

# THE SEMIMICROSCOPIC DESCRIPTION OF THE SIMPLEST PHOTONUCLEAR REACTIONS WITH GIANT DIPOLE RESONANCE EXCITATION

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The main goal of the presented talk is the further development of the semimicroscopic approach for better description of the simplest photonuclear reactions in the energy region of the giant dipole resonance (GDR). Various versions of the semimicroscopic approach based on the continuum random-phase approximation (cRPA) were intensively used during last years for the studies of these reactions (see, e.g., [1, 2]). Saying of the simplest photonuclear reactions we mean, first of all, the total photoabsorption and the direct or inverse single-particle reactions. Though during these studies the important results have been obtained, all of them are characterized by one essential shortcoming. It appeared that it is impossible to obtain the proper position of the GDR energy in the framework of cRPA in the form adopted in the finite Fermi-system theory (FFST) using the residual Landau-Migdal partial-hole interaction in the following form:

$$F^{ph}(\mathbf{r}, \mathbf{r}') = C\delta(\mathbf{r} - \mathbf{r}') [f(\mathbf{r}) + f'(\mathbf{r})\vec{\tau} \cdot \vec{\tau}'] \quad , \quad (1)$$

where  $f(\mathbf{r})$  and  $f'(\mathbf{r})$  are the dimensionless parameters determining the intensities of the isoscalar and isovector interactions, respectively.

The decision of this problem is prompted by the general principles of the FFST. In the first edition of Migdal book [3] the following relations between the energy ( $\omega_M$ ), the integrated total photoabsorption cross section ( $\sigma_{E1,int}$ ) and parameter  $f'_1$  determining the intensity of momentum-dependent forces have been obtained in a model way:

$$\omega_M^2 = \omega_0^2 \left(1 + \frac{2}{3}f'_1\right) \quad , \quad \sigma_{E1,int} = \sigma_0 \left(1 + \frac{2}{3}f'_1\right) \quad . \quad (2)$$

Here  $\omega_0$  and  $\sigma_0$  are the GDR energy and its integrated cross section, respectively, calculated without account of momentum-dependent forces. As follows from Eq.(2) these forces should play an important role in the GDR formation and its properties.

The first study of the simplest photonuclear reactions in the GDR energy region with the momentum-dependent forces account has been carried out in Ref. [4], where the rather satisfactory description of the experimental photoabsorption and partial ( $n, \gamma$ )-reaction cross sections for some nuclei has been obtained, in particularly, for <sup>208</sup>Pb isotope. However, the calculated energy behaviour of the total photoabsorption cross section for this nucleus is characterized by the feature similar to the resonance peak splitting in the strongly deformed nuclei. Besides, the approach used in Ref. [4] has the certain shortcomings. For instance, introduction of the smearing parameter (the mean doorway-state spreading width)

$$I(\omega) = \alpha(\omega - \Delta)^2 / [1 + (\omega - \Delta)^2 / B^2] \quad (3)$$

(its form is similar to that proposed in Ref. [5]) leads in calculations to the appearance of the additional terms  $\mp(i/2)I(\omega)I(r)$  in the potential  $U(x)$  used for the calculations of the cRPA certain ingredients (Green's functions and continuum-state wave functions). As  $I(r)$  is usually used the Woods-Saxon function  $f_{WS}(r, R^*, a)$ . In the mentioned approach [4] the cutoff radius  $R^*$  was chosen as  $R^* = 2R$  in the calculations of the Green's functions and  $R^* = R$  in the calculations of the partial reaction amplitudes. Such choice of  $R^*$  is not quite justified and makes the used approach rather inconsistent. It is necessary to notice also that the smearing parameter  $I(\omega)$  is an imaginary part of the polarization operator  $\Pi(\omega) = -(i/2)I(\omega) + \text{Re } \Pi(\omega)$ , which determines the relaxation of the particle-hole degree of freedom. The quantity  $\text{Re } \Pi(\omega)$  which can be determined through  $I(\omega)$  with the help of the certain dispersion relation [6] also gives a contribution  $\text{Re } \Pi(\omega)I(r)$  to the potential  $U(x)$ .

Therefore the account of the above made remarks has been chosen as the first step to improve the approach of Ref. [4]. As it is well known in the FFST the effective fields  $V$  satisfy to the following equations (in symbolic form):

$$V = V_0 + F\mathcal{A}V \quad , \quad (4)$$

where  $V_0$  is the external field,  $\mathcal{A}$  is the response function and  $F$  is the residual particle-hole interaction. In the presented approach the isovector part of this interaction is chosen in the form having the separable momentum-dependent part [7]:

$$F(1, 2) \rightarrow \left( F' \delta(\vec{r}_1 - \vec{r}_2) + \frac{k'}{mA} (\vec{p}_1 \vec{p}_2) \right) (\vec{\tau}_1 \vec{\tau}_2) \quad , \quad (5)$$

where  $F' = f' \cdot 300 \text{ MeV fm}^3$ ,  $k'$  is the dimensionless intensity of the momentum-dependent separable forces,  $m$  is the nucleon mass and  $A$  is the number of nucleons. In the case of the GDR excitation the external field  $V_0(x)$  is taken as following:

$$V_0(x) = -\frac{1}{2} r Y_1(\Omega) \tau^{(3)} \quad ,$$

('x' means the set of space, spin and isospin variables). Assuming the operator equality  $\vec{p} = m d\vec{r}/dt$ , in the cRPA from Eqs. (4),(5) the following relations may be obtained after separation of isobaric and spin-angular variables:

$$\tilde{V}(r, \omega) = V(r, \omega) + V_k(r, \omega) \quad , \quad (6)$$

$$V(r, \omega) = r + \frac{2F'}{r^2} \int [\mathcal{A}(r, r', \omega) + \mathcal{A}_k(r, r', \omega)] V(r', \omega) dr' \quad , \quad (7)$$

$$V_k(r, \omega) = \frac{k\omega^2}{1 + k' - \omega^2 k \int r \mathcal{A}(r, r', \omega) r' dr dr'} \cdot r \int r \mathcal{A}(r, r', \omega) V(r', \omega) dr dr' \quad , \quad (8)$$

$$\mathcal{A}_k(r, r', \omega) = \frac{k\omega^2}{1 + k' - \omega^2 k \int r \mathcal{A}(r, r', \omega) r' dr dr'} \int \mathcal{A}(r, r', \omega) r' dr' \int r \mathcal{A}(r, r', \omega) dr \quad . \quad (9)$$

Here  $k = \frac{8\pi m}{3\hbar^2 A} k'$ . The small correction to the equality  $\vec{p} = m d\vec{r}/dt$  caused by the spin-orbit part of the nuclear mean field can be neglected [7]. All  $\omega$ -dependent single -particle quantities in Eqs. (6)-(9) are determined by the potential, including the above mentioned additional terms.

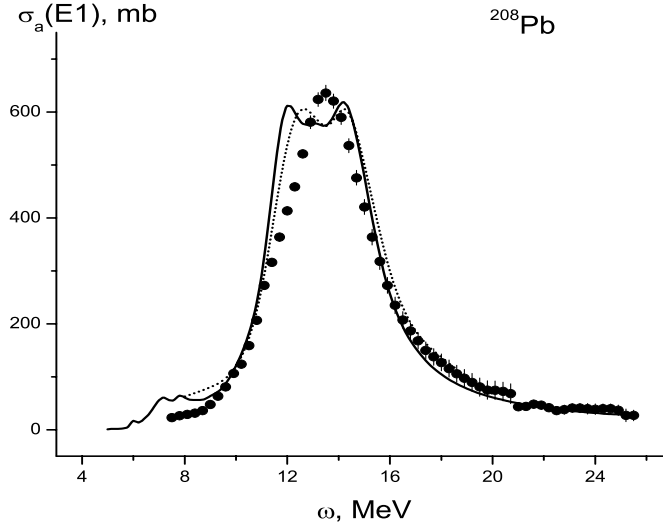
Using Eqs. (6)-(9) it is possible to find the dipole strength function

$$S(\omega) = -\frac{1}{\pi} \text{Im} \int r \mathcal{A}(r, r', \omega) \tilde{V}(r', \omega) dr dr' \quad , \quad (10)$$

where  $\omega$  is a photon energy and, hence, the total photoabsorption cross section in the energy region of the GDR:

$$\sigma(\omega) = \frac{8\pi^3}{3} \frac{e^2}{\hbar c} \omega S(\omega) \quad (11)$$

Unfortunately the results for photoabsorption cross section in  $^{208}\text{Pb}$  isotope obtained in this version of the presented approach happened to be practically similar to those obtained in Ref. [4] (Fig. 1).



**Fig 1.** The total photoabsorption cross section in  $^{208}\text{Pb}$  calculated in the first version of the presented approach (solid line) in comparison with corresponding results of Ref.[4] (dotted line) and the available experimental data [8].

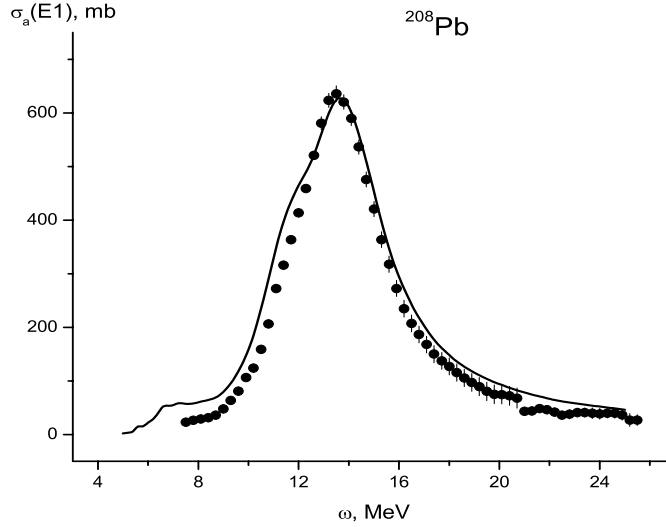
Due to that an attempt has been made to reconsider one of the principal foundation of the approach: to change the potential used for the calculations of all cRPA ingredients. This procedure has been carried out in Ref. [9] on the base of new phenomenological potential proposed in Ref. [10]. In this potential the mean field  $U(x)$  consists of pure nuclear parts, containing the isoscalar and isovector spin-orbit interaction, Coulomb field, the symmetry energy and is written in the following form:

$$U(x) = U_0(r) + (U_0^{SO}(r) + \frac{1}{2}U_1^{SO}(r)\tau^{(3)})\vec{l}\vec{s} + \frac{1}{2}v(r)\tau^{(3)} + \frac{1}{2}(1 - \tau^{(3)})U_C(r) \quad (12)$$

The space dependences of central field  $U_0(r)$  and spin-orbit interactions  $U_0^{SO}$  and  $U_1^{SO}$  are determined by the Woods-Saxon function  $f_{WS}(r, R^*, a)$  and its derivative  $df_{WS}(r)/dr$ , respectively. As to the symmetry potential  $v(r)$  and Coulomb field  $U_C(r)$ , they are calculated in a self-consistent way. The choice of all parameters determining every part of the potential  $U(x)$  is made by means of the minimization of the differences between calculated position of the energy levels and the experimental ones (for the details of the total procedure and obtained

values of potential parameters see Ref. [9]). It is worthwhile to notice that obtained value of the isovector constant  $f'$  is equal to 1.09 in contrast to the previously used value 1.0 (see, e.g., Ref. [4]).

The approach developed on the base of this new potential includes only two parameters which should be treated as the adjustable ones: the intensity of the momentum-dependent forces  $k'$  and the intensity of the smearing parameter  $\alpha$  (the values of  $\Delta$  and  $B$  are taken to be the same as in Ref. [4]). To determine them the experimental data on the photoabsorption cross section for  $^{208}\text{Pb}$  isotope [8] has been chosen as the most authentic ones. The corresponding calculations of the  $\sigma(\omega)$  has been carried out on the base of Eqs. (6)-(11) using the newly defined potential. The comparison of the obtained results with the experimental data in Fig. 2 allows to define the values of  $\alpha$  and  $k'$ :  $\alpha = 0.1 \text{ MeV}^{-1}$ ,  $k' = 0.4$ .



**Fig 2.** The calculated total photoabsorption cross section for  $^{208}\text{Pb}$  (solid line) in comparison with available experimental data [8].

To check the possibilities of the presented approach the available partial  $(n, \gamma)$  reaction cross sections for  $^{208}\text{Pb}$  isotope have been also investigated in the energy region of the GDR. The reaction amplitude which has been calculated on the base of Eqs. (6)-(9) has the following form:

$$M_c(\omega) = \langle (\lambda) || Y_1 || (\mu) \rangle \int \chi_{\varepsilon(\lambda)}^{(+)}(r) \tilde{V}(r, \omega) \chi_{\mu}(r) dr \quad , \quad (13)$$

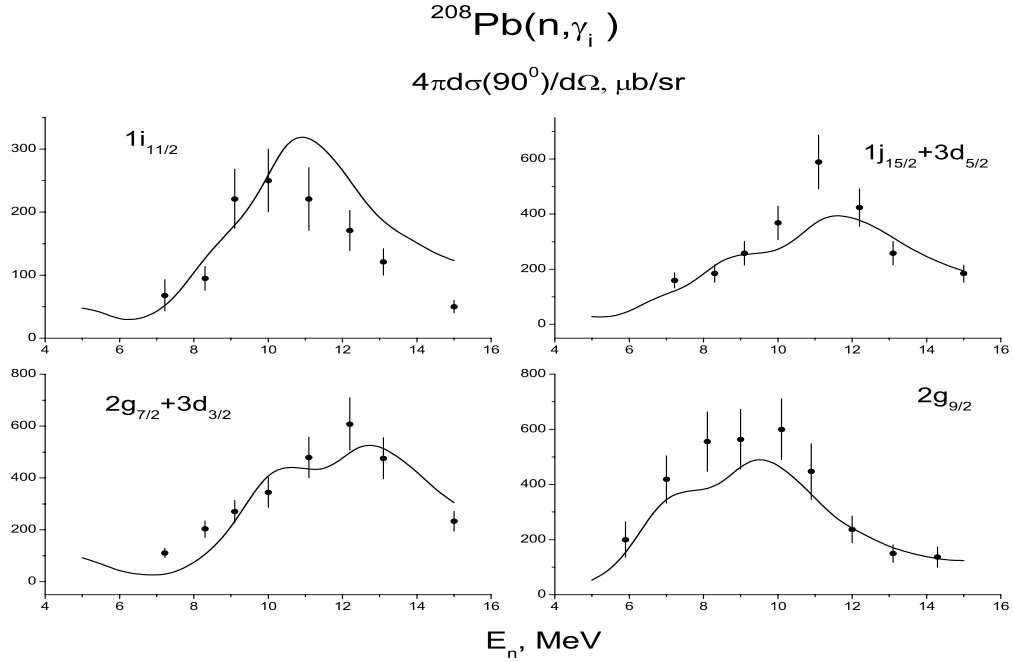
where  $c = \{\varepsilon(\lambda), \mu\}$  is the set of the reaction-channel quantum numbers,  $\mu = \{\varepsilon_{\mu}, (\mu)\}$  is the set of quantum numbers for single-particle state,  $\chi_{\mu}(r)$  is the radial wave function of the populated bound state with energy  $\varepsilon_{\mu}$  and  $\chi_{\varepsilon(\lambda)}^{(+)}(r)$  is the corresponding to this channel and normalized to the  $\delta$ -function of energy the scattering radial wave function of incident nucleon with energy  $\varepsilon = \omega + \varepsilon_{\mu}$  (for other notations see, e.g. [4]). The bound state wave function is calculated with new potential (12) while the scattering wave function - with potential (12) and additional terms connected with  $I(\omega)I(r)$  and  $\Pi(\omega)I(r)$  (see above). The cross sections of these reactions can be then presented in the following form:

$$\frac{d\sigma_c(\varepsilon, \theta)}{d\Omega} = \frac{8}{3} \pi^{5/2} \frac{e^2}{\hbar c} \omega \cdot \frac{\omega^2}{2mc^2\varepsilon} \sum_L A_L(\varepsilon, \mu) P_L(\cos\theta) \quad , \quad (14)$$

$$A_L(\varepsilon, \mu) = \sum_{(\lambda)(\lambda')} i^{l-l'} (-1)^{j_\mu+j} (1 - 111|L0) W(jj'11; Lj_\mu) <(\lambda)||Y_L||(\lambda') > M_c(\omega) M_c^*(\omega) \quad .$$

Here,  $P_L(\cos\theta)$  are the Legendre polynomials,  $(1-111|L0)$  and  $W(jj'11; Lj_\mu)$  are the Clebsch-Gordan coefficient and Racah coefficient, respectively.

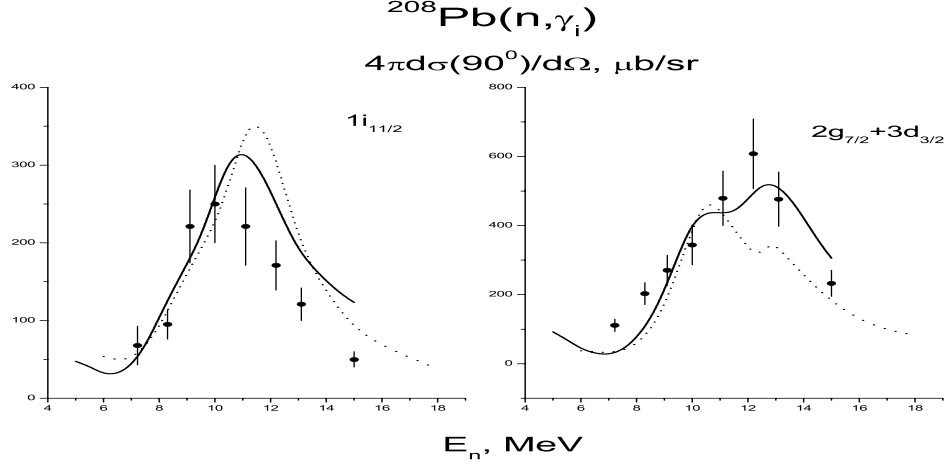
The results of these cross section calculations are shown in Figs. 3,4 in comparison with the available experimental data [11] and some results from Ref. [4]. As it is seen from Figs. 3,4 the presented approach based on the use of new potential [9] allows us to describe



**Fig 3.** The calculated on the base of the presented approach the partial  $(n,\gamma)$  reaction cross sections at  $90^\circ$  to some single-particle states in  $^{209}\text{Pb}$  in comparison with the available experimental data [11].

satisfactorily the studied photonuclear reactions: (i) the splitting of the GDR peak in  $^{208}\text{Pb}$  is absent; (ii) the obtained cross sections of partial  $(n,\gamma)$  reactions appreciably better agree with the available experimental data than those from Ref. [4]. The energy-averaged (due to the use of the smearing parameter  $I(\omega)$ ) differential cross sections for neutron radiative capture,  $d\sigma_\mu/d\Omega$ , are calculated without use of any adjustable parameters. Each calculation of the cross section is carried out with account of the experimental spectroscopic factor value,  $S_\mu$ , of the final product-nucleus single-particle state populated after the capture. The corresponding experimental data are taken from Ref. [12].

As a conclusion one could say the following. In the presented approach a reasonable description of the experimental data on the total photoabsorption cross section and the partial  $(n,\gamma)$  reactions for  $^{208}\text{Pb}$  isotope is obtained. In the future it is planned to expand the such investigations on the other medium-heavy mass nuclei having the similar structure. Besides, it is supposed to carry out these studies on the base of the potential [9] using the new set of parameters. Finally, it is planned to perform in a framework of the presented approach some predictive investigations, for instance, of  $(\gamma,n)$  reactions where there are some appropriate experimental data.



**Fig 4.** The calculated partial cross sections at  $90^\circ$  for neutron radiative capture (solid lines) to some single-particle states in  $^{209}\text{Pb}$  in comparison with the corresponding results of Ref. [4] (dotted lines) and the available experimental data [11].

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