Further Investigation of α +¹²C and α +¹⁶O Elastic Scattering

Sh. Hamada

Abstract—The current work aims to study the rainbow likestructure observed in the elastic scattering of alpha particles on both ¹²C and ¹⁶O nuclei. We reanalyzed the experimental elastic scattering angular distributions data for α +¹²C and α +¹⁶O nuclear systems at different energies using both optical model and double folding potential of different interaction models such as: CDM3Y1, DDM3Y1, CDM3Y6 and BDM3Y1. Potential created by BDM3Y1 interaction model has the shallowest depth which reflects the necessity to use higher renormalization factor (*N_r*). Both optical model and double folding potential of different interaction models fairly reproduce the experimental data.

Keywords—Nuclear rainbow, elastic scattering, optical model, double folding, density distribution

PACS number(s)-21.60.Gx, 24.10.Ht.

I. INTRODUCTION

CTUDY of nuclear interactions is of special interest as it Ocould provide us with useful information about nuclear structure and mechanism of interaction. One of the most interesting features which could be observed in study of nuclear reactions is nuclear rainbow phenomenon. This phenomenon was first realized in the scattering of alpha particles [1]. Similar effects were seen in the scattering of the light heavy ions 6Li [2], [3], in the elastic scattering of some nuclear system such as ${}^{16}O+{}^{12}C$ [4]-[6], ${}^{12}C+{}^{12}C$ [7]-[12], $^{16}\text{O}+^{16}\text{O}$ [13], [14] and also in ³He and α -nucleus elastic and inelastic scattering [15]-[18]. The elastic scattering angular distributions for these nuclear systems reveal refractive features, such as rainbow scattering patterns and broad interference minima "Airy minima" [19]. Nuclei are well known to have wave properties and they may suffer from being diffracted, refracted and also from interference. Consequently, the nucleus-nucleus scattering may display rainbow features depending on: The scattering conditions, energy of the incident projectile and the binding structure of the projectile and target. The rainbow could be observed well in the elastic scattering whereas there is no change in the total flux. Therefore, the nuclear rainbow should be strongest in the elastic scattering channel especially in nuclear systems with small absorption.

In the present work, we reanalyzed both $\alpha^{+12}C$ and $\alpha^{+16}O$ elastic scattering angular distributions using both optical model (OM) and double folding potential (DFP) of different interaction models such as: CDM3Y1, DDM3Y1, CDM3Y6

and BDM3Y1. All these different interaction models are essentially based on the basic form of M3Y Reid or M3Y Paris, but the main difference between them lies in the values of the parameters used in calculating density dependent function $F(\rho)$. We have extracted the renormalization factor (N_r) for $(\alpha^{+12}C \text{ and } \alpha^{+16}O)$ nuclear systems at the different concerned energies within the framework of the different concerned interaction models.

In the previous studies for α +¹⁶O, Michel et al. [20] presented a systematic OM analysis for ${}^{16}O(\alpha, \alpha){}^{16}O$ elastic scattering with a new parametrization for the real part of the potential, whereas for the imaginary part they used a squared Woods-Saxon form factor. The extracted potential in Michel's work gives a precise description of α^{+16} O elastic scattering data in the energy range between 30 and 150 MeV. H. Abele et al. [21] analyzed α +¹⁶O elastic scattering data using an early version for DFP and the renormalization factor allowed to be varied from 1.395 at the lowest energy down to 1.275 at the highest energy. M. El-Azab Farid [15] used the single folding cluster (SFC) [22] and double folding cluster (DFC) [23], [24] models to analyze the experimental data for α +¹⁶O elastic scattering and he obtained a reasonable fitting. H. Abele et al. [21] analyzed the α +¹²C elastic scattering, the real part of the potential was calculated using DFP with DDM3Y model of NN interaction and the imaginary potential was expressed in terms of a Fourier-Bessel function of six terms. Khallaf et al. [25] analyzed this nuclear system using the JLM NN interaction.

In the current work, we tried to analyze α +¹²C and α +¹⁶O elastic scattering data using OM and DFP of different models of interaction and to extract N_r from the different concerned interaction models. Based on the theoretical calculation, rainbow phenomenon and Airy minima are well reproduced and presented.

The rest of the paper is organized as follows. In Section II the theoretical analysis of the experimental data is presented and Section III is devoted to the results and discussion. The summary is given in Section IV.

II. THEORETICAL ANALYSIS

A. Elastic Scattering in the Framework of OM

The elastic scattering angular distributions for both α +¹⁶C and α +¹⁶O are firstly reanalyzed within the framework of the standard OM of the nucleus. The interaction potential could be written as:

Sh. Hamada is with Faculty of Science, Tanta University, Tanta, Egypt (e-mail: sh.m.hamada@science.tanta.edu.eg).

v

$$U(\mathbf{r}) = V_{\rm C} - V_0 \left[1 + \exp\left(\frac{r - R_{\rm V}}{a_{\rm V}}\right) \right]^{-1} - iW_0 \left[1 + \exp\left(\frac{r - R_{\rm W}}{a_{\rm W}}\right) \right]^{-1}$$
(1)

with radius:

$$R_i = r_i (A_T^{1/3}), i = V, W, C$$

This potential consists of three parts: a- Real volume part, b-Imaginary volume part and c- Coulomb potential of a uniform charged sphere. Real and imaginary parts of the potential are taken to be of Woods-Saxon form factor.

In OM calculations, the differential cross section could be easily calculated as the square of scattering amplitude $\frac{d\sigma}{d\Omega} = |f(\theta, \varphi)|^2$, and the elastic nucleus-nucleus scattering

amplitude can be expressed in the form:

$$f(\theta) = f_c(\theta) + \frac{i}{2k} \sum_{l} (2l+1) \exp(2i\sigma_l) (1-S_l) P_l(\cos\theta)$$
(2)

where $f_c(\theta)$ is the amplitude of the Coulomb scattering, σ_l is the Coulomb phase shift, k is the wave number, S_l is the scattering matrix element for the *l*-th partial wave, and $P_l(\cos\theta)$ is the Legendre polynomial.

It is frequently found that many sets of parameters give good fits to the data, and the question then arises whether any one of these sets is more physical than the others, and if so which is to be preferred. These parameter ambiguities are of two main types, discrete and continuous. Discrete ambiguities refer to regions of parameter space that give acceptable fits separated by unacceptable regions. Continuous ambiguities refer to combinations of parameters that may be simultaneously varied being subjected to some constraint without significantly affecting the fit. The existence of these and other more complicated parameter ambiguities means that it is not possible to establish the optical potential (OP) by phenomenological analyses alone. It is necessary to start by constraining the potential as closely as possible by physical requirements before parameter optimization. So, it is preferable to use more microscopic potential such as folding potential.

B. Elastic Scattering in the Framework of Double Folding

The DFP for $\alpha^{+12}C$ and $\alpha^{+16}O$ at the different concerned energies are prepared using code DFMSPH [26]. In the double folding (DF) calculations, the real part of the potential is calculated from a more fundamental basis by the DF method in which the nucleon-nucleon interaction potential is folded into the densities of both the projectile and target nuclei:

$$V^{DF}(\vec{R}) = \iint \rho_{p}(\vec{r}_{1})\rho_{t}(\vec{r}_{2})V_{NN}(\left|\vec{R}+\vec{r}_{2}-\vec{r}_{1}\right|)d^{3}\vec{r}_{1}d^{3}\vec{r}_{2}$$
(3)

where $\rho_p(\vec{r}_1)$ and $\rho_t(\vec{r}_2)$ are the nuclear matter density distributions of both the projectile and target nuclei

respectively, and $V_{NN}(\vec{R} + \vec{r_2} - \vec{r_1})$ is the effective NN interaction potential which gives the corresponding values of the nuclear incompressibility K in the Hartree-Fock (HF) calculation of nuclear matter and it was taken to be of the CDM3Y1 and CDM3Y6 form based on the M3Y-Paris potential:

$$v_{\rm D}(s) = 11061.625 \frac{\exp(-4s)}{4s} - 2537.5 \frac{\exp(-2.5s)}{2.5s},$$

$$E_{\rm EX}(s) = -1524.25 \frac{\exp(-4s)}{4s} - 518.75 \frac{\exp(-2.5s)}{2.5s} - 7.8474 \frac{\exp(-0.7072s)}{0.7072s}$$
(4)

and also BDM3Y1 and DDM3Y1 form based on the M3Y-Reid potential:

$$v_{\rm D}(s) = 7999.0 \frac{\exp(-4s)}{4s} - 2134.25 \frac{\exp(-2.5s)}{2.5s},$$

$$v_{\rm EX}(s) = -4631.38 \frac{\exp(-4s)}{4s} - 1787.13 \frac{\exp(-2.5s)}{2.5s} - 7.8474 \frac{\exp(-0.7072s)}{0.7072s}$$
(5)

The M3Y-Paris and M3Y- Reid interactions are scaled by an explicit density-dependent function $F(\rho)$:

$$v_{\rm D(EX)}(\rho, s) = F(\rho)v_{\rm D(EX)}(s), \qquad S = \left| \vec{R} + \vec{r}_2 - \vec{r}_1 \right|$$
(6)

 $v_{\rm D}$ and $v_{\rm EX}$ are the direct and exchange components of the M3Y-Paris and M3Y- Reid, ρ is the nuclear matter density. The density-dependent function is expressed as:

$$F(\rho) = C[1 + \alpha \exp(-\beta \rho) - \gamma \rho^{n}], \qquad (7)$$

The parameters C, α , β , γ , n for the different concerned interaction models listed in Table I were taken from [27],

| TABLEI | | | | | | | | | |
|--|--------|--------|----------------------------|--------------------------|---|------------|--|--|--|
| PARAMETERS OF DENSITY DEPENDENCE FUNCTION $F(P)$ | | | | | | | | | |
| Interaction Model | С | α | β (fm ³) | $\gamma ({\rm fm}^{3n})$ | n | K (MeV) | | | |
| CDM3Y6 | 0.2658 | 3.8033 | 1.4099 | 4.0 | 1 | 252 | | | |
| CDM3Y1 | 0.3429 | 3.0232 | 3.5512 | 0.5 | 1 | 188 | | | |
| BDM3Y1 | 1.2253 | 0.0 | 0.0 | 1.5124 | 1 | 232 | | | |
| DDM3Y1 | 0.2845 | 3.6391 | 2.9605 | 0.0 | 0 | 171 | | | |

The density distributions of α , ¹²C and ¹⁶O are expressed in modified form of the Gaussian shape as: a $\rho(r) = \rho_0 (1 + wr^2) \exp(-\beta r^2)$, the parameters (ρ_0, w, β) for α particles, ¹²C and ¹⁶O are listed in Table II.

III. RESULTS AND DISCUSSION

A. $\alpha + {}^{12}C$ Nuclear System

Both the phenomenological (OM) and semi microscopic (DF) calculations are performed using code FRESCO and code SFRESCO [30]. The comparisons between the experimental data for α +¹²C nuclear system at energies (104, 120, 139, 145, 166 and 172.5 MeV) [31]-[35] and the theoretical calculations within the framework of OM are shown in Fig. 1. The radius parameter for real volume part of potential (r_V) was fixed at 1.245 fm, radius parameter for the imaginary volume part of potential (r_W) was fixed at 1.6 fm, and radius parameter for the Coulomb part of potential (r_C) was fixed at 1.25 fm. The optimal OP parameters are listed in Table II.

Table II. As shown in Fig. 1, the most significant contribution for the cross section comes from the farside component. The farside components at energies 104, 120, 139, 145, 166 and 172.5 MeV predicted the position of Airy minimum to be around 24.8°, 23.7°, 23.2°, 21.3°, 20.9° and 19.8° respectively. The significant increase in cross section at backward angles at energy 104 MeV might indicate that other non-scattering events, like the elastic alpha cluster transfer process, still contribute to the elastic cross section at this energy.

Nearside and farside components of cross sections are calculated using the same the potential parameters listed in



Fig. 1 The comparison between the experimental data for α +¹²C elastic scattering at energies (104, 120, 139, 145, 166 and 172.5 MeV) and the theoretical calculations using OM



Fig. 2 The variation of reaction cross section with energy

We have calculated the reaction cross section (σ_R) for α +¹²C nuclear system at the different concerned energies using both OM and DFP of different interaction models as shown in Fig.

2. The energy dependence of reaction cross section can be expressed as: $\sigma = 1854 - 2.128$ E.

The quality of the fitting of OM and DF calculations to the experimental data can be estimated using the χ^2 - method, which is represented as:

$$\chi^{2} = \frac{1}{N} \sum_{i=1}^{N} \left[\frac{\sigma_{i}^{t}(\theta) - \sigma_{i}^{e}(\theta)}{\Delta \sigma_{i}^{e}(\theta)} \right]^{2} = \frac{1}{N} \sum_{i=1}^{N} \chi_{i}^{2}$$
(8)

where σ^e and σ^t are the experimental and theoretical differential cross sections of the elastic scattering for *i*-th scattering angle, $\Delta\sigma^e$ is the error of the experimental differential cross sections at these angles, N is the number of measurements. χ^2/N values are listed in Table II, the less the value χ^2 is, the better is the description of the experimental data in terms of the selected theoretical representation.

World Academy of Science, Engineering and Technology International Journal of Chemical and Molecular Engineering Vol:10, No:12, 2016

| TABLE II | | | | | | | | | | |
|---|------------|----------------------|------------------|---------|------------------|--------|-----------------|---------------------------|---------------------------|-------|
| OP AND DF PARAMETERS FOR α-PARTICLES ELASTICALLY SCATTERED ON ¹² C AT DIFFERENT ENERGIES | | | | | | | | | | 2/11 |
| E (MeV) | Model | V _o (MeV) | $a_{\rm v}$ (fm) | W₀(MeV) | $a_{\rm w}$ (fm) | Nr | σ_R (mb) | Jv (MeV.fm ³) | Jw (MeV.fm ³) | χ²/Ν |
| 172.5 | OM | 90.38 | 0.772 | 22.02 | 0.663 | | 1510 | 315.4 | 125 | 3.02 |
| | DF- CDM3Y1 | | | 22.02 | 0.663 | 1.111 | 1506 | 287.5 | 125 | 1.74 |
| | DF- DDM3Y1 | | | 22.02 | 0.663 | 1.166 | 1505 | 282.0 | 125 | 1.51 |
| | DF- CDM3Y6 | | | 22.02 | 0.663 | 1.181 | 1513 | 287.38 | 125 | 1.54 |
| | DF-BDM3Y1 | | | 22.02 | 0.663 | 1.238 | 1512 | 268.56 | 125 | 1.51 |
| 166 | OM | 90.43 | 0.745 | 21.29 | 0.6347 | | 1460 | 306.38 | 118.38 | 16.11 |
| | DF- CDM3Y1 | | | 21.29 | 0.6347 | 1.118 | 1498 | 305.08 | 118.38 | 10.28 |
| | DF-DDM3Y1 | | | 21.29 | 0.6347 | 1.172 | 1495 | 285.94 | 118.38 | 8.97 |
| | DF- CDM3Y6 | | | 21.29 | 0.6347 | 1.1804 | 1505 | 290.8 | 118.38 | 9.97 |
| | DF-BDM3Y1 | | | 21.29 | 0.6347 | 1.236 | 1502 | 272.7 | 118.38 | 9.99 |
| 145 | OM | 92.77 | 0.772 | 21.14 | 0.69 | | 1581 | 323.7 | 122.4 | 4.16 |
| | DF- CDM3Y1 | | | 21.14 | 0.69 | 1.0827 | 1570 | 317.6 | 122.4 | 2.32 |
| | DF-DDM3Y1 | | | 21.14 | 0.69 | 1.1352 | 1566 | 300.28 | 122.4 | 2.43 |
| | DF- CDM3Y6 | | | 21.14 | 0.69 | 1.137 | 1577 | 303.02 | 122.4 | 2.25 |
| | DF-BDM3Y1 | | | 21.14 | 0.69 | 1.191 | 1573 | 286.6 | 122.4 | 2.45 |
| 139 | OM | 93.65 | 0.712 | 25.68 | 0.626 | | 1465 | 306.1 | 129.8 | 2.89 |
| | DF- CDM3Y1 | | | 25.68 | 0.493 | 1.041 | 1459 | 321.5 | 129.8 | 6.62 |
| | DF-DDM3Y1 | | | 25.68 | 0.493 | 1.08 | 1459 | 304.56 | 129.8 | 6.64 |
| | DF- CDM3Y6 | | | 25.68 | 0.493 | 1.099 | 1461 | 306.6 | 129.8 | 7.05 |
| | DF-BDM3Y1 | | | 25.68 | 0.493 | 1.1366 | 1460 | 290.73 | 129.8 | 7.24 |
| 120 | OM | 96.12 | 0.772 | 20.98 | 0.673 | | 1631 | 335.4 | 119.9 | 3.39 |
| | DF- CDM3Y1 | | | 20.98 | 0.673 | 1.0574 | 1604 | 333.56 | 119.9 | 2.39 |
| | DF-DDM3Y1 | | | 20.98 | 0.673 | 1.0917 | 1597 | 318.44 | 119.9 | 2.96 |
| | DF- CDM3Y6 | | | 20.98 | 0.673 | 1.1068 | 1610 | 318.21 | 119.9 | 2.58 |
| | DF-BDM3Y1 | | | 20.98 | 0.673 | 1.1413 | 1603 | 304.17 | 119.9 | 3.21 |
| 104 | OM | 100.6 | 0.769 | 19.55 | 0.694 | | 1689 | 349.9 | 113.5 | 2.72 |
| | DF- CDM3Y1 | | | 19.55 | 0.694 | 1.0645 | 1652 | 344.08 | 113.5 | 2.44 |
| | DF-DDM3Y1 | | | 19.55 | 0.694 | 1.0976 | 1645 | 330.7 | 113.5 | 2.45 |
| | DF- CDM3Y6 | | | 19.55 | 0.694 | 1.1193 | 1660 | 328.34 | 113.5 | 2.43 |
| | DF-BDM3Y1 | | | 19.55 | 0.694 | 1.1413 | 1645 | 315.98 | 113.5 | 1.38 |



Fig. 3 Energy dependence of real (Jv) and imaginary (Jw) volume integral

OP and DFP can conveniently be characterized by the real and imaginary volume integrals per nucleon:

$$Jv = -\frac{4\pi}{A_P A_t} \int_0^\infty V(r) r^2 dr , \quad J_w = -\frac{4\pi}{A_P A_t} \int_0^\infty W(r) r^2 dr$$
(9)

The J_{ν} values in our work ranged between 268.56 and 335.4 MeV.fm³ and J_{ν} between 113.5 and 129.8 MeV.fm³ as listed in Table II which agree well with the previously existing theoretical estimations [15]. The energy dependence of real volume integral in this work could be expressed as $J_{\nu} = 399.5$ -0.543E, and for imaginary volume integral $J_{\nu} = 103.84$ + 0.109E as shown in Fig. 3.

The dependence of V₀ and W₀ values for ${}^{12}C(\alpha, \alpha){}^{12}C$ on energy are shown in Figs. 4 (a) and (b) and can be approximated by:

(a)
$$V = 113.9 - 0.14E$$
,

(b) W = 17.16+0.027E.

The comparison between the experimental data for α +¹²C at energies (104, 120, 139, 145, 166 and 172.5 MeV) and the DF calculations of different interaction models: CDM3Y1, DDM3Y1, CDM3Y6 and BDM3Y1 are shown in Figs. 5, 6, 7 and 8 respectively. Optimal potential parameters for α +¹²C nuclear system, also those from DF of different interaction models are listed in Table II. The obtained renormalization factor (*N_r*) for α +¹²C nuclear system is in the range 1.041-1.238. The optimal parameters for the imaginary part of potential obtained from OM were kept constant during the DF calculations. The interaction potential in this case has the shape:

World Academy of Science, Engineering and Technology International Journal of Chemical and Molecular Engineering Vol:10, No:12, 2016

$$U(R) = V_{\rm C} + Nr V^{DF}(R) - iW(R)$$
⁽¹⁰⁾



Fig. 4 (a) The relationship between real potential depth and energy



Fig. 4 (b) The relationship between imaginary potential depth and energy



Fig. 5 The comparison between the experimental data for α +¹²C at energies (104, 120, 139, 145, 166 and 172.5 MeV) and the theoretical calculations using DFP of interaction model CDM3Y1



Fig. 6 The same as Fig. 5 but for interaction model DDM3Y1



Fig. 7 The same as Fig. 5 but for interaction model CDM3Y6



Fig. 8 The same as Fig. 5 but for interaction model BDM3Y1

The Airy minimum could be observed for ${}^{12}C(\alpha,\alpha){}^{12}C$ elastic scattering at the aforementioned energies. The characteristic features of the falloff of the cross section beyond the rainbow angle in the experimental angular distributions are

well reproduced. As shown in Fig. 1, with increasing the energy of the incident projectile, the position of rainbow angle is shifted toward small angles. The angular position of rainbow angle at the different concerned energies is shown in Fig. 9.

As we discussed before, the calculations in this work are performed using different four models of interaction: CDM3Y1, DDM3Y1, CDM3Y6 and BDM3Y1. The used DFP consists of two parts direct part and exchange part. Fig. 10 shows the variation of the potential depth with radius for the direct and exchange parts of the potential and also their sum (direct + exchange) at E=172.5 MeV, this figure also shows variation of the potential depth with radius at the different concerned alpha particles energies. Fig. 11 shows how the potential is created by BDM3Y1 model of interaction is shallower in comparison with the rest which reflect the necessity to use higher renormalization factor.



Fig. 9 Angular position of the Airy minimum for α +¹²C elastic scattering at energies 104, 120, 139, 145, 166 and 172.5 MeV. The lines are only to guide the eye



Fig. 10 (a) The variation of the potential depth with radius for both direct and exchange parts of the potential and also their sum (direct + exchange) at E=172.5 MeV; (b) the variation of the potential depth with radius at different energies

B. α +¹⁶*O* Nuclear System

The comparison between the experimental data for α +¹⁶O nuclear system at energies 49.5, 69.5, 80.7 and 104 MeV [20], [33], [36] and the theoretical calculations within the framework of OM is shown in Fig. 12. The radius parameter for real volume part of potential (r_V) was fixed at 1.36 fm, radius parameter for imaginary volume part of potential (r_W)

was fixed at 1.73 fm and radius parameter for the Coulomb part of potential (r_c) was fixed at 1.25 fm.

The comparisons between the experimental data for α +¹⁶O at energies 49.5, 69.5, 80.7 and 104 MeV and the DF calculations of different interaction models: CDM3Y1, DDM3Y1, CDM3Y6 and BDM3Y1 are shown in Figs. 13, 14, 15 and 16 respectively. Optimal OP parameters for α +¹⁶O nuclear system, also those from DF of different interaction models are listed in Table III. The obtained N_r for α +¹⁶O

nuclear system is in the range of 1.1157-1.3312. The imaginary part of the potential was taken in the form of standard Woods-Saxon form factor, the same optimal parameters for the imaginary part of potential obtained from OM calculations were kept constant during the DF calculations, and the used interaction potential has the same shape as in (10).

The nuclear rainbow phenomenon could be clearly $^{16}O(\alpha,\alpha)^{16}O$ observed in elastic scattering at the aforementioned energies. The characteristic features of the falloff of the cross section beyond the rainbow angle in the experimental angular distributions are well reproduced. Both OM and DF of different interaction models: CDM3Y1, DDM3Y1, CDM3Y6 and BDM3Y1 could reasonably reproduce the experimental data. The angular distribution data and the OM calculations at E=49.5 MeV showed that the Airy minimum is \approx 78°, but the DF calculations expected the minimum to be $\approx 75^{\circ}$ and this may be due to the existence of a valley of two minima: the first minimum at 78° as shown in Fig. 12 and the second probably at 70° (not shown in the experimental data). DF calculations expected that the position of Airy minimum is $\approx 48^{\circ}$ at *E*=69.5 MeV as shown in Figs. 13-16, while from the OM calculations it is around 40° as shown in Fig. 12, the minimum at $\approx 48^{\circ}$ could not be observed from experimental data. At energies 80.7 and 104 MeV, both OM and DF calculations expected the same position for Airy minimum (35° at E=80.7 MeV and 31° at E=104 MeV). As shown in Figs. 12-16, with increasing the energy of the incident projectile, the position of rainbow angle is shifted toward small angles. The angular position of rainbow angle at the different concerned energies is shown in Fig. 17.







Fig. 12 The comparison between the experimental data for α +¹⁶O at energies (49.5, 69.5, 80.7 and 104 MeV) and the theoretical calculations using OM

TABLE III

| OP AND DF PARAMETERS FOR α -PARTICLES ELASTICALLY SCATTERED ON ¹⁶ O AT DIFFERENT ENERGIES | | | | | | | | | | |
|---|------------|----------------------|-----------------------|----------------------|------------------|--------|-----------------|---------------------------|---------------------------|------------|
| E (MeV) | Model | V _o (MeV) | $a_{\rm v}({\rm fm})$ | W _o (MeV) | $a_{\rm w}$ (fm) | Nr | σ_R (mb) | Jv (MeV.fm ³) | Jw (MeV.fm ³) | χ^2/N |
| 49.5 | OM | 109.63 | 0.6 | 8.94 | 0.739 | | 1871 | 376.18 | 62.22 | 10.99 |
| | DF- CDM3Y1 | | | 8.94 | 0.739 | 1.1228 | 1872 | 382.18 | 62.22 | 7.6 |
| | DF-DDM3Y1 | | | 8.94 | 0.739 | 1.1474 | 1861 | 375.87 | 62.22 | 7.97 |
| | DF- CDM3Y6 | | | 8.94 | 0.739 | 1.1931 | 1887 | 365.14 | 62.22 | 8.41 |
| | DF-BDM3Y1 | | | 8.94 | 0.739 | 1.2143 | 1874 | 359.03 | 62.22 | 8.3 |
| 69.5 | OM | 99.99 | 0.715 | 15.58 | 0.648 | | 2006 | 376.66 | 102.9 | 5.78 |
| | DF- CDM3Y1 | | | 15.58 | 0.648 | 1.1209 | 1983 | 367.75 | 102.9 | 3.59 |
| | DF-DDM3Y1 | | | 15.58 | 0.648 | 1.1458 | 1973 | 359.0 | 102.9 | 2.33 |
| | DF- CDM3Y6 | | | 15.58 | 0.648 | 1.1782 | 1990 | 350.89 | 102.9 | 3.53 |
| | DF-BDM3Y1 | | | 15.58 | 0.648 | 1.2002 | 1979 | 342.92 | 102.9 | 2.59 |
| 80.7 | OM | 101.26 | 0.714 | 16.7 | 0.556 | | 1923 | 381.13 | 105.08 | 2.74 |
| | DF- CDM3Y1 | | | 16.7 | 0.556 | 1.1307 | 1894 | 359.89 | 105.08 | 2.12 |
| | DF-DDM3Y1 | | | 16.7 | 0.556 | 1.1637 | 1885 | 349.69 | 105.08 | 1.51 |
| | DF- CDM3Y6 | | | 16.7 | 0.556 | 1.191 | 1903 | 343.25 | 105.08 | 1.61 |
| | DF-BDM3Y1 | | | 16.7 | 0.556 | 1.2225 | 1894 | 333.91 | 105.08 | 1.24 |
| 104 | OM | 93.25 | 0.731 | 17.25 | 0.541 | | 1873 | 356.07 | 107.74 | 3.24 |
| | DF- CDM3Y1 | | | 17.25 | 0.541 | 1.1157 | 1823 | 343.9 | 107.74 | 2.82 |
| | DF-DDM3Y1 | | | 17.25 | 0.541 | 1.1717 | 1829 | 327.82 | 107.74 | 2.2 |
| | DF- CDM3Y6 | | | 17.25 | 0.541 | 1.2687 | 1832 | 304.59 | 107.74 | 2.19 |
| | DF-BDM3Y1 | | | 17.25 | 0.541 | 1.3312 | 1837 | 290.02 | 107.74 | 1.79 |

World Academy of Science, Engineering and Technology International Journal of Chemical and Molecular Engineering Vol:10, No:12, 2016



Fig. 13 The comparison between the experimental data for α +¹⁶O at energies (49.5, 69.5, 80.7 and 104 MeV) and the theoretical calculations using DFP of interaction model CDM3Y1



Fig. 14 The same as Fig. 13 but for interaction model DDM3Y1



Fig. 15 The same as Fig. 13 but for interaction model CDM3Y6



Fig. 16 The same as Fig. 13 but for interaction model BDM3Y1



Fig. 17 Angular position of the Airy minimum for α +¹⁶O elastic scattering at energies 49.5, 69.5, 80.7 and 104 MeV. The lines are only to guide the eye



Fig. 18 The variation of reaction cross section with energy

We have calculated the reaction cross section (σ_R) at the different concerned energies using both OM and DF of different interaction models as shown in Fig. 18. The energy

dependence of reaction cross section can be expressed as $\sigma = 1854 - 2.128 \text{ E}$.

The J_{ν} values from both OP and DFP in our work ranged between 290.02 and 382.18 MeV.fm³ and decreased with increasing energy while the $J_{\nu\nu}$ ranges between 62.22 and 107.74 MeV.fm³ and increases with increasing energy as shown in Table III which agree well with the previously existing theoretical estimations [18], [19]. The energy dependence of real volume integral in this work could be expressed as $J_{\nu} = 399.15$ -0.354E, and for imaginary volume integral $J_{\mu\nu} = 34.68$ +0.787E as shown in Fig. 19.



Fig. 19 Energy dependence of real (J_v) and imaginary (J_w) volume integral

The dependence of V₀ (depth for the real volume part of potential) and W₀ (depth for the imaginary volume part of potential) for ¹⁶O(α , α)¹⁶O nuclear system on energy are shown in Figs. 20 (a) and (b) and can be approximated by (a) V = 122.47-0.28E, (b) W = 3.468+0.146E.



Fig. 20 (a) Energy dependence of real potential depth



Fig. 20 (b) Energy dependence of imaginary potential depth

Fig. 21 shows the variation of the potential depth with radius for the direct and exchange parts of the potential and also their sum (Direct + Exchange) for α +¹⁶O nuclear system at E_{α} =49.5 MeV using DF-CDM3Y1. Fig. 22 shows that the potential created by BDM3Y1 model of interaction has the shallowest depth in comparison with the rest, which reflects the necessity to use higher renormalization factor.

IV. SUMMARY

We reanalyzed the available experimental angular distribution data for (α +¹²C and α +¹⁶O) nuclear systems using both OP and DFP of different interaction models. Refractive features and nuclear rainbow phenomenon is presented and well reproduced by theoretical calculations. Analysis of experimental data using DFP of different interaction models: CDM3Y1, CDM3Y6, DDM3Y1 and BDM3Y1 showed that, the potential created by BDM3Y1 model of interaction has the shallowest depth which reflects the necessity to use higher renormalization factor, and the potential created by CDM3Y1 model of interaction is the deepest which reflects the necessity to use smaller renormalization factor. Both (OP) and (DFP) of different interaction models reasonably reproduce the experimental data. The obtained N_r using all the aforementioned interaction models for α +¹²C nuclear system is in the range 1.041 – 1.238. For α +¹⁶O, N_r is in the range 1.1157 - 1.3312. Theoretical calculations showed that DF-CDM3Y1 is the most preferable interaction model in analysis $\alpha^{+12}C$ and $\alpha^{+16}O$ nuclear systems as it gives us the closest values of N_r to unity.

World Academy of Science, Engineering and Technology International Journal of Chemical and Molecular Engineering Vol:10, No:12, 2016



Fig. 21 The variation of the potential depth with radius for both Direct and Exchange parts of the potential and also their sum (Direct + Exchange) at E_{α} =49.5 MeV



Fig. 22 The variation of the potential depth with radius for α +¹⁶O folding potential at E_{α} =49.5 MeV using different interaction models: CDM3Y1, DDM3Y1, CDM3Y6 and BDM3Y1

References

- D. A. Goldberg, S. M. Smith and G. F. Burdzik, Phys. Rev. C 10 (1974) 1362.
- [2] R. M. DeVries, D. A. Goldberg, J. W. Watson, M. S. Zisman, and J. G. Cramer, Phys. Rev. Lett. 39 (1977) 450.
- [3] P. Schwandt, S. Kailas, W. W. Jacobs, M. D. Kaitchuck, W. Ploughe, and P. P. Singh, Phys. Rev. C 21 (1980) 1656.
- [4] A. A. Ogloblin, Yu. A. Glukhov, W. H. Trzaska, A. S. Dem'yanova, S. A. Goncharov, R. Julin, S. V. Klebnikov, M. Mutterer, M. V. Rozhkov, V. P. Rudakov, G. P. Tiorin, Dao T. Khoa, and G. R. Satchler, Phys. Rev. C 62 (2000) 044601.
- [5] M. P. Nicoli, F. Haas, R. M. Freeman, S. Szilner, Z. Basrak, A. Morsad, G. R. Satchler, and M. E. Brandan, Phys. Rev. C 61 (2000) 034609.
- [6] Yu. A. Glukhov, S. A. Goncharov, A. S. Dem'yanova, A. A. Ogloblin, M. V. Rozhkov, V. P. Rudakov, V. Trashka, J. IZV 65 (2001) 647.
- [7] M. Buenerd, P. Martin, R. Bertholet, C. Guet, M. Maurel, J. Mougey, H. Nifenecker, J. Pinston, P. Perrin, F. Schussler, J. Jullen, J. P. Bondorf, L. Carlen, H. Å. Gustafsson, B. Jakobsson, T. Johansson, P. Kristiansson, O. B. Nielsen, A. Oskarsson, I. Otterlund, H. Ryde, B. Schröder, and G. Tibell, Phys. Rev. C 26 (1982) 1299.

- [8] S. Kubono, K. Morita, M.H. Tanaka, M. Sugitani, H. Utsunomiya, H. Yonehara, M.-K. Tanaka, S. Shimoura, E. Takada, M. Fukuda, K. Takimoto, Phys. Lett. B. 127 (1983) 19.
- [9] H. G. Bohlen, X. S. Chen, J. G. Cramer, P. Fröbrich, B. Gebauer, H. Lettau, A. Miczaika, W. von Oertzen, R. Ulrich, T. Wilpert, Z. Phys. A. 322 (1985) 241.
- [10] A. J. Cole, W. D. M. Rae, M. E. Brandan, A. Dacal, B. G. Harvey, R. Legrain, M. J. Murphy, R. G. Stokstad, Phys. Rev. Lett. 47 (1981) 1705.
- [11] M. Buenerd, A. Lounis, J. Chauvin, D. Lebrun, P. Martin, G. Duhamel, J.C. Gondrand, P. De Saintignon, Nucl. Phys. A. 424 (1984) 313.
- [12] J.Y. Hostachy, M. Buenerd, J. Chauvin, D. Lebrun, Ph. Martin, B. Bonin, G. Bruge, J.C. Lugol, L. Papineau, P. Roussel, J. Arvieux, C. Cerruti, Phys. Lett. B. 184 (1987) 139.
- [13] C.C. Sahm, T. Murakami, J.G. Cramer, A.J. Lazzarini, D.D. Leach, D.R. Tieger, R.A. Loveman, W.G. Lynch, M.B. Tsang, J. Van der, Phys. Rev. C 34 (1986) 2165.
- [14] M. E. Brandan, Phys. Rev. Lett. 60 (1988) 784.
- [15] M. El-Azab Farid, Z. M. M. Mahmoud, and G. S. Hassan, Phys. Rev. C 64 (2001) 014310.
- [16] Dao T. Khoa, Phys. Rev. C 63 (2001) 034007.
- [17] Sh. Hamada, Y. Hirabayashi, N. Burtebayev, S Ohkubo, Phys. Rev. C 87, 024311 (2013).

World Academy of Science, Engineering and Technology International Journal of Chemical and Molecular Engineering Vol:10, No:12, 2016

- [18] Y. Hirabayashi and S. Ohkubo, Phys. Rev. C 88 014314 (2013).
- [19] D. A. Goldberg and S. M. Smith, Phys. Rev. Lett. 33 (1974) 715.
- [20] F. Michel, J. Albinski, P. Belery, Th. Delbar, Gh. Grégoire, B. Tasiaux, and G. Reidemeister, Phys. Rev. C 28(1983) 1904.
- [21] H. Abele, H. J. Hauser, A. Körber, W. Leitner, R. Neu, H. Plappert, T. Rohwer, G. Staudt, M. Strasser, S. Weite, and M. Walz, Z. Phys. A 326, 373 (1987).
- [22] F. Michel, Phys. Lett. 60B (1976) 229.
- [23] Z. Majka, H. J. Gils, and H. Rebel, Phys. Rev. C 25 (1982) 2996.
- [24] M. El-Azab Farid, J. Phys. G 16 (1990) 461.
- [25] S. A. E. Khallaf, A. M. Amry, and S. R. Mokhtar, Phys. Rev.C 56 (1997) 2093.
- [26] I. I. Gontchar and M. V. Chushnyakova, Comput. Phys. Commun. 181 (2010) 168.
- [27] Dao T. Khoa, G.R. Satchler, W. von Oertzen, Phys. Rev. C 56 (1997) 954.
- [28] S. Qing-biao, F. Da-chun and Z. Yi-Zhong, Phys. Rev. C 43(1991) 2773.
- [29] G.R. Satchler, W.G. Love, Phys. Rep. 55 (1979) 183.
- [30] I. J. Thompson, Comput. Phys. Rep. 7 (1988) 167.
- [31] B. Tatischeff, I. Brissauud, Nucl. Phys. A. 155 (1970) 89.
- [32] S. Wiktor, C. Mayer-Böricke, A. Kiss, M. Rogge, P. Turek, H. Dabrowski, Acta Phys. Pol. B 12 (1981) 491.
- [33] G. Hauser, R. Löhken, H. Rebel, G. Schatz, G.W. Schweimer, J. Specht, Nucl. Phys. A. 128 (1969) 81.
- [34] M. Yasue, T. Tanabe, F. Soga, J. Kokame, F. Shimokoshi, J. Kasagi, Y. Toba, Y. Kadota, T. Ohsawa, K. Furuno, Nucl. Phys. A. 394 (1983) 14.
- [35] S. M. Smith, G. Tibell, A. A. Cowley, D. A. Goldberg, H. G. Pugh, W. Reichart, N.S. Wall, Nucl. Phys. A. 207 (1973) 273.
- [36] M. Reed thesis, Lawrence Berkeley National Laboratory, Berkeley, CA, USA (1968).