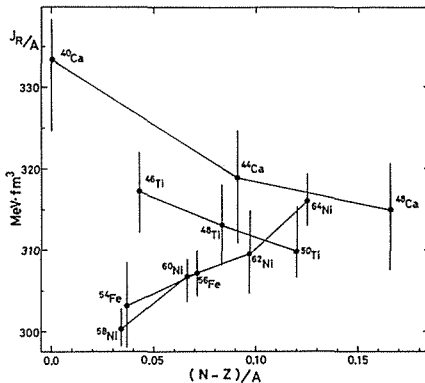


Fig. 16. Observed J_A/A -values are plotted versus $(N-Z)/A$ for f - p shell nuclei. The definition of the uncertainties is the same as in Fig. 14.

the slope must be constant. The observed behavior in Fig. 16 is anomalous in this respect. As shown in Fig. 15, the global JLM microscopic model predicts always the positive slope in this mass number region. So it cannot explain the anomalous isotope dependence.

Using our mean square matter radius data in Sec. 5 and the diffuseness value obtained by fitting to the electron scattering data we now have a new density distributions for each target nuclei. The density distribution is assumed to be the Fermi type.

$$\rho(r) = \frac{\rho_0}{1 + \exp((r - R_m)/a_m)}$$

where the R_m value is calculated from the formula;

$$\langle r^2 \rangle_{\text{nucleon}} + \langle r^2 \rangle_{\text{matt}} = \frac{3}{5} R_m^2 + \frac{7}{5} \pi^2 a_m^2$$

and $a_m = 0.470$ is obtained in Sec. 5 by fitting to the electron scattering data. $\langle r^2 \rangle_{\text{nucleon}}$ is the mean square radius of a nucleon and is estimated to be as same as $\langle r^2 \rangle_{\text{proton}}$ (0.64 fm^2).

Using this density distribution and the JLM model, the volume integrals of the real central part of the optical potential are calculated for each target nuclei. The calculated J_R/A -values are shown in Fig. 17. The main part of the anomalous isotope effect is reproduced by the JLM model calculations using our new matter distribution data. The origin of the anomalous isotope effect consist of two parts, the density dependence of the effective interaction and the shell closure effect. For nuclei with $N=28$ the size of the nucleus is contracted comparative to their neighbouring nuclei. Since the matter density for the $N=28$ nuclei increases, the strength of the effective interaction is reduced according to its density dependence. Thus for nuclei $A \leq 53$ the J_A/A value takes the negative slope versus the mass number. But in the figure we notice differences between calculation and experimental data, which will be explained only by the nuclear structure dependent effective interaction.

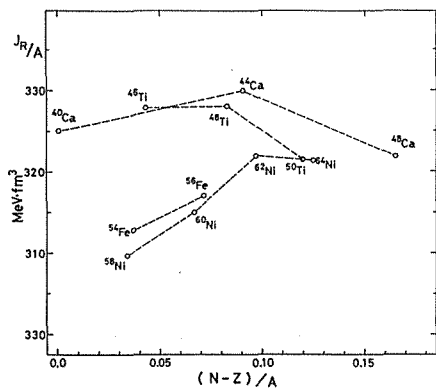


Fig. 17. Calculated J_R/A -values are plotted versus $(N-Z)/A$ for f - p shell nuclei.

11. Conclusions

We have systematically measured polarized proton elastic scattering from 25 targets at 65 MeV. An optical potential analysis gave good fits to both cross section data and analyzing power data. By plotting the mean square radius of the real central part of the optical potential versus $A^{2/3}$, we have obtained the global systematics of the MSR of the potential as $\langle r^2 \rangle_{\text{pot}} = (6.42 \pm 0.21) + (0.937 \pm 0.012) A^{2/3} \text{ fm}^2$. Using the simple folding model and by comparing with the charge distribution obtained by electron scattering, we found that the effective interaction range has a mass number dependence of the form

$$\langle r^2 \rangle_{\text{int}} = (4.24 \pm 0.24) + (0.132 \pm 0.013) A^{2/3} \text{ fm}^2.$$

Assuming this mass number dependence of the effective interaction range, we have obtained root mean square radii of the point nucleon distribution, which are in accord with the high energy LAMPF data and reflect the shell effect and the individual characteristics of the target nuclei. This mass number dependence of $\langle r^2 \rangle_{\text{int}}$ was shown to be different for different projectiles such as d , ${}^3\text{He}$ or α .

For nuclei of $A < 58$, the J_R/A value decreases as the mass number of the target increases. This rapid decrease was interpreted as evidence of the density dependence of the effective interaction. The JLM model explains both energy and A dependence of the J_R/A but cannot explain the value of $\langle r^2 \rangle_{\text{pot}}$, and its A -dependence.

On the other hand the BR calculation reproduces our $\langle r^2 \rangle_{\text{pot}}$ -values but could not predict the J_R/A -values, especially the energy dependence of J_R/A . At present, each of the two global theories could explain the experimental results only partially, but is found to be an effective guide in clarifying nuclear many body dynamics. The anomalous isotope dependence of J_R/A is explained qualitatively by the JLM density dependent interaction and the matter radius obtained from our experiment.

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University and the polarized ion source team under Prof. K. Nishimura of Kyoto University. Their wonderful operation of the cyclotron and the polarized ion source are greatly appreciated. Sincere thanks are also expressed to Prof. H. Ogata, Prof. K. Hosono and Dr. N. Matsuoka of RCNP, Osaka University for their hospitalities during the experiment. The author owes thanks to Prof. S. Nagata of Miyazaki University for fruitful discussions on the microscopic optical potential. He is also grateful to Mrs. E. Takahashi and other secretaries for their patience in typing the manuscript.

^{44}Ca			^{46}Ca			^{48}Ti		
Angle	Cross Section (mb/sr)	Analyzing Power	Angle	Cross Section (mb/sr)	Analyzing Power	Angle	Cross Section (mb/sr)	Analyzing Power
15.16	1581.3 ± 2.0	0.0002 ± 0.0020	24.49	42.30 ± 0.84	0.160 ± 0.020	15.35	1793.1 ± 8.3	-0.0095 ± 0.0063
16.18	1227.8 ± 3.9	-0.0234 ± 0.0050	25.51	36.12 ± 0.78	0.544 ± 0.019	16.37	1396.4 ± 5.2	-0.0338 ± 0.0050
17.72	827.7 ± 1.4	-0.0575 ± 0.0030	26.53	37.97 ± 0.80	0.775 ± 0.020	17.39	1067.7 ± 6.4	-0.0426 ± 0.0080
19.25	505.5 ± 2.5	-0.0937 ± 0.0070	27.55	44.19 ± 0.86	0.831 ± 0.016	18.42	771.4 ± 3.9	-0.0771 ± 0.0066
20.28	355.45 ± 0.93	-0.1262 ± 0.0041	28.57	54.69 ± 0.96	0.760 ± 0.018	19.44	535.1 ± 3.2	-0.1125 ± 0.0084
21.30	227.13 ± 1.02	-0.1487 ± 0.0060	29.58	64.62 ± 1.04	0.678 ± 0.014	20.46	373.6 ± 2.2	-0.1440 ± 0.0080
22.32	149.49 ± 0.47	-0.1593 ± 0.0051	30.60	71.68 ± 1.10	0.6026 ± 0.0096	21.48	242.1 ± 2.2	-0.165 ± 0.012
23.34	88.16 ± 0.57	-0.1398 ± 0.0090	33.14	77.90 ± 0.74	0.473 ± 0.011	22.50	147.0 ± 1.1	-0.176 ± 0.010
24.37	56.63 ± 0.29	0.0076 ± 0.0079	35.69	66.63 ± 0.49	0.3875 ± 0.0080	23.52	88.98 ± 0.83	-0.116 ± 0.013
25.39	35.75 ± 0.36	0.334 ± 0.013	38.24	46.13 ± 0.40	0.2990 ± 0.0092	24.55	54.59 ± 0.51	0.0025 ± 0.0127
26.41	29.51 ± 0.27	0.670 ± 0.013	40.77	27.46 ± 0.31	0.247 ± 0.012	25.57	38.30 ± 0.43	0.381 ± 0.015
27.43	33.60 ± 0.25	0.828 ± 0.014	43.31	13.81 ± 0.16	0.300 ± 0.012	26.59	35.41 ± 0.52	0.704 ± 0.018
28.45	38.26 ± 0.24	0.843 ± 0.012	45.85	7.650 ± 0.134	0.499 ± 0.018	27.61	37.93 ± 0.43	0.825 ± 0.016
29.48	44.95 ± 0.33	0.771 ± 0.014	48.39	5.952 ± 0.103	0.885 ± 0.016	28.63	45.51 ± 0.42	0.826 ± 0.014
30.50	50.95 ± 0.27	0.743 ± 0.011	50.92	6.230 ± 0.086	0.988 ± 0.013	29.65	52.74 ± 0.50	0.762 ± 0.015
33.05	59.90 ± 0.22	0.5586 ± 0.0080	53.46	6.361 ± 0.095	0.948 ± 0.014	30.67	60.89 ± 0.49	0.664 ± 0.012
35.60	54.67 ± 0.29	0.4594 ± 0.0095	55.99	5.915 ± 0.073	0.871 ± 0.012	33.22	67.64 ± 0.52	0.515 ± 0.011
38.15	41.60 ± 0.13	0.3760 ± 0.0065	58.52	4.696 ± 0.075	0.801 ± 0.015	35.77	59.33 ± 0.49	0.419 ± 0.012
40.70	26.81 ± 0.20	0.318 ± 0.011	61.04	3.220 ± 0.054	0.711 ± 0.015	38.32	43.11 ± 0.42	0.342 ± 0.013
43.24	14.783 ± 0.076	0.3098 ± 0.0082	63.57	2.190 ± 0.051	0.689 ± 0.022	40.86	26.63 ± 0.23	0.273 ± 0.012
45.79	7.757 ± 0.039	0.405 ± 0.014	66.09	1.299 ± 0.028	0.575 ± 0.020	43.41	14.05 ± 0.17	0.270 ± 0.016
48.33	4.874 ± 0.039	0.689 ± 0.013	68.61	0.958 ± 0.026	0.635 ± 0.026	45.95	8.050 ± 0.104	0.422 ± 0.016
50.87	4.295 ± 0.041	0.946 ± 0.015	71.13	0.826 ± 0.022	0.761 ± 0.023	48.49	5.195 ± 0.084	0.768 ± 0.020
53.41	4.211 ± 0.036	0.996 ± 0.015				51.03	4.827 ± 0.070	1.001 ± 0.017
55.94	4.063 ± 0.036	0.944 ± 0.015				56.10	4.450 ± 0.060	0.994 ± 0.018
58.48	3.378 ± 0.033	0.860 ± 0.015				61.16	2.674 ± 0.038	0.779 ± 0.018
61.01	2.649 ± 0.023	0.822 ± 0.014				63.69	1.777 ± 0.043	0.683 ± 0.029
66.07	1.154 ± 0.014	0.697 ± 0.017				66.22	1.147 ± 0.024	0.675 ± 0.026
71.11	0.604 ± 0.010	0.688 ± 0.021				68.74	0.860 ± 0.021	0.648 ± 0.030

⁴⁶ Ti				⁴⁶ Ti				⁵⁶ Fe			
Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power	
15.33	1644.1 ± 3.5	-0.155 ± 0.0028		15.32	1670.8 ± 6.6	-0.0225 ± 0.0052		15.30	2050.3 ± 5.7	-0.0162 ± 0.0037	
16.35	1281.1 ± 3.1	-0.0360 ± 0.0032		16.34	1286.4 ± 5.8	-0.0429 ± 0.0059		16.32	1522.2 ± 4.9	-0.0394 ± 0.0044	
17.38	964.4 ± 2.7	-0.0614 ± 0.0037		17.36	956.7 ± 3.6	-0.0700 ± 0.0050		17.33	1171.1 ± 4.3	-0.0717 ± 0.0050	
18.40	684.5 ± 2.3	-0.0864 ± 0.0044		18.38	688.4 ± 3.0	-0.1019 ± 0.0059		18.35	809.2 ± 3.6	-0.0966 ± 0.0061	
19.42	495.7 ± 2.4	-0.1012 ± 0.0063		19.40	475.8 ± 2.5	-0.1279 ± 0.0071		19.37	567.0 ± 3.0	-0.1528 ± 0.0073	
20.44	327.9 ± 1.2	-0.1462 ± 0.0051		20.42	312.9 ± 2.0	-0.1430 ± 0.0087		20.39	357.4 ± 1.7	-0.1770 ± 0.0067	
21.46	212.1 ± 1.5	-0.1648 ± 0.0097		21.44	193.6 ± 1.6	-0.195 ± 0.011		21.41	223.3 ± 1.3	-0.2295 ± 0.0085	
22.48	131.35 ± 0.77	-0.1693 ± 0.0078		22.46	115.6 ± 1.2	-0.172 ± 0.014		22.43	130.10 ± 0.83	-0.2206 ± 0.0089	
23.50	76.69 ± 0.93	-0.087 ± 0.016		23.48	69.31 ± 0.78	-0.081 ± 0.015		22.94	96.94 ± 0.72	-0.184 ± 0.010	
24.52	49.80 ± 0.47	0.096 ± 0.012		24.50	43.96 ± 0.62	0.185 ± 0.019		23.45	73.90 ± 0.63	-0.115 ± 0.012	
25.54	36.88 ± 0.45	0.448 ± 0.016		25.52	34.68 ± 0.43	0.575 ± 0.016		24.47	47.17 ± 0.50	0.185 ± 0.014	
26.56	35.68 ± 0.40	0.753 ± 0.015		26.54	36.43 ± 0.44	0.803 ± 0.016		25.48	38.70 ± 0.45	0.662 ± 0.016	
27.58	40.15 ± 0.42	0.822 ± 0.015		27.56	45.52 ± 0.49	0.808 ± 0.015		25.99	40.85 ± 0.47	0.780 ± 0.016	
28.60	48.77 ± 0.33	0.781 ± 0.012		28.58	54.92 ± 0.54	0.753 ± 0.014		26.50	44.04 ± 0.48	0.872 ± 0.015	
29.62	57.18 ± 0.51	0.707 ± 0.013		29.60	64.23 ± 0.58	0.674 ± 0.013		27.52	55.26 ± 0.47	0.840 ± 0.014	
30.64	63.36 ± 0.38	0.631 ± 0.010		30.62	70.84 ± 0.61	0.603 ± 0.012		28.54	68.96 ± 0.47	0.741 ± 0.012	
33.19	68.28 ± 0.57	0.494 ± 0.012		33.16	75.68 ± 0.65	0.451 ± 0.012		29.55	82.19 ± 0.51	0.644 ± 0.011	
35.74	60.47 ± 0.53	0.391 ± 0.012		35.71	63.72 ± 0.27	0.3548 ± 0.0067		30.57	89.93 ± 0.54	0.567 ± 0.010	
38.28	42.26 ± 0.32	0.301 ± 0.010		38.25	45.46 ± 0.41	0.297 ± 0.012		33.11	95.52 ± 0.40	0.4180 ± 0.0085	
40.83	24.96 ± 0.20	0.242 ± 0.011		40.79	26.00 ± 0.17	0.2293 ± 0.0089		35.66	77.07 ± 0.36	0.3228 ± 0.0082	
43.37	13.24 ± 0.15	0.279 ± 0.015		43.33	13.59 ± 0.16	0.262 ± 0.015		38.20	53.84 ± 0.30	0.2411 ± 0.0094	
45.91	7.265 ± 0.070	0.474 ± 0.013		45.87	7.610 ± 0.084	0.516 ± 0.015		40.74	30.58 ± 0.23	0.179 ± 0.011	
48.45	5.361 ± 0.060	0.822 ± 0.015		48.41	5.820 ± 0.068	0.899 ± 0.015		43.27	16.09 ± 0.17	0.269 ± 0.016	
50.99	5.139 ± 0.059	1.006 ± 0.015		50.95	6.012 ± 0.058	0.988 ± 0.014		45.81	9.759 ± 0.091	0.600 ± 0.015	
53.52	5.236 ± 0.060	0.970 ± 0.015		53.48	6.077 ± 0.070	0.955 ± 0.015		48.34	8.460 ± 0.085	0.892 ± 0.017	
56.05	4.724 ± 0.057	0.879 ± 0.016		56.01	5.522 ± 0.056	0.861 ± 0.014		50.88	8.283 ± 0.068	0.999 ± 0.015	
58.58	3.731 ± 0.042	0.810 ± 0.015		58.54	4.302 ± 0.049	0.785 ± 0.015		53.41	8.097 ± 0.068	0.941 ± 0.015	
61.11	2.634 ± 0.035	0.751 ± 0.017		61.07	3.020 ± 0.041	0.684 ± 0.017		55.94	7.017 ± 0.063	0.835 ± 0.015	
63.64	1.764 ± 0.029	0.661 ± 0.020		63.59	1.908 ± 0.033	0.660 ± 0.021		58.46	5.358 ± 0.055	0.755 ± 0.017	
66.16	1.133 ± 0.023	0.618 ± 0.025		66.12	1.240 ± 0.027	0.593 ± 0.026		60.99	3.674 ± 0.042	0.659 ± 0.018	
68.69	0.827 ± 0.020	0.674 ± 0.029		68.64	0.990 ± 0.024	0.636 ± 0.028		63.51	2.421 ± 0.344	0.587 ± 0.022	
71.21	0.696 ± 0.015	0.746 ± 0.025		71.16	0.873 ± 0.018	0.753 ± 0.024		66.04	1.738 ± 0.024	0.592 ± 0.021	
								68.56	1.431 ± 0.022	0.626 ± 0.023	
								71.07	1.264 ± 0.019	0.734 ± 0.021	

⁵⁶ Fe				⁵⁸ Co				⁵⁸ Ni			
Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power	
14.27	2444.8 ± 8.0	-0.0161 ± 0.0051		15.37	1794.9 ± 4.9	-0.0523 ± 0.0039		14.26	2578.8 ± 11.1	-0.0199 ± 0.0063	
16.30	1417.3 ± 8.6	-0.0550 ± 0.0093		16.39	1311.4 ± 4.2	-0.0759 ± 0.0045		16.29	1503.1 ± 6.1	-0.0705 ± 0.0059	
18.34	700.1 ± 3.0	-0.1253 ± 0.0068		17.41	946.6 ± 3.5	-0.1165 ± 0.0055		18.33	746.1 ± 3.0	-0.1439 ± 0.0062	
20.38	286.0 ± 1.9	-0.243 ± 0.011		18.43	627.4 ± 1.2	-0.1511 ± 0.0061		20.36	303.7 ± 1.9	-0.2368 ± 0.0095	
21.40	165.9 ± 1.5	-0.280 ± 0.014		19.44	407.4 ± 1.6	-0.1982 ± 0.0031		21.38	174.1 ± 1.2	-0.282 ± 0.010	
22.41	90.24 ± 0.88	-0.188 ± 0.015		20.46	250.60 ± 0.74	-0.2451 ± 0.0050		22.40	96.55 ± 0.89	-0.243 ± 0.013	
23.43	53.21 ± 0.63	0.0028 ± 0.0176		21.48	141.55 ± 0.97	-0.266 ± 0.010		23.42	53.62 ± 0.66	-0.0074 ± 0.0175	
24.45	39.04 ± 0.54	0.502 ± 0.021		22.49	80.62 ± 0.42	-0.1603 ± 0.0075		24.43	38.86 ± 0.49	0.465 ± 0.018	
25.47	39.65 ± 0.51	0.797 ± 0.019		23.51	48.59 ± 0.57	0.138 ± 0.016		24.94	38.71 ± 0.19	0.692 ± 0.010	
26.48	51.49 ± 0.58	0.861 ± 0.018		24.53	40.75 ± 0.52	0.647 ± 0.017		25.45	41.34 ± 0.58	0.842 ± 0.020	
28.52	77.82 ± 0.64	0.662 ± 0.014		25.54	48.32 ± 0.40	0.815 ± 0.013		25.96	46.05 ± 0.61	0.873 ± 0.019	
30.55	95.90 ± 0.64	0.522 ± 0.012		26.56	61.72 ± 0.64	0.775 ± 0.015		26.47	51.12 ± 0.65	0.865 ± 0.018	
33.09	94.30 ± 0.51	0.3896 ± 0.0091		27.58	76.99 ± 0.51	0.705 ± 0.011		27.48	65.95 ± 0.73	0.768 ± 0.017	
35.63	72.10 ± 0.36	0.2965 ± 0.0084		28.59	90.88 ± 0.77	0.605 ± 0.013		28.50	82.23 ± 0.82	0.684 ± 0.015	
38.17	45.42 ± 0.35	0.193 ± 0.012		29.61	98.70 ± 0.57	0.5401 ± 0.0097		30.53	102.34 ± 0.79	0.511 ± 0.012	
40.71	24.36 ± 0.21	0.196 ± 0.014		30.63	103.07 ± 0.83	0.484 ± 0.012		33.07	97.67 ± 0.50	0.3741 ± 0.0089	
43.24	12.47 ± 0.15	0.345 ± 0.018		31.64	102.44 ± 0.58	0.4258 ± 0.0091		35.61	75.59 ± 0.44	0.2698 ± 0.0090	
45.78	8.271 ± 0.087	0.738 ± 0.017		33.67	89.67 ± 0.54	0.3346 ± 0.0092		38.15	47.63 ± 0.29	0.1999 ± 0.0091	
48.31	7.911 ± 0.093	0.960 ± 0.018		35.70	67.30 ± 0.39	0.2544 ± 0.0085		40.68	26.35 ± 0.19	0.176 ± 0.010	
50.84	8.217 ± 0.071	0.947 ± 0.015		38.24	40.18 ± 0.30	0.194 ± 0.011		43.22	13.66 ± 0.12	0.325 ± 0.013	
53.37	7.643 ± 0.075	0.886 ± 0.017		40.77	20.76 ± 0.21	0.207 ± 0.014		45.75	9.169 ± 0.077	0.735 ± 0.014	
55.90	6.198 ± 0.062	0.788 ± 0.016		43.31	11.26 ± 0.13	0.457 ± 0.016		48.28	8.896 ± 0.088	0.978 ± 0.016	
58.43	4.376 ± 0.057	0.690 ± 0.022		45.84	8.736 ± 0.089	0.829 ± 0.015		50.82	8.982 ± 0.076	0.951 ± 0.015	
60.95	2.879 ± 0.042	0.601 ± 0.022		48.37	8.949 ± 0.050	0.956 ± 0.012		53.34	8.513 ± 0.067	0.869 ± 0.014	
63.48	1.885 ± 0.029	0.561 ± 0.024		50.90	8.949 ± 0.079	0.917 ± 0.014		55.87	6.796 ± 0.060	0.768 ± 0.014	
66.00	1.333 ± 0.024	0.594 ± 0.026		53.43	7.961 ± 0.074	0.816 ± 0.014		58.40	4.852 ± 0.041	0.670 ± 0.013	
68.52	1.185 ± 0.019	0.716 ± 0.023		55.96	6.135 ± 0.058	0.752 ± 0.014		60.92	3.188 ± 0.041	0.562 ± 0.018	
71.04	1.104 ± 0.018	0.843 ± 0.023		58.48	4.199 ± 0.048	0.666 ± 0.016		63.44	2.148 ± 0.024	0.564 ± 0.016	
				61.01	2.659 ± 0.031	0.586 ± 0.016		65.96	1.587 ± 0.020	0.623 ± 0.018	
				63.53	1.778 ± 0.026	0.607 ± 0.019		68.48	1.404 ± 0.016	0.734 ± 0.016	
				66.05	1.448 ± 0.020	0.708 ± 0.018		71.00	1.318 ± 0.015	0.827 ± 0.016	
				71.08	1.240 ± 0.017	0.897 ± 0.017					
				76.11	0.896 ± 0.013	0.968 ± 0.018					

^{60}Ni				^{60}Ni				^{60}Ni			
Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power	
14.25	2776.7 ± 28.7	-0.0153 ± 0.0048		13.33	3351.6 ± 17.8	-0.0094 ± 0.0039		13.32	3256.0 ± 11.9	-0.0189 ± 0.0034	
16.28	1553.8 ± 16.7	-0.0809 ± 0.0068		14.34	2585.3 ± 14.3	-0.0334 ± 0.0043		14.34	2472.6 ± 9.5	-0.0406 ± 0.0038	
18.32	725.7 ± 7.8	-0.1531 ± 0.0069		16.38	1369.8 ± 8.45	-0.1028 ± 0.0061		16.37	1284.2 ± 5.9	-0.1048 ± 0.0056	
19.34	459.8 ± 5.2	-0.2213 ± 0.0087		18.41	594.7 ± 4.2	-0.2082 ± 0.0082		18.40	537.6 ± 3.0	-0.2227 ± 0.0080	
20.35	273.7 ± 3.9	-0.285 ± 0.016		19.43	365.0 ± 2.6	-0.2624 ± 0.0087		19.42	312.6 ± 2.5	-0.279 ± 0.012	
21.37	152.5 ± 1.9	-0.295 ± 0.012		20.44	202.5 ± 1.8	-0.305 ± 0.011		20.43	170.7 ± 1.6	-0.353 ± 0.015	
22.38	80.28 ± 1.35	-0.235 ± 0.021		21.46	107.7 ± 1.2	-0.303 ± 0.015		21.45	85.54 ± 0.89	-0.295 ± 0.015	
23.40	49.38 ± 0.85	0.1447 ± 0.021		22.48	57.74 ± 0.68	-0.074 ± 0.016		22.46	49.82 ± 0.58	0.090 ± 0.018	
24.42	41.59 ± 0.89	0.686 ± 0.026		23.49	40.78 ± 0.56	0.510 ± 0.019		23.48	41.23 ± 0.60	0.627 ± 0.021	
24.93	46.31 ± 0.52	0.819 ± 0.012		24.51	45.88 ± 0.60	0.848 ± 0.018		24.49	53.25 ± 0.84	0.832 ± 0.024	
25.43	51.93 ± 0.76	0.880 ± 0.017		25.52	62.84 ± 0.71	0.861 ± 0.016		25.51	74.70 ± 0.72	0.792 ± 0.016	
26.45	69.70 ± 0.12	0.819 ± 0.021		26.54	83.41 ± 1.04	0.736 ± 0.018		26.52	95.05 ± 0.94	0.722 ± 0.017	
28.48	103.53 ± 1.60	0.602 ± 0.018		28.57	116.29 ± 1.27	0.564 ± 0.015		28.55	127.64 ± 1.33	0.514 ± 0.016	
30.51	120.11 ± 1.32	0.4656 ± 0.0097		30.50	127.56 ± 0.90	0.421 ± 0.010		30.48	132.92 ± 0.70	0.3836 ± 0.0087	
33.05	110.36 ± 1.37	0.334 ± 0.012		30.60	125.39 ± 1.19	0.424 ± 0.013		30.58	134.37 ± 0.91	0.3866 ± 0.0099	
35.59	79.19 ± 0.92	0.2577 ± 0.0098		33.04	108.92 ± 0.80	0.3108 ± 0.095		33.02	107.27 ± 0.62	0.2887 ± 0.0085	
38.13	46.80 ± 0.60	0.169 ± 0.012		35.57	73.27 ± 0.61	0.208 ± 0.011		35.55	68.54 ± 0.47	0.1971 ± 0.0099	
40.66	23.70 ± 0.30	0.165 ± 0.012		38.11	40.07 ± 0.42	0.144 ± 0.014		38.09	35.89 ± 0.27	0.1277 ± 0.011	
43.20	12.33 ± 0.17	0.409 ± 0.015		39.12	29.73 ± 0.35	0.145 ± 0.017		39.10	26.63 ± 0.23	0.140 ± 0.013	
45.73	9.678 ± 0.13	0.854 ± 0.015		40.64	19.45 ± 0.28	0.202 ± 0.020		40.62	16.64 ± 0.16	0.247 ± 0.014	
48.26	10.05 ± 0.15	0.982 ± 0.017		43.17	10.85 ± 0.17	0.574 ± 0.022		43.15	10.29 ± 0.12	0.704 ± 0.018	
50.79	10.19 ± 0.14	0.936 ± 0.015		45.70	10.07 ± 0.14	0.939 ± 0.020		44.67	10.11 ± 0.12	0.909 ± 0.018	
53.32	9.07 ± 0.12	0.824 ± 0.015		47.22	10.40 ± 0.13	0.994 ± 0.018		45.68	10.40 ± 0.11	0.985 ± 0.017	
55.84	7.005 ± 0.102	0.728 ± 0.016		48.23	10.89 ± 0.14	0.952 ± 0.018		46.69	10.99 ± 0.11	1.007 ± 0.017	
58.37	4.845 ± 0.073	0.615 ± 0.017		50.76	10.76 ± 0.13	0.889 ± 0.018		48.21	11.74 ± 0.12	0.953 ± 0.017	
60.89	2.985 ± 0.053	0.579 ± 0.021		53.29	8.875 ± 0.11	0.793 ± 0.018		50.74	11.16 ± 0.11	0.829 ± 0.017	
63.41	1.936 ± 0.031	0.547 ± 0.019		55.82	6.488 ± 0.093	0.679 ± 0.021		53.26	9.043 ± 0.094	0.754 ± 0.016	
65.93	1.604 ± 0.028	0.672 ± 0.020		58.34	4.055 ± 0.057	0.591 ± 0.020		55.79	5.973 ± 0.066	0.612 ± 0.017	
68.45	1.495 ± 0.025	0.799 ± 0.019		60.86	2.470 ± 0.036	0.580 ± 0.021		58.31	3.674 ± 0.042	0.567 ± 0.017	
70.97	1.425 ± 0.022	0.877 ± 0.018		63.38	1.777 ± 0.030	0.617 ± 0.024		60.84	2.238 ± 0.025	0.549 ± 0.017	
				65.90	1.515 ± 0.024	0.729 ± 0.022		63.35	1.767 ± 0.022	0.674 ± 0.019	
				68.42	1.500 ± 0.024	0.863 ± 0.022		65.87	1.673 ± 0.018	0.859 ± 0.017	
								68.39	1.636 ± 0.022	0.913 ± 0.019	
								70.91	1.526 ± 0.021	0.935 ± 0.020	

^{100}Mo				^{144}Sm				^{144}Sm (continued)			
Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power		Angle	Cross Section (mb/sr)	Analyzing Power	
14.15	2863.1 ± 26.9	-0.110 ± 0.015		13.60	4245.2 ± 18.5	-0.1488 ± 0.0089		49.32	5.866 ± 0.081	0.385 ± 0.020	
16.17	974.9 ± 9.1	-0.335 ± 0.015		14.10	3283.5 ± 16.3	-0.1895 ± 0.0077		50.33	5.998 ± 0.071	0.707 ± 0.017	
18.19	183.7 ± 2.2	-0.695 ± 0.018		15.11	1728.1 ± 11.8	-0.270 ± 0.011		51.33	6.789 ± 0.075	0.945 ± 0.017	
19.20	57.55 ± 0.86	-0.778 ± 0.022		16.12	800.9 ± 5.7	-0.416 ± 0.012		52.34	7.992 ± 0.082	0.981 ± 0.016	
20.21	40.82 ± 0.41	0.603 ± 0.016		17.13	315.8 ± 2.5	-0.541 ± 0.013		54.35	9.695 ± 0.090	0.909 ± 0.015	
21.22	86.52 ± 1.05	0.886 ± 0.018		18.13	146.0 ± 1.3	-0.100 ± 0.013		56.35	9.564 ± 0.080	0.767 ± 0.014	
22.23	153.96 ± 1.56	0.697 ± 0.016		19.14	172.3 ± 1.5	0.530 ± 0.014		58.36	7.601 ± 0.071	0.640 ± 0.014	
24.25	278.85 ± 2.43	0.444 ± 0.014		19.64	214.1 ± 2.4	0.588 ± 0.016		60.37	5.111 ± 0.059	0.433 ± 0.017	
26.27	315.50 ± 2.37	0.311 ± 0.012		20.15	282.0 ± 2.4	0.580 ± 0.014		62.38	3.043 ± 0.036	0.294 ± 0.020	
26.27	315.14 ± 2.82	0.294 ± 0.014		21.15	406.4 ± 3.3	0.454 ± 0.013		64.38	1.931 ± 0.033	0.270 ± 0.027	
27.79	284.24 ± 2.74	0.232 ± 0.015		22.16	502.3 ± 3.2	0.378 ± 0.010		66.39	1.699 ± 0.027	0.556 ± 0.022	
28.29	270.79 ± 2.39	0.201 ± 0.014		23.17	562.4 ± 4.8	0.267 ± 0.013		68.40	1.904 ± 0.023	0.899 ± 0.017	
30.31	179.64 ± 1.78	0.076 ± 0.016		24.17	549.7 ± 4.7	0.192 ± 0.013		70.40	2.086 ± 0.024	0.971 ± 0.017	
32.83	78.22 ± 0.84	-0.079 ± 0.017		25.18	524.4 ± 4.6	0.127 ± 0.013		72.41	1.972 ± 0.024	0.978 ± 0.017	
34.35	40.63 ± 0.43	-0.170 ± 0.017		26.19	444.3 ± 3.5	0.064 ± 0.012		74.41	1.701 ± 0.022	0.965 ± 0.018	
35.35	25.16 ± 0.26	-0.058 ± 0.016		27.20	351.3 ± 3.8	-0.024 ± 0.016					
36.36	17.49 ± 0.22	0.237 ± 0.020		28.20	263.2 ± 1.9	-0.110 ± 0.010					
37.88	15.32 ± 0.19	0.814 ± 0.019		29.21	177.3 ± 1.5	-0.213 ± 0.013					
39.39	19.43 ± 0.22	0.947 ± 0.018		30.21	107.86 ± 0.82	-0.340 ± 0.012					
40.40	23.18 ± 0.25	0.921 ± 0.017		31.22	58.02 ± 0.50	-0.508 ± 0.013					
42.92	29.53 ± 0.34	0.708 ± 0.018		31.72	41.69 ± 0.38	-0.566 ± 0.014					
45.44	25.96 ± 0.26	0.535 ± 0.016		32.23	29.55 ± 0.36	-0.546 ± 0.018					
50.47	8.918 ± 0.109	0.256 ± 0.019		33.23	17.08 ± 0.19	-0.061 ± 0.016					
55.51	3.459 ± 0.068	0.776 ± 0.026		34.24	17.60 ± 0.25	0.779 ± 0.020					
				35.25	25.74 ± 0.23	0.990 ± 0.015					
				36.25	37.06 ± 0.45	0.889 ± 0.017					
				38.26	56.53 ± 0.55	0.645 ± 0.015					
				40.28	59.73 ± 0.31	0.4482 ± 0.0089					
				42.29	46.77 ± 0.28	0.2746 ± 0.0093					
				43.79	34.05 ± 0.24	0.137 ± 0.010					
				44.30	29.10 ± 0.18	0.0826 ± 0.0091					
				46.31	14.57 ± 0.13	-0.110 ± 0.013					
				47.31	9.816 ± 0.091	-0.151 ± 0.014					
				48.32	6.837 ± 0.087	0.015 ± 0.019					

^{208}Pb			^{209}Bi		
Angle	Cross Section (mb/sr)	Analyzing Power	Angle	Cross Section (mb/sr)	Analyzing Power
13.17	6644.7 ± 9.3	-0.1224 ± 0.0027	13.33	6329.9 ± 12.3	-0.1249 ± 0.0033
14.17	3484.6 ± 6.8	-0.1648 ± 0.0037	14.34	3364.9 ± 8.9	-0.1646 ± 0.0045
15.18	1795.0 ± 4.9	-0.1426 ± 0.0044	14.84	2462.5 ± 7.6	-0.1613 ± 0.0051
16.18	1034.3 ± 3.7	0.0025 ± 0.0052	15.18	1950.7 ± 6.8	-0.1591 ± 0.0052
17.19	884.0 ± 3.4	0.2444 ± 0.0066	15.34	1776.9 ± 6.5	-0.1482 ± 0.0057
18.19	966.9 ± 3.6	0.3079 ± 0.0068	15.85	1272.9 ± 5.5	-0.0955 ± 0.0064
19.20	1116.9 ± 3.8	0.2867 ± 0.0064	16.18	1088.5 ± 3.6	-0.0302 ± 0.0045
20.20	1222.1 ± 4.0	0.2325 ± 0.0057	16.35	1034.6 ± 4.9	0.0084 ± 0.0070
21.21	1207.0 ± 4.0	0.1639 ± 0.0053	16.68	913.0 ± 3.3	0.0779 ± 0.0050
22.21	1106.3 ± 3.8	0.0910 ± 0.0051	17.19	845.0 ± 3.2	0.1971 ± 0.0057
23.22	929.8 ± 3.5	0.0166 ± 0.0054	17.35	851.8 ± 4.5	0.2250 ± 0.0083
24.22	709.2 ± 3.0	-0.0590 ± 0.0062	17.69	853.8 ± 3.2	0.2905 ± 0.0063
25.23	492.3 ± 1.8	-0.1751 ± 0.0057	18.19	896.9 ± 3.3	0.3066 ± 0.0061
26.23	301.2 ± 1.4	-0.3384 ± 0.0081	18.69	969.6 ± 3.4	0.3071 ± 0.0061
27.24	156.8 ± 1.0	-0.531 ± 0.011	19.20	1036.1 ± 3.5	0.2950 ± 0.0058
28.24	66.59 ± 0.47	-0.850 ± 0.013	20.20	1138.3 ± 3.7	0.2468 ± 0.0054
28.74	43.14 ± 0.38	-0.925 ± 0.015	20.37	1170.5 ± 5.3	0.2320 ± 0.0073
29.24	30.01 ± 0.31	-0.727 ± 0.016	22.21	1068.2 ± 3.6	0.1070 ± 0.0049
30.25	28.36 ± 0.31	0.502 ± 0.016	24.22	701.1 ± 1.4	-0.0552 ± 0.0030
30.75	37.42 ± 0.31	0.762 ± 0.014	26.23	309.08 ± 0.96	-0.3173 ± 0.0056
31.25	51.92 ± 0.37	0.831 ± 0.013	27.40	148.84 ± 1.19	-0.560 ± 0.013
31.76	66.78 ± 0.66	0.811 ± 0.016	28.24	75.20 ± 0.39	-0.811 ± 0.011
32.26	80.25 ± 0.46	0.741 ± 0.013	28.91	41.52 ± 0.57	-0.908 ± 0.019
32.76	99.68 ± 0.93	0.693 ± 0.015	29.24	31.28 ± 0.25	-0.832 ± 0.013
34.27	128.36 ± 0.71	0.484 ± 0.010	30.25	25.17 ± 0.22	0.271 ± 0.013
36.27	127.52 ± 0.71	0.2794 ± 0.0088	30.41	27.32 ± 0.51	0.433 ± 0.027
38.28	87.84 ± 0.59	0.051 ± 0.010	30.92	35.72 ± 0.46	0.806 ± 0.019
39.29	63.59 ± 0.50	-0.083 ± 0.011	31.42	48.72 ± 0.48	0.819 ± 0.015
40.29	43.00 ± 0.41	-0.264 ± 0.015	31.92	64.34 ± 0.62	0.797 ± 0.015
41.29	24.33 ± 0.22	-0.462 ± 0.013	32.76	88.27 ± 0.21	0.6865 ± 0.0084
42.30	13.55 ± 0.15	-0.577 ± 0.017	35.27	128.16 ± 0.36	0.3931 ± 0.0059
43.30	8.479 ± 0.130	-0.313 ± 0.022	35.43	131.64 ± 1.11	0.379 ± 0.013
44.31	8.113 ± 0.090	0.488 ± 0.016	37.78	98.34 ± 0.70	0.1175 ± 0.0098
45.31	11.192 ± 0.122	0.912 ± 0.016	41.96	17.60 ± 0.22	-0.593 ± 0.018
46.31	15.921 ± 0.159	0.958 ± 0.017	42.46	13.32 ± 0.20	-0.659 ± 0.021
47.32	19.987 ± 0.199	0.872 ± 0.016	42.80	11.09 ± 0.13	-0.546 ± 0.017
48.32	22.723 ± 0.212	0.750 ± 0.016	43.47	8.146 ± 0.152	-0.295 ± 0.027
50.33	23.864 ± 0.218	0.569 ± 0.015	44.30	7.855 ± 0.138	0.386 ± 0.024
52.33	17.468 ± 0.167	0.333 ± 0.014	45.31	10.471 ± 0.101	0.857 ± 0.014
54.34	10.042 ± 0.142	0.0082 ± 0.205	46.48	15.32 ± 0.21	0.931 ± 0.019
55.34	6.627 ± 0.081	-0.145 ± 0.018	47.82	20.39 ± 0.22	0.833 ± 0.016
56.35	4.505 ± 0.060	-0.265 ± 0.019	50.33	23.19 ± 0.17	0.557 ± 0.011
57.35	3.202 ± 0.056	-0.134 ± 0.026	52.83	15.54 ± 0.16	0.276 ± 0.014
58.35	2.992 ± 0.055	0.178 ± 0.026	55.34	6.922 ± 0.083	-0.145 ± 0.017
60.36	3.705 ± 0.061	0.856 ± 0.021	56.51	4.374 ± 0.091	-0.288 ± 0.030
61.36	4.463 ± 0.067	0.960 ± 0.020	57.85	2.965 ± 0.032	-0.043 ± 0.015
62.36	4.926 ± 0.070	0.974 ± 0.019	60.36	2.628 ± 0.060	0.849 ± 0.021
64.37	5.272 ± 0.092	0.875 ± 0.022	61.52	4.388 ± 0.083	0.952 ± 0.023
66.37	4.325 ± 0.066	0.660 ± 0.022	62.86	4.943 ± 0.078	0.935 ± 0.020
68.38	2.814 ± 0.047	0.468 ± 0.024	65.37	4.773 ± 0.058	0.813 ± 0.016
70.38	1.516 ± 0.035	0.212 ± 0.033	67.87	3.065 ± 0.050	0.506 ± 0.022
			70.38	1.531 ± 0.031	0.104 ± 0.028
			72.88	0.935 ± 0.018	0.136 ± 0.027
			75.38	1.082 ± 0.016	0.667 ± 0.020

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