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Primordial $\alpha + d \rightarrow {}^{6}\text{Li} + \gamma$ reaction and second Lithium puzzle

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During the Big Bang, ⁶Li was synthesized via the ²H(α, γ)⁶Li reaction. After almost 25 years of the failed attempts to measure the ${}^{2}H(\alpha,\gamma)^{6}Li$ reaction in the lab at the Big Bang energies, just recently the LUNA collaboration presented the first successful measurements at two different Big Bang energies [M. Anders et al., Phys. Rev. Lett. 113, 042501 (2014)]. In this paper we will discuss how to improve the accuracy of the direct experiment. To this end the photon's angular distribution is calculated in the potential model. It contains contributions from electric dipole and quadrupole transitions and their interference, which dramatically changes the photon's angular distribution. The calculated distributions at different Big Bang energies have a single peak at $\sim 50^{\circ}$. These calculations provide the best kinematic conditions to measure the ${}^{2}H(\alpha,\gamma)^{6}Li$ reaction. The expressions for the total cross section and astrophysical factor are also derived by integrating the differential cross section over the photon's solid angle. The LUNA data are in excellent agreement with our calculations using a potential approach combined with a well established asymptotic normalization coefficient for ${}^{6}\text{Li} \rightarrow \alpha + d$. Comparisons of the available experimental data for the S_{24} astrophysical factor and different calculations are presented. The Big Bang lithium isotopic ratio ${}^{6}\text{Li}/{}^{7}\text{Li} = (1.5 \pm 0.3) \times 10^{-5}$ following from the LUNA data and the present analysis are discussed in the context of the disagreement between the observational data and the standard Big Bang model, which constitutes the second Lithium problem.

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I. INTRODUCTION

The primordial nuclei were formed during the first 20 minutes after the Big Bang. Among them the lithium isotopes, ⁷Li and a much smaller amount of ⁶Li, were synthesized. Later on, cosmic rays, novae and pulsations of AGB stars were the main generators of the ⁷Li isotope, and ⁶Li was formed mainly by cosmic rays. In 1982, two important papers [1, 2] for the first time noted that metal-poor $(-2.4 \leq [Fe/H] \leq -1.4)$, warm $(5700 \leq T_{eff} \leq 6250 \text{ K})$ dwarf stars demonstrated remarkably constant ⁷Li abundance (Spite plateau), which does not depend on metallicity and effective temperature. It was quite a surprising observation because depletion of lithium over such a broad temperature range should be significant. Because it was impossible to explain the existence of the Spite plateau over a wide range of temperatures, it was suggested that no depletion of ⁷Li took place in the observed dwarf warm stars and that the constant abundance of ⁷Li is the primordial one. However, this interpretation of the Spite plateau was periodically challenged. For example, in [3, 4] the meltdown of the Spite plateau was discovered in some low metallicity stars.

Explanation of the lithium abundance in low metallicity stars in the halo of our Galaxy where the lithium abundance demonstrates independence on metallicity, forming the Spite plateau, became one of the hot topics in modern cosmology/nuclear astrophysics. Note that the observations of the primordial lithium are restricted to white dwarfs because the loosely bound ⁷Li nuclei are easily destroyed by the ⁷Li(p, α)⁴He reaction when the temperature exceeds 2.6×10^6 K. For this reason, red giants cannot be used to determine the lithium primordial abundance.

In the standard Big Bang nucleosynthesis model, ⁷Li is formed right after the Big Bang, together with ¹H, ²H, ³He and ⁴He. The primordial reactions start from the deuteron formation $p + n \rightarrow d + \gamma$. The deuteron's yield depends on the primordial baryon/photon ratio η_B . Because the deuterons are seeds, which are necessary to synthesize heavier elements, the abundance of heavier elements, and lithium in particular, also depends on η_B . ²H and ⁷Li are two primordial nuclei which is most sensitive to η_B .

The abundance of the primordial ⁷Li within the framework of the standard Big Bang scenario, calculated using the extended reaction network and nine years of WMAP results [5], is ⁷Li/H = 5.13×10^{-10} . It is ⁷Li/H = $(4.56 - 5.34) \times 10^{-10}$ [6] based on the Planck results [7] plus the information about the lensing potential and ground-based high resolution experiments. The latter is considered to be the most up-to-date estimation of the ⁷Li isotope abundance within the standard Bing Bang scenario. This abundance remains significantly higher than more recent observations in metal poor halo stars [4]: ⁷Li/H = $1.58^{+0.35}_{-0.28} \times 10^{-10}$. The shortage of the observed ⁷Li compared to the standard Big Bang predictions represents the so-called first Lithium puzzle.

During the Big Bang a very small amount of ⁶Li was synthesized via the ²H(α, γ)⁶Li reaction. Later, ⁶Li was mostly formed by cosmic rays. The primordial ⁶Li is assumed to be present in the gas from which the stars were formed. Unlike most of the other elements, when ⁶Li is synthesized inside the stars by hydrostatic nucleosynthesis, it is quickly destroyed. But in the atmosphere of the halo metal-poor warm dwarfs, the primordial ⁶Li can survive for 13 billion years not being affected by cosmic rays, although its survival can be questioned. ⁷Li is used to help determine the primordial Big Bang ⁶Li abundance. First, the presence of ⁶Li constrains the destruction of ⁷Li, because ⁶Li is more easily destroyed than ⁷Li. Besides, if ⁶Li was formed before the formation of the stars, then the same is true for ⁷Li.

Stellar ⁷Li abundance is usually determined from the resonance line at 670.8 nm but in exceptional cases also from the weaker line at 610.4 nm. The isotope ⁶Li can be detected through the isotopic shift in the Li I 670.8 nm line. The distortion of the line profile is very small and therefore requires very high quality spectra. Ref. [8] reported for the first time the detection of a high abundance of ⁶Li in very metal-poor stars. The authors concluded that the observed ⁶Li was formed during Big-Bang nucleosynthesis. The detection of ⁶Li was based on the fact noted above, that the presence of ⁶Li in the stellar atmosphere causes an asymmetry in the Li 670.8 nm line. The average ⁶Li/⁷Li isotopic ratio in the nine stars, in which ⁶Li was detected, was ⁶Li/⁷Li ~ 5×10^{-2} [8]. Such a high isotopic ratio of the primordial lithium isotopes in the metal-poor stars contradicts the Big-Bang based model predictions ⁶Li/⁷Li ~ 10^{-5} [6] and cannot be explained by the galactic cosmic rays. This disagreement between the observations and Big Bang predictions of the lithium isotopic ratio constitutes the second Lithium problem.

Later it was pointed out in [9] that the line asymmetry caused by convection in the photospheres of metal-poor stars is practically indistinguishable from the asymmetry produced by a weak ⁶Li distortion of a symmetric ⁷Li line. Hence, the ⁶Li abundance obtained in [8] could be significantly overestimated, and the result obtained in [8] can be considered only as an upper limit of the lithium isotopic ratio. In Ref. [10] the lithium isotopic ratio was reanalyzed within the framework of the 3D, non-local thermodynamic equilibrium (NLTE) model. The authors came to the conclusion that "the observational support for a significant and non-standard ⁶Li production source in the early universe is substantially weakened by our findings" [10], which opens a way to a hope that the primordial abundance of ⁶Li calculated in the standard Big Bang nucleosynthesis can be eventually reconciled with observational data.

The yields of the observed and predicted primordial ⁷Li are established quite well [6]. If the observed ⁶Li is primordial (a Big Bang product) then its abundance is determined by the ²H(α, γ)⁶Li reaction. The first successful attempt to measure the ²H(α, γ)⁶Li reaction was reported in [11] where residual ⁶Li nuclei were detected. The astrophysical factor was measured in the vicinity of the first resonance ⁶Li(3⁺) at the relative α -d energy E = 0.712 MeV and at higher energies. But no data were obtained at Big Bang energies, $30 \leq E \leq 400$ keV. In [12] the astrophysical $S_{24}(E)$ factor was also measured only at the resonance energy, using in-beam spectroscopy.

In Ref. [13] for the first time, an attempt was made to measure the astrophysical factor at the Big Bang energies, using the Coulomb breakup of ⁶Li at 26 MeV/A energy on a ²⁰⁸Pb target. However, only an upper limit was established. The failure of this indirect attempt to measure the $S_{24}(E)$ astrophysical factor could be anticipated because the *E*1 transition, which usually dominates, is suppressed in the case under consideration: the effective charge for the dipole transition is very small owing to practically the same charge/mass ratio for α -particle and deuteron. Because the Coulomb dissociation cross section is dominated by the *E*2 transition, the obtained data may be considered only as an upper limit. After that, another unsuccessful attempt to measure the $S_{24}(E)$ factor ended with an upper limit $S_{24}(53 \text{ keV}) < 2.0 \times 10^{-7}$ MeVb and a pessimistic conclusion that it would be impossible to measure directly $S_{24}(E)$ at Big Bang energies [14].

The second attempt to use the indirect Coulomb dissociation technique was made in [15], where the breakup of ⁶Li ions at 150 MeV/A on a ²⁰⁸Pb target was measured. However, in this case, the breakup was dominated by nuclear breakup, which overwhelmed the Coulomb breakup. Hence, no information about $S_{24}(E)$ was extracted from the analysis of the breakup data. Further, in Ref. [15] the astrophysical factor was calculated using a two-body potential model (see below). Finally, after almost 25 years of failed attempts to measure the ²H(α, γ)⁶Li reaction at the Big Bang energies, just recently the LUNA collaboration presented the first successful measurements at two different Big Bang energies [16]. Definitely it is a remarkable achievement in the studies of Big Bang nucleosynthesis.

In this work we discuss the astrophysical ${}^{2}H(\alpha, \gamma){}^{6}Li$ reaction within the framework of the potential approach and impact on experimental measurements. For the first time, we present the angular distribution of the photons emitted in this direct radiative capture. Although the photon differential cross section is being derived for the ${}^{2}H(\alpha, \gamma){}^{6}Li$ process, it can be applied for any direct electric radiative capture reaction. The calculated angular distributions provide the best kinematics to be used in the measurement of the emitted photons, which differ from the one used in the LUNA experiment. Optimal kinematics will allow one to decrease significantly the uncertainty of direct measurements of the ${}^{2}H(\alpha, \gamma){}^{6}Li$ process compared to the uncertainties in the LUNA experiment. By integrating the differential cross section over the photon solid angle, the total cross section and astrophysical factor of the direct radiative capture are derived. The calculations of the photon's angular distribution and astrophysical $S_{24}(E)$ factor are done in the potential model using the well determined asymptotic normalization coefficient for the virtual decay ${}^{6}Li \rightarrow \alpha + d$. The primordial ${}^{6}Li$ abundance is presented.

II. PHOTON DIFFERENTIAL CROSS SECTIONS, TOTAL CROSS SECTIONS AND ASTROPHYSICAL S-FACTORS

A. Photon angular distribution in direct radiative capture

In this section the expression for the angular distribution of the photons emitted in the $\alpha(d, \gamma)^6$ Li direct radiative capture is derived and further simplified in the subsequent section. This result can help to improve future experiments on this reaction by decreasing their uncertainties. Often photon angular distributions are not discussed in the papers dealing with measurements of the astrophysical factors. That is why we believe it is timely to do it. Besides by integrating the photon differential cross section over the photon's solid angle the total cross section and the astrophysical factor can be derived.

We consider the photon angular distribution taking into account the spin-orbit interaction in the initial state. Hence, the initial scattering wave function depends on the initial α – d relative orbital angular momentum l_i , the channel spin s and the total angular momentum in the initial channel J_i . In the case under consideration $s = J_d$, where $J_d = 1$ is the spin of the deuteron. The differential cross section of the emitted photons with momentum \mathbf{k}_{γ} and helicity $\lambda = \pm 1$ in the electromagnetic transition from the initial continuum state l_i , s, J_i to the final state l_f , s, J_f in the center-of-mass of ⁶Li is given by

$$\frac{\mathrm{d}\sigma_{\lambda}}{\mathrm{d}\Omega} \sim k_{\gamma}^{2} \left| -\frac{1}{c} \int \mathrm{d}\mathbf{r} \left\langle \varphi_{^{6}\mathrm{Li}}(\zeta_{\alpha}, \zeta_{d}; \mathbf{r}_{\alpha \, d}) \left| \mathbf{\hat{J}}(\mathbf{r}) \right| \Psi_{\mathbf{k}}^{(+)}(\zeta_{\alpha}, \zeta_{d}; \mathbf{r}_{\alpha \, d}) \right\rangle \cdot \mathbf{A}_{\lambda \, \mathbf{k}_{\gamma}}^{*}(\mathbf{r}) \right|^{2}.$$
(1)

Here, $\mathbf{A}_{\lambda \mathbf{k}_{\gamma}}(\mathbf{r})$ is the vector-potential of the photon with helicity λ and momentum \mathbf{k}_{γ} at coordinate \mathbf{r} . The initial wave function is

$$\Psi_{\mathbf{k}}^{(+)}(\zeta_{\alpha},\,\zeta_{d};\,\mathbf{r}_{\alpha\,d}) = \varphi_{\alpha}\left(\zeta_{\alpha}\right)\varphi_{d}(\zeta_{d})\,\psi^{(+)}(\mathbf{k},\,\mathbf{r}_{\alpha\,d}),\tag{2}$$

 $\varphi(\zeta_i)$ is the bound-state wave function of nucleus *i* with the set of the internal coordinates ζ_i , which includes spinisospin variables. $\psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d})$ is the α – d scattering wave function in the initial state, $\mathbf{r}_{\alpha d}$ is the radius-vector We use the long wavelength approximation, which is valid for $k_{\gamma} R_{\alpha d} \ll 1$. Here $R_{\alpha d}$ is the effective $\alpha - d$ distance determined so that distances $r \sim R_{\alpha d}$ give the dominant contribution to the amplitude of the direct radiative capture. The long electromagnetic wavelength of the emitted radiation allows us to approximate the charge current density by the current density of the point-like α -particle and deuteron neglecting their internal structure:

$$\hat{\mathbf{J}}(\mathbf{r}) = \frac{Z_d e}{2 m_d} \left[\delta(\mathbf{r} - \mathbf{r}_d) \, \hat{\mathbf{p}}_d + \hat{\mathbf{p}}_d \, \delta(\mathbf{r} - \mathbf{r}_d) \right] + \frac{Z_\alpha e}{2 m_\alpha} \left[\delta(\mathbf{r} - \mathbf{r}_\alpha) \, \hat{\mathbf{p}}_\alpha + \hat{\mathbf{p}}_\alpha \, \delta(\mathbf{r} - \mathbf{r}_\alpha) \right], \tag{3}$$

where $\hat{\mathbf{p}}_i = -i\hbar\partial/\partial\mathbf{r}_i$ is the momentum operator, $\mathbf{r}_d = -(m_\alpha/m_{\alpha d})\mathbf{r}_{\alpha d}$ and $\mathbf{r}_\alpha = (m_d/m_{\alpha d})\mathbf{r}_{\alpha d}$ are the coordinates of the centers of mass of the deuteron and alpha-particle, respectively, m_i and Z_i are the mass and atomic number of nucleus *i* and $m_{ij} = m_i + m_j$. We neglect here the spin contribution to the current density because below we consider only the electric transitions which are largely due to the charge current.

Now the overlap function of the bound-state wave functions of ⁶Li, α -particle and deuteron can be introduced:

$$I_{l_f s J_f}(\mathbf{r}_{\alpha d}) = \left\langle \varphi_{\alpha}(\zeta_{\alpha}) \varphi_d(\zeta_d) \middle| \varphi_{^6\mathrm{Li}}(\zeta_{\alpha}, \zeta_d; \mathbf{r}_{\alpha d}) \right\rangle$$

= $\sum_{m_{l_f} m''_s} \left\langle l_f m_{l_f} s m''_s \middle| J_f M_f \right\rangle Y_{l_f m_{l_f}}(\hat{\mathbf{r}}_{\alpha d}) \chi_{s m''_s} I_{l_f s J_f}(r_{\alpha d}),$ (4)

where $I_{l_f s J_f}(r_{\alpha d})$ is the radial overlap function, $\langle l_f m_{l_f} s m_s'' | J_f M_f \rangle$ is the Clebsch-Gordan coefficient, l_f is the α - d relative orbital angular momentum in the bound state, $J_f = 1$ is the spin of the ground state of ⁶Li; $\chi_{s m_s''}$ is the spin wave function describing the state with the channel spin s and its projection m_s'' , $\hat{\mathbf{r}} = \mathbf{r}/\mathbf{r}$ is the unit vector. The integration in the matrix element $\langle \varphi_{\alpha}(\zeta \alpha) \varphi_d(\zeta_d) | \varphi_{^6\text{Li}}(\zeta_{\alpha}, \zeta_d; \mathbf{r}_{\alpha d}) \rangle$ is taken over all the internal coordinates ζ_{α} and ζ_d making the overlap function depending only on the radius-vector $\mathbf{r}_{\alpha d}$.

In the peripheral region the radial overlap function is given by

$$I_{l_f s J_f}(r_{\alpha d}) \stackrel{r_{\alpha d} > r_0}{\approx} C_{l_f s J_f} W_{-\eta, l_f + 1/2}(2 \kappa r_{\alpha d}) / r_{\alpha d}, \tag{5}$$

where $C_{l_f s J_f}$ is the asymptotic normalization coefficient (ANC) for the virtual decay ⁶Li $\rightarrow \alpha + d$ expressed in fm^{-1/2}, $W_{-\eta_f, l_f+1/2}(2 \kappa r_{\alpha d})$ is the Whittaker function determining the radial shape of the overlap function beyond of the $\alpha - d$ nuclear interaction region, $\eta_f = (Z_{\alpha} Z_d e^2/\hbar c)(\mu_{\alpha d} c/\hbar)(1/\kappa)$ is the Coulomb $\alpha - d$ bound-state parameter, $\kappa = \sqrt{2 \mu_{\alpha d} c^2 \varepsilon}/\hbar c$ is the $\alpha - d$ bound-state wave number expressed in fm⁻¹. The radial overlap function is expressed in fm^{-3/2}. r_0 is the channel radius, which is selected so that at $r_{\alpha d} > r_0$ the nuclear interaction between the deuteron and α -particle is negligible.

The matrix element in (1) now can be rewritten as

$$\frac{1}{c} \int d\mathbf{r} \left\langle \varphi_{^{6}\mathrm{Li}}(\zeta_{\alpha}, \zeta_{d}; \mathbf{r}_{\alpha d}) \left| \hat{\mathbf{J}}(\mathbf{r}) \right| \Psi^{(+)}(\zeta_{\alpha}, \zeta_{d}; \mathbf{r}_{\alpha d}) \right\rangle \cdot \mathbf{A}_{\lambda \mathbf{k}_{\gamma}}^{*}(\mathbf{r}) \\
= \frac{1}{c} \int d\mathbf{r} \left\langle I_{l_{f} s J_{f}}(\mathbf{r}_{\alpha d}) \left| \hat{\mathbf{J}}(\mathbf{r}) \right| \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle \cdot \mathbf{A}_{\lambda \mathbf{k}_{\gamma}}^{*}(\mathbf{r}).$$
(6)

To simplify further this matrix element we need to use the multipole expansion of the vector potential [17, 18]:

$$\mathbf{A}_{\lambda \mathbf{k}_{\gamma}}(\mathbf{r}) = \frac{1}{2\pi} \sqrt{\frac{\hbar c}{k_{\gamma}}} \mathbf{e}_{\lambda \mathbf{k}_{\gamma}} e^{i \mathbf{k}_{\gamma} \cdot \mathbf{r}} = \frac{1}{2\sqrt{2\pi} k_{\gamma}} \sum_{LM} \sqrt{2L+1} D_{M\lambda}^{L}(\varphi,\theta,0) \times \left(\mathbf{A}_{ek_{\gamma}LM}(\mathbf{r}) + \lambda \mathbf{A}_{mk_{\gamma}LM}(\mathbf{r})\right).$$
(7)

Here, $\mathbf{e}_{\lambda \mathbf{k}_{\gamma}}$ is the unit polarization vector of the plane wave, which is orthogonal to the photon momentum \mathbf{k}_{γ} , $\mathbf{A}_{ek_{\gamma}LM}(\mathbf{r})$ and $\mathbf{A}_{mk_{\gamma}LM}(\mathbf{r})$ are the electric and magnetic multipoles, correspondingly. In the system $z \| \mathbf{k}_{\gamma}$ the helicity of the circularly polarized photon $\lambda = \pm 1$. $D_{M\lambda}^{L}(\varphi, \theta, 0)$ is the Wigner *D*-function, *L* is the multipolarity of the transition. In Eq. (7) only the electric multipoles $\mathbf{A}_{ek_{\gamma}LM}(\mathbf{r})$ will be taken into account because for the reaction under consideration the contribution of the magnetic multipoles $\mathbf{A}_{m k_{\gamma} L M}(\mathbf{r})$ is negligible [19]. Following Ref. [17], $\mathbf{A}_{e\,k_{\gamma}\,L\,M}(\mathbf{r})$ can be rewritten as

$$\mathbf{A}_{e\,k_{\gamma}\,L\,M}(\mathbf{r}) = -2\,i^{L}\,\sqrt{\frac{\hbar\,c}{k_{\gamma}}} \left[\boldsymbol{\nabla}_{\mathbf{r}} \times \left(j_{L}(k_{\gamma}\,r)\,\mathbf{Y}_{LM}^{L}(\hat{\mathbf{r}}) \right) \right] \\ = 2\,i^{L-1}\,\sqrt{\hbar\,c\,k_{\gamma}} \left[\sqrt{\frac{L+1}{2\,L+1}}\,j_{L-1}(k_{\gamma}\,r)\,\mathbf{Y}_{LM}^{L-1}(\hat{\mathbf{r}}) - \sqrt{\frac{L}{2\,L+1}}\,j_{L+1}(k_{\gamma}\,r)\,\mathbf{Y}_{LM}^{L+1}(\hat{\mathbf{r}}) \right], \tag{8}$$

where $\mathbf{Y}_{LM}^{\tilde{L}}(\hat{\mathbf{r}})$ is the vector spherical harmonics [17, 18] and $j_L(k_{\gamma} r)$ is the spherical Bessel function. Now the matrix element (6) can be reduced to

$$\frac{1}{c} \int d\mathbf{r} \left\langle I_{l_{f}sJ_{f}}(\mathbf{r}_{\alpha d}) \left| \hat{\mathbf{J}}(\mathbf{r}) \right| \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle \cdot \mathbf{A}_{\lambda \mathbf{k}_{\gamma}}^{*}(\mathbf{r}) \\
= \sqrt{\frac{\hbar}{2\pi c k_{\gamma}}} \sum_{LM} i^{-L+1} \sqrt{2L+1} \left(D_{M\lambda}^{L}(\varphi, \theta, 0) \right)^{*} \int d\mathbf{r} \left\langle I_{l_{f}sJ_{f}}(\mathbf{r}_{\alpha d}) \left| \hat{\mathbf{J}}(\mathbf{r}) \right| \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle \\
\times \left[\sqrt{\frac{L+1}{2L+1}} j_{L-1}(k_{\gamma} r) \left(\mathbf{Y}_{LM}^{L-1}(\hat{\mathbf{r}}) \right)^{*} - \sqrt{\frac{L}{2L+1}} j_{L+1}(k_{\gamma} r) \left(\mathbf{Y}_{LM}^{L+1}(\hat{\mathbf{r}}) \right)^{*} \right] \\
\approx \sqrt{\frac{\hbar}{2\pi c k_{\gamma}}} \sum_{LM} \frac{i^{-L+1} k_{\gamma}^{L-1}}{(2L-1)!!} \sqrt{L+1} \left(D_{M\lambda}^{L}(\varphi, \theta, 0) \right)^{*} \int d\mathbf{r} \left\langle I_{l_{f}sJ_{f}}(\mathbf{r}_{\alpha d}) \left| \hat{\mathbf{J}}(\mathbf{r}) \right| \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle \\
\times r^{L-1} \left(\mathbf{Y}_{LM}^{L-1}(\hat{\mathbf{r}}) \right)^{*}.$$
(9)

In the long wavelength approximation $k_{\gamma} r \ll 1$, $j_L(k_{\gamma} r) \approx (k_{\gamma} r)^L/(2L+1)!!$. Hence, the lowest partial waves dominate and the term containing $j_{L+1}(k_{\gamma} r) \approx (k_{\gamma} r)^{L+1}/(2L+3)!!$ is small compared to the term containing $j_{L-1}(k_{\gamma} r) \approx (k_{\gamma} r)^{L-1}/(2L-1)!!$ and can be neglected. Taking into account that [20]

$$\boldsymbol{\nabla}_{\mathbf{r}}[r^{L} Y_{LM}(\hat{\mathbf{r}})] = \sqrt{L(2L+1)} r^{L-1} \mathbf{Y}_{LM}^{L-1}(\hat{\mathbf{r}}), \qquad (10)$$

Equation (9) can be reduced to

$$\frac{1}{c} \int d\mathbf{r} \left\langle I_{l_{f}s J_{f}}(\mathbf{r}_{\alpha d}) | \hat{\mathbf{J}}(\mathbf{r}) | \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle \mathbf{A}_{\lambda \mathbf{k}_{\gamma}}^{*}(\mathbf{r}) \\
\approx \sqrt{\frac{\hbar}{2 \pi c k_{\gamma}}} \sum_{LM} \frac{i^{-L+1} k_{\gamma}^{L-1}}{(2L+1)!!} \sqrt{\frac{(L+1)(2L+1)}{L}} \left(D_{M \lambda}^{L}(\varphi, \theta, 0) \right)^{*} \\
\int d\mathbf{r} \left\langle I_{l_{f}s J_{f}}(\mathbf{r}_{\alpha d}) | \hat{\mathbf{J}}(\mathbf{r}) | \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle \nabla_{\mathbf{r}} [r^{L} \left(Y_{LM}(\hat{\mathbf{r}}) \right)^{*}].$$
(11)

Integrating by parts and using the static current conservation

$$\boldsymbol{\nabla}_{\mathbf{r}} \hat{\mathbf{J}}(\mathbf{r}) = i \, k_{\gamma} \, c \, \hat{\rho}(\mathbf{r}), \tag{12}$$

where

$$\hat{\rho}(\mathbf{r}) = Z_d \, e \, \delta(\mathbf{r} - \mathbf{r}_d) + Z_\alpha \, e \, \delta(\mathbf{r} - \mathbf{r}_\alpha) \tag{13}$$

is the charge density operator, one gets

$$\frac{1}{c} \int d\mathbf{r} \left\langle I_{l_{f}sJ_{f}}(\mathbf{r}_{\alpha d}) | \hat{\mathbf{J}}(\mathbf{r}) | \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle \mathbf{A}_{\lambda \mathbf{k} \gamma}^{*}(\mathbf{r}) \\
\approx \frac{1}{2\pi} \sqrt{\frac{\hbar c}{2 k_{\gamma}}} \sum_{LM} \frac{i^{-L} k_{\gamma}^{L}}{(2L-1)!!} \sqrt{\frac{L+1}{L}} \left(D_{M \lambda}^{L}(\varphi, \theta, 0) \right)^{*} \\
\times \left\langle I_{l_{f}sJ_{f}}(\mathbf{r}_{\alpha d}) | \left(\hat{Q}_{LM}^{(e)}(\mathbf{r}_{\alpha d}) \right)^{*} | \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle.$$
(14)

Here,

$$\hat{Q}_{LM}^{(e)}(\mathbf{r}_{\alpha d}) = \sqrt{\frac{4\pi}{2L+1}} \int d\mathbf{r} \,\hat{\rho}(\mathbf{r}) \, r^L \, Y_{LM}(\hat{\mathbf{r}}), \tag{15}$$

is the electric static 2^L moment operator.

Thus the initial matrix element (6) containing $\mathbf{A}_{\lambda \mathbf{k}_{\gamma}}^{*}(\mathbf{r})$ after the multipole expansion and series of transformations is reduced to the matrix element, which is expressed in terms of the electric charge density operator. This is possible due to Siegert's theorem [21].

Equation (1) for the differential cross section of the electric transition takes the form

$$\frac{\mathrm{d}\sigma_{\lambda}}{\mathrm{d}\Omega} \sim \left| -\frac{1}{2\pi} \sqrt{\frac{\hbar c}{2 k_{\gamma}}} \sum_{LM} \frac{i^{-L} k_{\gamma}^{L+1}}{(2L-1)!!} \sqrt{\frac{L+1}{L}} \left(D_{M\lambda}^{L}(\varphi,\theta,0) \right)^{*} \times \left\langle I_{l_{f} s J_{f}}(\mathbf{r}_{\alpha d}) \left| \left(\hat{Q}_{LM}^{(e)}(\mathbf{r}_{\alpha d}) \right)^{*} \left| \psi^{(+)}(\mathbf{k},\mathbf{r}_{\alpha d}) \right\rangle \right|^{2} \right|^{2}$$
(16)

In the case under consideration the dominant contribution comes from the electric dipole (L = 1) and electric quadrupole (L = 2) transitions. Because the sum over multipoles L is incoherent the interference of the dipole and quadrupole amplitudes should be taken into account.

Integrating over \mathbf{r} in Eq. (15) one gets

$$\hat{Q}_{LM}^{(e)}(\mathbf{r}_{\alpha\,d}) = \sqrt{\frac{4\,\pi}{2\,L+1}} \, e \, Z_{eff(L)} \, r_{\alpha\,d}^L \, Y_{LM}(\hat{\mathbf{r}}_{\alpha\,d}). \tag{17}$$

 $e Z_{eff(L)}$ is the effective charge for the electric transition of the multipolarity L, where

$$Z_{eff(L)} = \mu_{\alpha d}^L \left(\frac{Z_\alpha}{m_\alpha^L} + (-1)^L \frac{Z_d}{m_d^L} \right).$$
⁽¹⁸⁾

To derive Eq. (17) we took into account that $Y_{LM}(-\hat{\mathbf{r}}_{\alpha d}) = (-1)^L Y_{LM}(\hat{\mathbf{r}}_{\alpha d})$. The improvement of the leading order of the long wavelength approximation leads to the replacement of $r_{\alpha d}^L$ in Eq. (17) by more refined expressions [22]. For the dipole transition $r_{\alpha d}$ in Eq. (17) should be replaced by

$$O_1(r_{\alpha d}) = \frac{3}{y^3} \left[(y^2 - 2) \sin y + 2y \cos y \right] r_{\alpha d}$$
(19)

and for the quadrupole transition $r_{\alpha d}^2$ should be replaced by

$$O_2(r_{\alpha d}) = \frac{15}{y^5} \left[(5y^2 - 12) \sin y + (12 - y^2) y \cos y \right] r_{\alpha d}^2, \tag{20}$$

where $y = k_{\gamma} r_{\alpha d}$.

The initial scattering wave function with spin-orbit interaction is given by

$$\psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) = 4 \pi \sum_{J_i} \sum_{l_i} i^{l_i} \psi^{(+)}_{l_i s J_i}(k, r_{\alpha d}) \sum_{m_{l_i} m_s} \langle l_i m_{l_i} s m_s | J_i M_i \rangle Y_{l_i m_{l_i}}(\hat{\mathbf{r}}_{\alpha d}) \\ \times \chi_{s m_s} \sum_{m'_{l_i} m'_s} \langle l_i m'_{l_i} s m'_s | J_i M_i \rangle Y^*_{l_i m'_{l_i}}(\hat{\mathbf{k}}).$$
(21)

It is assumed that the projection M_i of J_i is fixed. For $z \| \mathbf{k} \ Y_{l_i m'_{l_i}}(\hat{\mathbf{k}}) = \sqrt{(2l_i + 1)/4\pi} \, \delta_{m'_{l_i} 0}$ and, hence, $m'_s = M_i$. Then

$$\psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) = \sum_{J_i} \sum_{l_i} i^{l_i} \sqrt{4 \pi (2 l_i + 1)} \psi^{(+)}_{l_i s J_i}(k, r_{\alpha d}) \sum_{m_{l_i} m_s} \langle l_i m_{l_i} s m_s | J_i M_i \rangle Y_{l_i m_{l_i}}(\hat{\mathbf{r}}_{\alpha d}) \\ \times \chi_{s m_s} \langle l_i 0 s M_i | J_i M_i \rangle.$$
(22)

The asymptotic behavior of the radial scattering wave function is taken in the form

$$\psi_{l_i \, sJ_i}^{(+)}(k, \, r_{\alpha \, d}) \approx \frac{1}{2 \, i \, r_{\alpha \, d}} e^{-i \, \delta_{l_i \, sJ_i}} \left[I_{l_i}(k, r_{\alpha \, d}) - e^{2 \, i \, \delta_{l_i \, sJ_i}} \, O_{l_i}(k, \, r_{\alpha \, d}) \right]. \tag{23}$$

$$I_{l_i}(k, r_{\alpha d}) = G_{l_i}(k, r_{\alpha d}) - i F_{l_i}(k, r_{\alpha d})$$
(24)

7

and

$$O_{l_i}(k, r_{\alpha d}) = G_{l_i}(k, r_{\alpha d}) + i F_{l_i}(k, r_{\alpha d})$$
(25)

are the incoming and outgoing spherical waves expressed in terms of the regular, $F_{l_i}(k, r_{\alpha d})$, and singular, $G_{l_i}(k, r_{\alpha d})$, Coulomb solutions of the radial Schrödinger equation. $\delta_{l_i s J_i}$ is the scattering phase shift. Inserting Eqs (4) and (22) into the matrix element of Eq. (16) one finds that

$$\left\langle I_{l_{f}s J_{f}}(\mathbf{r}_{\alpha d}) \left| \hat{Q}_{LM}^{(e)}(\mathbf{r}) \right| \psi^{(+)}(\mathbf{k}, \mathbf{r}_{\alpha d}) \right\rangle = \sqrt{\frac{4\pi}{2L+1}} \sum_{J_{i}} \sum_{l_{i}} \sum_{l_{i}} i^{l_{i}} \sqrt{4\pi (2l_{i}+1)}$$

$$\times \sum_{m_{l_{i}}m_{s}m_{l_{f}}} \left\langle l_{f} m_{l_{f}} s m_{s} \left| J_{f} M_{f} \right\rangle \left\langle l_{i} m_{l_{i}} s m_{s} \right| J_{i} M_{i} \right\rangle \left\langle l_{i} 0 s M_{i} \right| J_{i} M_{i} \right\rangle$$

$$\times \left\langle I_{l_{f}s J_{f}}(r_{\alpha d}) Y_{l_{f}m_{l_{f}}}(\hat{\mathbf{r}}_{\alpha d}) \left| e Z_{eff(L)} r_{\alpha d}^{L} Y_{LM}^{*}(\hat{\mathbf{r}}_{\alpha d}) \right| Y_{l_{i} m_{l_{i}}}(\hat{\mathbf{r}}_{\alpha d}) \psi^{(+)}_{l_{i} s J_{i}}(k, r_{\alpha d}) \right\rangle$$

$$= \sqrt{4\pi (2l_{f}+1)} e Z_{eff(L)} \sum_{J_{i}} \sum_{l_{i}} \sum_{m_{l_{i}}m_{s}m_{l_{f}}} i^{l_{i}} \left\langle l_{f} m_{l_{f}} s m_{s} \right| J_{f} M_{f} \right\rangle \left\langle l_{i} m_{l_{i}} s m_{s} \left| J_{i} M_{i} \right\rangle$$

$$\times \left\langle l_{i} 0 s M_{i} \left| J_{i} M_{i} \right\rangle \left\langle l_{f} 0 L0 \left| l_{i} 0 \right\rangle \left\langle l_{f} m_{l_{f}} L M \left| l_{i} m_{l_{i}} \right\rangle R_{l_{f} s L J_{f} l_{i} J_{i}}(k),$$

$$(26)$$

$$R_{l_{f}sLJ_{f}l_{i}J_{i}}(k) = \int_{0}^{\infty} \mathrm{d}r_{\alpha d} r_{\alpha d}^{L+2} I_{l_{f}sJ_{f}}(r_{\alpha d}) \psi_{l_{i}sJ_{i}}^{(+)}(k, r_{\alpha d}).$$
(27)

When deriving Eq. (26) it was taken into account that $\langle \chi_{sm''_s} | \chi_{sm_s} \rangle = \delta_{m''_s m_s}$ and [20]

$$\int d\Omega Y_{l_f m_{l_f}}^*(\hat{\mathbf{r}}_{\alpha d}) Y_{LM}^*(\hat{\mathbf{r}}_{\alpha d}) Y_{l_i m_{l_i}}(\hat{\mathbf{r}}_{\alpha d}) = \sqrt{\frac{(2 \, l_f + 1) \, (2 \, L + 1)}{4 \, \pi \, (2 \, l_i + 1)}} \, \langle l_f \, 0 \, L \, 0 \big| l_i \, 0 \rangle \\ \times \, \langle l_f \, m_{l_f} \, L \, M \big| l_i \, m_{l_i} \rangle \,.$$
(28)

Now we are able to rewrite the expression for the photon differential cross section including all the kinematical factors. If the polarization of the initial and final nuclei (in the case under consideration deuteron and ⁶Li) and of the photon are not measured then the differential cross section takes the form

$$\frac{d\sigma}{d\Omega} = \frac{1}{4} \frac{(2l_f + 1)}{(2J_d + 1)(2J_\alpha + 1)} \frac{(\hbar c)^3}{\mu_{\alpha d} c^2} \frac{k}{E^2} \frac{e^2}{\hbar c} \sum_{M_i M_f} \sum_{J'_i J_i} \sum_{L'L} i^{L'-L} Z_{eff(L')} Z_{eff(L)} \\
\times \sqrt{\frac{(L'+1)(L+1)}{L'L}} \frac{k_{\gamma}^{L'+L+1}}{(2L'-1)!! (2L-1)!!} \sum_{m'_{l_f} m_{l_f}} \sum_{m'_s m_s} \sum_{M' M} \sum_{\lambda = \pm 1} D_{M' \lambda}^{L'}(\varphi, \theta, 0) D_{M \lambda}^{L*}(\varphi, \theta, 0) \\
\times \sum_{l'_i l_i} \sum_{m'_{l_i} m_{l_i}} i^{l_i - l'_i} \left\langle l_f m'_{l_f} s m'_s | J_f M_f \right\rangle \left\langle l_f m_{l_f} s m_s | J_f M_f \right\rangle \left\langle l_f m_{l_f} s m_s | J_f M_f \right\rangle \left\langle l_f m_{l_f} s m_s | J_f M_f \right\rangle \left\langle l_f 0 L' 0 | l'_i 0 \right\rangle \left\langle l_f 0 L 0 | l_i 0 \right\rangle \\
\times \left\langle l_f m'_{l_f} L' M' | l'_i m'_{l_i} \right\rangle \left\langle l_f m_{l_f} L M | l_i m_{l_i} \right\rangle R_{l_f s L' J_f l'_i J_i}^*(k) R_{l_f s L J_f l_i J_i}(k).$$
(29)

Equation (29) can be further simplified taking into account that [23]

$$\sum_{m_{l_f} m_s m_{l_i}} \left\langle l_f m_{l_f} s m_s \big| J_f M_f \right\rangle \left\langle l_i m_{l_i} s m_s \big| J_i M_i \right\rangle \left\langle l_f m_{l_f} L M \big| l_i m_{l_i} \right\rangle$$
$$= (-1)^{s+J_f+l_i+L} \sqrt{(2J_f+1)(2l_i+1)} \left\langle J_f M_f L M \big| J_i M_i \right\rangle \left\{ \begin{array}{c} l_f s J_f \\ J_i L l_i \end{array} \right\}, \tag{30}$$

where $\left\{ \begin{array}{c} l_f \, s \, J_f \\ J_i \, L \, l_i \end{array} \right\}$ is the 6*j*-symbol [23].

Then

$$\frac{d\sigma}{d\Omega} = \frac{1}{4} \frac{(2\,l_f + 1)(2\,J_f + 1)}{(2\,J_d + 1)(2\,J_d + 1)} \frac{(\hbar\,c)^3}{\mu_{\alpha\,d}\,c^2} \frac{k}{E^2} \frac{e^2}{\hbar\,c} \sum_{M_i\,M_f} \sum_{J'_i\,J_i} \sum_{L'L} i^{L'-L} Z_{eff(L')} Z_{eff(L)} \\
\times \sqrt{\frac{(L'+1)(L+1)}{L'L}} \frac{k_{\gamma}^{L'+L+1}}{(2\,L'-1)!!(2\,L-1)!!} \sum_{M'\,M} \sum_{\lambda=\pm 1} D_{M'\,\lambda}^{L'}(\varphi,\,\theta,0) D_{M\,\lambda}^{L^*}(\varphi,\,\theta,0) \\
\times \sum_{l'_i\,l_i} i^{l_i-l'_i} \langle l'_i\,0\,s\,M_i \big| J'_i\,M_i \rangle \langle l_i\,0\,s\,M_i \big| J_i\,M_i \rangle \langle l_f\,0\,L'\,0 \big| l'_i0 \rangle \langle l_f\,0\,L\,0 \big| l_i\,0 \rangle \\
\times (-1)^{l'_i+l_i+L'+L} \sqrt{(2\,l'_i+1)(2\,l_i+1)} \langle J_f\,M_f\,L'\,M' \big| J'_i\,M_i \rangle \langle J_f\,M_f\,L\,M \big| J_i\,M_i \rangle \\
\times \left\{ \begin{array}{l} l_f\,s\,J_f \\ J'_i\,L'\,l'_i \end{array} \right\} \left\{ \begin{array}{l} l_f\,s\,J_f \\ J_i\,L\,l_i \end{array} \right\} R^*_{l_f\,s\,L'\,J_f\,l'_i\,J'_i}(k)\,R_{l_f\,s\,L\,J_f\,l_i\,J_i}(k).$$
(31)

From $\langle J_f M_f L' M' | J'_i M_i \rangle \langle J_f M_f L M | J_i M_i \rangle$ follows that M' = M and [20]

$$D_{M\lambda}^{L'}(\varphi,\theta,0) \left(D_{M\lambda}^{L}(\varphi,\theta,0) \right)^{*} = (-1)^{M-\lambda} D_{M\lambda}^{L'}(\varphi,\theta,0) D_{-M-\lambda}^{L}(\varphi,\theta,0)$$
$$= (-1)^{M-\lambda} \sum_{J} \left\langle L'M \ L - M \big| J0 \right\rangle \left\langle L'\lambda \ L - \lambda \big| J0 \right\rangle D_{00}^{J}(\varphi,\theta,0)$$
$$= -(-1)^{M} \sum_{J} \left\langle L'M \ L - M \big| J0 \right\rangle \left\langle L'\lambda \ L - \lambda \big| J0 \right\rangle P_{J}(\cos\theta)$$
(32)

and

$$\sum_{\lambda=\pm 1} \left\langle L'\lambda \ L - \lambda \middle| J0 \right\rangle = \left\langle L'1 \ L - 1 \middle| J0 \right\rangle \left[1 + (-1)^{L'+L-J} \right]. \tag{33}$$

Then

$$\frac{d\sigma}{d\Omega} = -\frac{1}{4} \frac{(2 \, l_f + 1)(2 \, J_f + 1)}{(2 \, J_d + 1)(2 \, J_\alpha + 1)} \frac{(\hbar \, c)^3}{\mu_\alpha \, d \, c^2} \frac{k}{E^2} \frac{e^2}{\hbar c} \sum_{M_i \, M_f} \sum_{J'_i \, J_i} \sum_{L'L} i^{L'-L} Z_{eff(L')} Z_{eff(L)} \sqrt{\frac{(L'+1)(L+1)}{L'L}} \\
\times \frac{k_{\gamma}^{L'+L+1}}{(2 \, L' - 1)!! \, (2 \, L - 1)!!} \sum_{M} (-1)^M \sum_{J} \langle L'M \, L - M | J \, 0 \rangle \langle L'1 \, L - 1 | J \, 0 \rangle \left[1 + (-1)^{L'+L-J} \right] \\
\times P_J(\cos \theta) \sum_{l'_i \, l_i} i^{l_i - l'_i} \langle l'_i \, 0 \, s \, M_i | J'_i \, M_i \rangle \langle l_i \, 0 \, s \, M_i | J_i \, M_i \rangle \langle l_f \, 0 \, L' \, 0 | l'_i \, 0 \rangle \langle l_f \, 0 \, L \, 0 | l_i \, 0 \rangle \\
\times (-1)^{l'_i + l_i + L' + L} \sqrt{(2 \, l'_i + 1)(2 \, l_i + 1)} \, \langle J_f \, M_f \, L' \, M | J'_i \, M_i \rangle \langle J_f \, M_f \, L \, M | J_i \, M_i \rangle \\
\times \left\{ \begin{array}{l} l_f \, s \, J_f \\ J'_i \, L' \, l'_i \end{array} \right\} \left\{ \begin{array}{l} l_f \, s \, J_f \\ J_i \, L \, l_i \end{array} \right\} R^*_{l_f \, s \, L' \, J_f \, l'_i \, J'_i}(k) \, R_{l_f \, s \, L \, J_f \, l_i \, J_i}(k).$$
(34)

Eq. (34) is quite general and can be applied for the analysis of the photon angular distribution in the direct radiative capture reactions contributed by electric transitions with different multipolarities L or with one dominant L. In Eq. (29) $\hbar c = 197.3$ MeV fm, $e^2/(\hbar c) = 1/137$, $\mu_{\alpha d} c^2$ and E are expressed in MeV, k_{γ} and k are expressed in fm⁻¹. Assuming that only L = 1 or L = 2 contribute one can easily derive differential cross sections for the electric dipole and quadrupole transitions.

Equation (34) can be further simplified for the ${}^{2}\mathrm{H}(\alpha, \gamma)^{6}\mathrm{Li}$ reaction, for which $l_{f} = 0, J_{f} = 1, s = 1, J_{\alpha} = 0$. For this reaction

$$\begin{cases} 0 s J_f \\ J_i L l_i \end{cases} = (-1)^{J_f + L + J_i} \frac{\delta_{s J_f} \delta_{l_i L}}{\sqrt{(2 J_f + 1)(2 L + 1)}}, \\ \begin{cases} 0 s J_f \\ J'_i L' l'_i \end{cases} = (-1)^{J_f + L' + J'_i} \frac{\delta_{s J_f} \delta_{l'_i L'}}{\sqrt{(2 J_f + 1)(2 L' + 1)}},$$
(35)

 $\langle 0 0 L 0 | l_i 0 \rangle = \delta_{l_i L}$ and $\langle 0 0 L' 0 | l'_i 0 \rangle = \delta_{l'_i L'}$.

Then for the differential cross section for the reaction under consideration we get

$$\frac{d\sigma}{d\Omega} = -\frac{1}{12} \frac{(\hbar c)^3}{\mu_{\alpha d} c^2} \frac{k}{E^2} \frac{e^2}{\hbar c} \sum_{M_i M_f} \sum_{J'_i J_i} \sum_{L'L} Z_{eff(L')} Z_{eff(L)} \sqrt{\frac{(L'+1)(L+1)}{L'L}} \\
\times \frac{k_{\gamma}^{L'+L+1}}{(2L'-1)!! (2L-1)!!} \sum_{M} (-1)^M \sum_{J} \langle L'M \ L-M | J0 \rangle \langle L'1 \ L-1 | J0 \rangle \\
\times \left[1 + (-1)^{L'+L-J}\right] P_J(\cos \theta) \langle L'0 \ J_f M_i | J'_i M_i \rangle \langle L0 \ J_f M_i | J_i M_i \rangle \\
\times \langle J_f M_f \ L' M | J'_i M_i \rangle \langle J_f M_f \ L M | J_i M_i \rangle R_{0 \ L' 1 \ J'_i}^*(k) R_{0 \ L 1 \ J_i}(k).$$
(36)

Equations (34) and (36) are our first main result.

B. Total cross sections

The total cross sections can be obtained by integrating the above differential cross sections over the photon's solid angle. Integrating Eq. (34) keeps only the term J = 0 what leads to L' = L. Then

$$\sum_{M_f M} \left\langle J_f M_f L M \middle| J'_i M_i \right\rangle \left\langle J_f M_f L M \middle| J_i M_i \right\rangle = \delta_{J_i J_i},\tag{37}$$

 $\langle LM \ L - M | 0 0 \rangle = (-1)^{L-M} \sqrt{1/(2L+1)}$ and $\langle L1 \ L - 1 | 0 0 \rangle = (-1)^{L-1} \sqrt{1/(2L+1)}$. From $\langle l_f 0 \ L 0 | l_i 0 \rangle$ follows that two subsequent l_i can differ by 2. At astrophysically relevant energies only minimal l_i dominate. Hence we can drop the sum over l_i assuming that each l_i is uniquely determined by L. Also

$$\langle l_i 0 \ s \ M_i | J_i \ M_i \rangle = (-1)^{s+M_i} \sqrt{\frac{2 \ J_i + 1}{2 \ l_i + 1}} \langle J_i \ -M_i \ s \ M_i | l_i \ 0 \rangle$$
(38)

and

$$\sum_{M_i} \left(\langle J_i - M_i \ s \ M_i | l_i \ 0 \rangle \right)^2 = 1.$$
(39)

Taking into account the above results the total cross section reduces to

$$\sigma = 2\pi \frac{(2l_f + 1)(2J_f + 1)}{(2J_d + 1)(2J_\alpha + 1)} \frac{(\hbar c)^3}{\mu_{\alpha d} c^2} \frac{k}{E^2} \frac{e^2}{\hbar c} \sum_{J_i} (2J_i + 1) \sum_L (Z_{eff(L)})^2 \frac{(L+1)(2L+1)}{L} \times \frac{k_{\gamma}^{2L+1}}{((2L+1)!!)^2} (\langle l_f \ 0 \ L \ 0 | l_i 0 \rangle)^2 \left[\left\{ \begin{array}{l} l_f \ s \ J_f \\ J_i \ L \ l_i \end{array} \right\} \right]^2 |R_{l_f \ s \ L \ J_f \ l_i \ J_i}(k)|^2.$$

$$(40)$$

The total cross section for the dipole (quadrupole) transition can be obtained from Eq. (40) by taking L = 1 (L = 2). The total cross section for the reaction under consideration takes the form ($l_f = 0, s = J_f, l_i = L$)

$$\sigma = \frac{2\pi}{3} \frac{(\hbar c)^3}{\mu_{\alpha d} c^2} \frac{k}{E^2} \frac{e^2}{\hbar c} \sum_{J_i} (2J_i + 1) \sum_L \left(Z_{eff(L)} \right)^2 \frac{L+1}{L} \frac{k_{\gamma}^{2L+1}}{\left((2L+1)!! \right)^2} \left| R_{0L1J_i}(k) \right|^2.$$
(41)

Equations (40) and (41) are our second main result.

The astrophysical factor is determined by

$$S(E) = E e^{2\pi \eta_i} \sigma(E).$$
(42)

Here, η_i is the Coulomb parameter in the initial state of the radiative capture process. Replacing $\sigma(E)$ by $\sigma_{Ei}(E)$, where i = 1, 2, we get the astrophysical factors for the dipole (E1) and quadrupole (E2) transitions, correspondingly.

C. Potential model

The most important quantity in calculations of the radiative capture reactions is the radial matrix element $R_{l_f s L} J_{f l_i} J_i(k)$, which is expressed in terms of the the initial and final nuclear wave functions. Different approaches were used to calculate the radial matrix elements. The most frequent used potential approach was based on the pioneering works [24, 25]. In the potential approach the initial scattering wave function is a solution of the Schrödinger equation with the $\alpha - d$ potential, which can be found from the fitting experimental elastic scattering phase shifts in the corresponding partial waves ($l_i = 1, 2$ in the case under consideration). The result is very sensitive to the choice of the final overlap function $I_{l_f s J_f}(r_{\alpha d})$. It was long ago recognized [26] that the ${}^2\mathrm{H}(\alpha, \gamma)^6\mathrm{L}$ reaction is peripheral at astrophysically relevant energies, that is, the overall normalization of the astrophysical factor at Big Bang energies $30 \leq E \leq 400$ keV is practically determined by the square of the ANC $C_{l_f s J_f}$.

In [27] the ⁶Li bound-state wave function was calculated within the framework of the multi-cluster dynamic model. Projection of this bound-state wave function on the two-body channel $\alpha + d$ channel gives the overlap function with correct tail. The two-body potential model was used in [26] to calculate the astrophysical factors for the electric dipole and quadrupole transitions and the total S(E) factor at energies $E \leq 500$ keV. In the two-body potential model the overlap function is replaced by the $\alpha - d$ bound-state wave function:

$$I_{l_f s J_f}(r_{\alpha d}) = S_{n_r l_f s J_f}^{1/2} \varphi_{n_r l_f s J_f}(r_{\alpha d}), \tag{43}$$

where $\varphi_{n_r l_f s J_f}(r_{\alpha d})$ is the α - d two-body bound-state wave function calculated in some phenomenological Woods-Saxon α - d plus Coulomb potential, $n_r = 1$ is the principal quantum number showing the number of the nodes of the radial bound-state wave function at $r_{\alpha d} > 0$. $S_{n_r l_f s J_f}$ is the spectroscopic factor of the configuration α + d in the ground state of ⁶Li. The tail of the bound-state wave function is given by

$$\varphi_{n_r l_f s J_f}(r_{\alpha d}) \stackrel{r_{\alpha d} \gg r_0}{\approx} b_{n_r l_f s J_f} W_{-\eta, l_f + 1/2}(2 \kappa r_{\alpha d}) / r_{\alpha d}, \tag{44}$$

where $b_{n_r l_f s J_f}$ is the single-particle ANC. The value of $b_{n_r l_f s J_f}$ depends on the adopted bound-state potential. The spectroscopic factor $S_{n_r l_f s J_f}$ reflects the fact that the overlap function is not an eigenfunction of any Hamiltonian and, hence, is not normalized to unity, in contrast to the bound-state wave function. Eq. (44) puts limitation on the spectroscopic factor for given $b_{n_r l_f s J_f}$.

The bound-state Woods-Saxon potential should be adjusted to obtain the experimental α – d binding energy (well-depth procedure). However, there are infinite number of such potentials because there are three fitting parameters: geometrical parameters, radius and diffuseness, and the well depth. The final adjustment can be done using the spectroscopic factor. The two-body potential model was also used in [15]. To find the α – d bound-state wave function the Woods-Saxon potential was adjusted to fit the experimental *s*-wave elastic scattering phase shift and to reproduce the experimental α – d binding energy. Since the experimental elastic scattering phase shift includes the many-body effects of the scattered nuclei, the same is true for the two-body potential, which fits the elastic scattering data. Hence, the spectroscopic factor in Eq. (44) should be set to $S_{n_r l_f s J_f} = 1$. However, there is again infinite number of the Woods-Saxon potentials, which differ by the most crucial quantity - the ANC (the inverse scattering problem theorem by Gel'fand-Levitan-Marchenko [28]). The potential adopted in [15] was one of the infinite set of the phase-equivalent potentials with the ANC, which exceeds the experimental ANC [29] and *ab initio* calculations [30] by \approx 18%. Hence, the normalization of the peripheral part of the S(E) factor calculated in [15] exceeded the correct one by \approx 38%. All these questions about ambiguity of the two-body bound-state potentials were addressed in details in [31].

The first full microscopic 6-body approach to calculate the final state ⁶Li bound-state wave function was developed in [19] using the variational Monte Carlo method. The projection of the ⁶Li on the two-body channel α +d has correct tail with the ANC close to the experimental one [29]. The calculated total S(E) factor is in a good agreement with direct measurements around 3⁺ resonance at E = 712 keV.

In hour work, to calculate the photon differential cross sections we used the potential model approach. To calculate the bound-state wave function, two different potentials were used. The first one is the Woods-Saxon potential with the geometrical parameters: radius $r_0 = 1.20$ fm and diffuseness a = 0.7 fm. The square of the single-particle ANC of the bound-state wave function generated by this potential is $b_{1011}^2 = 7.22$ fm⁻¹. To get the correct normalization of the leading asymptotic term of the final-state overlap function $I_{011}(r_{\alpha d})$, that is, the square of the ANC $C_{011}^2 = 5.29$ fm⁻¹, we have to introduce in Eq. (43) the spectroscopic factor $S_{1011} = 0.72$. This method is referred to as M1. The second method is similar to the one described in [31]. In this method, referred to as M2, the Woods-Saxon potential used in [15] was modified to generate the bound-state wave function with correct asymptotic behavior. In this case the spectroscopic factor is $S_{1011} = 1$, that is, the overlap function $I_{011}(r_{\alpha d})$ and bound-state wave function $\varphi_{1011}(r_{\alpha d})$ do coincide at all radii. Thus, both used overlap functions have the same asymptotic behavior being different in the internal region. In both methods the initial α -d scattering wave function is generated by the Woods-Saxon potential from [15]. Its parameters are adjusted to reproduce the experimental phase shifts in the partial waves $l_i = 1, 2$: the radial parameter is r = 1.25 fm, diffuseness a = 0.65 fm, the depth of the potential 56.7 MeV. At $l_i = 2$ this potential reproduces the 3⁺ resonance. To calculate the bound-state and scattering wave functions and the radial matrix elements we used the modified RADCAP code [32].

III. PHOTON ANGULAR DISTRIBUTION IN DIRECT RADIATIVE CAPTURE ${}^{2}H(\alpha, \gamma)^{6}Li$

The calculated photon angular distributions for the ${}^{2}H(\alpha, \gamma)^{6}$ Li direct radiative capture using both methods, M1 and M2, are shown in Fig 1 for 4 different Big Bang energies, E = 70, 100, 200 and 400 keV. As one can see, the dipole differential cross section has the peak at 90°. The quadrupole transition has two peaks, at 45° and 135°. However their interference dramatically changes the angular distribution generating one peak at $\approx 50^{\circ}$. Note that the exact location of the peak slightly depends on the energy. These calculations provide a recipe for the best experimental kinematics. Note that in the experiment performed by LUNA [16] the germanium detector was placed at a 90° angle with respect to the ion beam direction. At this angle the differential cross section is significantly smaller than at the peak value at $\approx 50^{\circ}$.

Another important conclusion is that both methods, M1 and M2, give practically indistinguishable results confirming that at low energies the reaction ${}^{2}H(\alpha, \gamma)^{6}$ Li is completely peripheral. It means that only the tail of the $\alpha - d$ bound-state wave function contributes to the reaction matrix element. Hence, to calculate the reaction matrix element it is enough to use any reasonable bound-state Woods-Saxon potential, which supports *s*-wave $\alpha - d$ bound state with the 1.47 MeV binding energy, and then to introduce a proper spectroscopic factor to provide correct normalization of the asymptotic term of the overlap function.

IV. ASTROPHYSICAL FACTOR

In Fig. 2 the experimental and calculated astrophysical $S_{24}(E)$ factors for the reaction ${}^{2}H(\alpha, d)^{6}Li$ are presented. In contrast to the differential cross section, the total astrophysical factor is given by the sum of the dipole and quadrupole astrophysical factors and does not contain their interference term. The potential model used in the present calculations with two different bound-state wave functions has been described in section II C. The expression for the astrophysical factor has been derived in section II B by integrating the photon's differential cross section over the photon's solid angle. Agreement between the LUNA data at two Big Bang energies and the potential model calculations based on the ANC provides a compelling evidence of the power of the ANC method. Note that the LUNA results are the first direct measurement of the ${}^{2}H(\alpha, \gamma){}^{6}Li$ cross section inside the Big Bang energy range.

Potential model, which we use here, allows us to reproduce the available direct data in the region of the first resonance, E = 0.712 MeV, and even at higher energies. The validity of the potential model at higher energies can be easily explained. At energies higher than Big Bang energies the quadrupole transition dominates. Owing to the presence of the factor $r_{\alpha d}^2$ the quadrupole radial matrix element is dominantly peripheral in the energy interval up to ~ 2 MeV. Hence, the used here potential model with correct normalization of the tail of the overlap function $I_{011}(r_{\alpha d})$ given by the ANC allows one to calculate the astrophysical factor in the broad energy interval $0 \le E \le 2$ MeV.

Note that the calculations from [33] and [15] at Big Bang energies are higher then presented here. For example, at 70 keV, which is the most effective Big Bang energy, $S_{24}(70\text{keV}) = 4.0 \text{ MeV}$ nb [33], $S_{24}(70\text{keV}) = 3.16 \text{ MeV}$ nb [15] and the present result is $S_{24}(70\text{keV}) = 2.58 \text{ MeV}$ nb. The insert in Fig. 2 shows the difference between different calculations of the $S_{24}(E)$ factors in the Big Bang energy interval. At higher energies calculations from [33] reproduce the data quite well while the results from [15] are systematically higher than the data before and after the resonance.

The accuracy of the long wavelength approximation in the case under consideration is quite high: a replacement of r^{L} in the integrand of the radial matrix elements (27) by $O_{1}(r)$, Eq. (19), for L = 1 and $O_{2}(r)$, Eq. (20), for L = 2 changes the astrophysical factor by only $\approx 1\%$. Note that two data points obtained by LUNA were extrapolated in [16] to other energies using calculations in [31]. The calculations in this paper using the method M2 are similar to calculations from [31] but performed with a different, more accurate code [32].

Hence the reaction rates calculated here and in [16] also agree. These reaction rates are significantly lower than the adopted reaction rate from [34] and systematically lower than the reaction rate adopted by NACRE [35]. For example, at $T_9 = 1$, which corresponds to E = 86.2 keV, the adopted NACRE reaction rate exceeds the calculated one in [31] by about 21.5%.



FIG. 1. (Color online) Angular distributions of the photons emitted in the direct radiative capture ${}^{2}H(\alpha,\gamma)^{6}Li$ at E = 70 keV (panel (a)), E = 100 keV (panel (b)), E = 200 keV (panel (c)) and E = 400 keV (panel (d)). All red (olive) lines are obtained using method M1 (M2). The red dashed line (short dashed olive line): the angular distribution calculated for the E1 transition; the red dotted line (short-dotted olive line): the E2 transition; the solid red line (dashed-dotted-dotted olive line): the total photon differential cross section, which is contributed by the sum of the electric dipole and quadrupole terms and their interference term.



FIG. 2. (Color online) Astrophysical $S_{24}(E)$ factors for the ${}^{2}H(\alpha, \gamma)^{6}Li$ reaction. Black dots are data from Ref. [13]; black crosses are data from Ref. [12]; black triangles are data from Ref. [11]. Two blue boxes are the LUNA experimental data reported at E = 94 and 134 keV [16] shown together with their uncertainties. The purple dashed-dotted line is the $S_{24}(E)$ astrophysical factor from Ref. [33]. The black dashed-dotted line is the $S_{24}(E)$ factor from Ref. [15]. All the red (olive) lines are our calculations obtained using model M1 (M2). The red dotted (olive short dotted), red dashed (olive short dashed) and red solid (olive dashed-dotted) lines are the dipole, quadrupole and total $S_{24}(E)$ factors, correspondingly, from the present calculations. Notations in the insert are the same.

V. ⁶Li/⁷Li ISOTOPIC PRIMORDIAL ABUNDANCE RATIO

Evidently that the present paper and LUNA's estimations of the Big Bang abundance of ⁶Li based on the reaction rate of the ²H(α, γ)⁶Li coincide. For the baryon-to-photon ratio 6.047×10^{-10} , which is within the interval determined by the Planck collaboration [36], the calculated primordial abundance of ⁶Li is ⁶Li/H = $(0.74 \pm 0.16) \times 10^{-14}$ [16] which is 34% lower than the abundance given in [34].

In the latest comprehensive analysis of the Big Bang nucleosynthesis the primordial abundance of ⁶Li was determined to be ⁶Li/H = $(0.90-1.77) \times 10^{-14}$ (Planck baryon-to-photon ratio was adopted) [6] and ⁶Li/H = $(1.23-1.32) \times 10^{-14}$ (WMAP baryon-to-photon ratio was taken into account) [37]. As we see, the central values of both results are twice as high as LUNA and present estimations. In both works [6, 37] the nuclear reaction rate from [15] was used claiming that this reaction rate was obtained from the ⁶Li Coulomb breakup. However, it was clearly stated in [15] that the attempt to determine the $S_{24}(E)$ factor from the Coulomb breakup failed and that a potential two-body model was used to calculate $S_{24}(E)$, which turns out to be ~ 30% higher than our and LUNA astrophysical factors [31] because a too large value of the ANC was used in [15]. Hence, the second claim in [6] that the calculated astrophysical factor in [15] and experimental LUNA astrophysical factor [16] "agree well" is also questionable and one of the reasons of high values of the ⁶Li primordial abundance obtained in [6, 37] is that the adopted reaction rates for the ²H(α, γ)⁶Li were based on results from [15].

Thus, by now the primordial abundance of ⁶Li has been established quite accurately. Taking into account the latest estimate of the ⁷Li abundance ⁷Li/H = $(5.1 \pm 0.4) \times 10^{-10}$ obtained from the most recent data on the ³He(α, γ)⁷Be reaction rate [38–40], the resulting isotopic ratio is ⁶Li/⁷Li = $(1.5 \pm 0.3) \times 10^{-5}$ [16]. This isotopic ratio is also the result of the present paper. The obtained from the LUNA experiment and indirect ANC method the Big Bang lithium isotopic ratio is lower than the previous estimates: 2.3×10^{-5} [37] and $(2 - 3.3) \times 10^{-5}$ [6]. However, invoking the reaction rate following from the present paper (or from [31]) and [16] will bring the result obtained in [6, 37] closer to our and LUNA estimations.

The established primordial lithium isotopic ratio is by three orders of magnitude lower then the upper limit determined from the lithium observational data in poor-metal, warm dwarf stars what constitutes the second lithium puzzle. However, the recent publication in Ref. [10] brings a hope that improving the accuracy of the observational ⁶Li data can resolve this puzzle without involving non-standard physics.

VI. SUMMARY

The analysis of the primordial ${}^{2}\text{H}(\alpha,\gamma)^{6}\text{Li}$ reaction is presented. First, the general expression for the angular distribution of the photons and specifically for the reaction under consideration is derived. After that the expressions for the total cross sections for the electric dipole and quadrupole transitions are obtained. The calculated photon's angular distribution, which takes into account the electric dipole and quadrupole transitions and their interference, exhibits the peak at $\approx 50^{\circ}$. It provides a recipe for the best experimental kinematics. Note that at the first direct measurements performed by LUNA [21], the germanium detector was placed at a 90° angle with respect to the ion beam direction, at which the cross section is significantly smaller than at the peak value. New measurements with a better geometry can significantly improve the accuracy of the data. Also the experimental and calculated $S_{24}(E)$ astrophysical factors are presented. Nice agreement between the LUNA data at two Big Bang energies and the potential model calculations based on the ANC proves the power of the ANC method.

The obtained primordial lithium isotopic ratio in [16] and here ${}^{6}\text{Li}/{}^{7}\text{Li} = (1.5 \pm 0.3) \times 10^{-5}$ is a very important result in understanding of the second lithium problem. In resolving this puzzle one needs to reconcile both the Big Bang model prediction of the lithium isotopic ratio and the observational data or to explain their three orders of magnitude difference. The better the accuracy of the Big Bang Li isotopes abundance prediction and the better the agreement with the observational data, the less there will be room for speculations. The results published by LUNA and in this work, ${}^{6}\text{Li}/{}^{7}\text{Li} = (1.5 \pm 0.3) \times 10^{-5}$, sets up quite a strong limit on the primordial isotopic ratio from the Big Bang model. The uncertainty of this ratio is contributed by only 8% uncertainty of the ${}^{7}\text{Li}$ abundance [41] and by 22% uncertainty of the ${}^{6}\text{Li}$ primordial abundance [16]. One of the main conclusions from our work is that the determined optimal kinematics can significantly improve the accuracy of the ${}^{2}\text{H}(\alpha, \gamma){}^{6}\text{Li}$ astrophysical S-factor and, hence, the standard Bing Bang ${}^{6}\text{Li}$ abundance.

But even existing predictions of the Big Bang isotopic ratio ${}^{6}\text{Li}/{}^{7}\text{Li} = (1.5 \pm 0.3) \times 10^{-5}$ puts quite a strong upper limit and much more accurate than the observational data. It looks like such a low value of the Big Bang lithium isotopic ratio makes the second lithium problem even more difficult to resolve. However, in [10] the lithium isotopic analysis in four halo metal-poor stars was revisited using, for the first time, a combined 3D and NLTE modeling technique. This upgraded model systematically reduces the Li isotopic ratio in all four analyzed stars significantly weakening validity of data requiring a significant non-standard primordial ${}^{6}\text{Li}$ production source. Hence, it is too

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