Validation of the ¹⁰Be Ground-State Molecular Structure Using ¹⁰Be $(p, p\alpha)^{6}$ He Triple Differential Reaction Cross-Section Measurements

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state transition was extracted for quasifree angle pairs (θ_p , θ_α) and compared to distorted-wave impulse approximation reaction calculations performed in a microscopic framework using successively the

Tohsaki-Horiuchi-Schuck-Röpke product wave function and the wave function deduced from antisymmetrized molecular dynamics calculations. The remarkable agreement between calculated and measured cross sections in both shape and magnitude validates the molecular structure description of the ¹⁰Be ground-state, configured as an α - α core with two valence neutrons occupying π -type molecular orbitals.

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Introduction.-The formation of structures inside a nucleus is an intriguing phenomenon driven in part by correlations coming from the details of the nucleon-nucleon interaction. Among the different partitioning possibilities within a given nucleus, α clustering has always been considered the most favorable due to the large binding energy of the α particle and its inert character. Consequently, nuclei composed of an integer number of α particles (the socalled self-conjugate nuclei) have initially focused clustering studies. The Ikeda diagram [1], which was proposed at the end of the 1960s, conveys the idea that the cluster degrees of freedom appear in the vicinity of the alpha emission threshold. The famous Hoyle state [2], the second 0⁺ state of ¹²C that plays a key role in the nucleosynthesis of elements heavier than helium is a typical example of such a cluster state [3]. It is located at an energy just above the 3α threshold in ¹²C. Its basic structure in three α particles is established, but its detailed nature is still an object of study. For example, this state can be described as a condensate of α particles occupying a large volume using a wave function of Tohsaki-Horiuchi-Schuck-Roepke (THSR) type [4]. Calculations in the antisymmetrized molecular dynamics (AMD) approach confirm the significant spatial extension of this state and suggest a structure dominated by a "loose" configuration of 3α [5]. In contrast, the ground state of ¹²C is described in the above models as rather compact, with mean-field-like structure, in accordance with the idea of the Ikeda diagram.

As compared to self-conjugate nuclei, the situation is different in neutron-rich light nuclei. For example, strong indications exist that in low-lying states of Be and B isotopes, including the ground state, adding neutrons to an N = Z core leads to spatially extended molecularlike structures in which valence neutrons orbit around the core composed of α particles [6,7]. A typical case is represented by the neutron-rich (N > 4) Be isotopes that were initially described as systems of $2\alpha + Xn$, the two alphas forming a dumbbell-shaped core in the intrinsic frame and X being the number of excess neutrons occupying molecular orbitals around this core [7]. The ground state of ⁹Be, the only stable beryllium isotope, may be described as a 3-body α -n- α molecular structure. The observed rotational band built on the ⁹Be ground-state is well-understood in terms of a π molecular orbital, while the band built on the first $(1/2^+)$ excited state at 1.68 MeV can be connected to a σ -type molecular structure [8-10]. This description of the cluster structure initially elaborated in molecular orbital models was later confirmed by mean-field-type approaches, namely AMD, from which the cluster structures emerge without the existence of clusters being presupposed. The next neutron-rich Be isotope is ¹⁰Be, an unstable nucleus whose structure is also expected to exhibit a marked molecular character. Experimental studies of this nucleus have revealed four rotational bands, corresponding to various cluster structures for excited states [8–14]. However, little is known about the cluster structure of its ground-state, apart from its associated rotational band. Breakup and neutronremoval reactions [15] provided evidence for the existence of ^{*x*}He + $^{A-x}$ He dicluster structures in 10,12,14 Be. The charge radius, which can be measured precisely, is directly related to the density distribution of protons, though not probing directly the cluster structure of ground-states. Among the beryllium isotopes, ¹⁰Be exhibits the smallest charge radius (2.36 fm) [16], which is consistent with AMD calculations that predict a minimum for N = 6 [17]. In these calculations, the ground-state of ¹⁰Be is described as a $2\alpha + 2n$ configuration, in which valence neutrons occupy the molecular attractive π orbital that produces a more compact 2α core, at variance with σ orbitals [6]. It should be noted that this configuration is also predicted in density functional theory calculations [18], a more general framework within which a large number of properties of nuclei can be reproduced. Namely, the dumbbell-shape structure of the alphas as well as the ring-shape π -type orbit of the neutrons emerge from the mean field.

In order to directly probe the spatial extension of α clusters in the ground state of ¹⁰Be and thus establish its overall α cluster molecular structure, we implemented a method based on cluster knockout reactions of the $(p, p\alpha)$ type. These reactions have been studied extensively with proton beams on stable targets from the 1970s until today, and were used essentially to extract the α cluster spectroscopic factors. These spectroscopic factors do not provide direct information on the spatial distribution of α particles in the ground-states of nuclei. On the other hand, the recent theoretical work [19] has demonstrated the high sensitivity of the triple differential cross section (TDX) of the ¹⁰Be($p, p\alpha$) reaction to the α -particle wave function hence to the geometric configuration of the α clusters in the ground-state of ¹⁰Be. In the present work we have performed for the first time the measurement of the TDX of the $(p, p\alpha)$ reaction in inverse kinematics using a beam of unstable nuclei of ¹⁰Be and in kinematical conditions covering the recoilless condition (zero momentum



FIG. 1. Schematic view of ${}^{10}\text{Be}(p, p\alpha)^6\text{He}$ reaction setup.

transfer). The measured TDX was further compared to reaction calculations carried out in a microscopic framework, including in particular the microscopic cluster wave function so as to infer the α -cluster molecular structure of the ¹⁰Be ground-state.

Experiment.-The experiment was performed at the Radioactive Isotope Beam Factory at RIKEN. A secondary beam of ¹⁰Be was produced at an energy of approximately 150 MeV/nucleon through projectile fragmentation of a 230 MeV/nucleon ¹⁸O beam impinging on a 15-mm-thick Be target and purified using the BigRIPS fragment separator [20]. The average ¹⁸O beam intensity was 500 pnA and the produced ¹⁰Be beam intensity was of 5×10^5 particles/second with a purity higher than 90%. ¹⁰Be beam particles were identified on an event-by-event basis. Figure 1 schematically shows the main components of the experimental setup around the secondary target. Beam ions were tracked by a set of two multiwire drift chambers (MWDC), BDC1 and BDC2 placed upstream of the target chamber. To minimize multiple scattering of recoil protons from the ¹⁰Be($p, p\alpha$) reaction in inverse kinematics, a 2-mm-thick pure solid hydrogen target (SHT) [21] was used as the reaction target. Recoil protons were detected using the recoil proton spectrometer (RPS) described in [22,23] in a two-arm configuration set at 60° with respect to the beam axis. Each arm was composed of three stages (MWDC, plastic scintillator, and NaI(Tl) rods) providing position and energy measurements that were used to reconstruct the scattering angle and total energy of the recoil protons. Given the energy range of the recoil protons for the reaction of interest in the present measurement (25-100 MeV), data from the elastic and inelastic channels that produce protons in the relevant energy range were used to perform the energy calibration of RPS. Knocked-out α clusters were measured by two telescopes composed of double-sided strip silicon detector of $62 \times 62 \text{ mm}^2$ active surface backed by CsI(Tl) modules from the FARCOS array [24] placed in the horizontal plane to cover the



FIG. 2. Excitation energy spectra for ${}^{10}\text{Be}(p, p\alpha){}^{6}\text{He}{}^*$ reaction. (a) Integrated over the full solid angle covered by proton and cluster detectors. The vertical dashed line indicates the 2-neutron separation energy (0.975 MeV). The red and blue solid lines show spectra corresponding to events gated by {}^{6}\text{He} and {}^{4}\text{He} residues, respectively. (b) Same for events corresponding to the quasifree angle pair ($\theta_p/\theta_\alpha = 65^\circ/ - 7.7^\circ$) with an angular bin size of $\pm 1^\circ$. See text for details.

angular range 4° - 12° . The scattering angle of helium residues emitted along the beam direction was determined using the MWDC (referred as FDC0 in Fig. 1) placed downstream the SHT, and their identification was performed using the SAMURAI spectrometer and its standard plastic hodoscopes [25].

Results.—Figure 2(a) displays the excitation energy spectra in ⁶He for the ¹⁰Be($p, p\alpha$) reaction obtained from the measured energy and angle of both the recoil proton and α cluster. For the purpose of the energy calibration of the telescopes used for the detection of knocked-out clusters, secondary beams of α particles were produced using BigRIPS in a dedicated run for obtaining the energy calibration while checking the homogeneity of the CsI crystals' response. The present triple coincidence measurement produces spectra with a rather low background. The ground state (g.s.) being the only bound state in ⁶He, the g.s. \rightarrow g.s. transition is easily separated by setting a gate on the ⁶He residues in SAMURAI. The corresponding peak (red histogram) is well fitted by a Gaussian function

centered at (-0.02 ± 0.03) MeV with a missing mass resolution $\sigma = 1.06$ MeV, which validates the calibrations of the proton and α -cluster detectors. The spectrum gated by ⁴He residues in SAMURAI (blue histogram) corresponds to excited states populated in the ¹⁰Be $(p, p\alpha)^{6}$ He^{*} \rightarrow ⁴He + 2*n* reaction channel. The resolution does not allow a clear identification of the populated states. The wellknown 2⁺ state can be seen at 1.797 MeV [26], but not well separated from other low-lying resonant states in the 2–6 MeV region reported in Refs. [26–31]. A less pronounced resonance at around 3.5 MeV is inferred to be the 2⁺₂ state, observed in Ref. [30], and predicted by theoretical calculations [32,33]. Figure 2(b) will be introduced in a later section.

The experimental TDXs for the $(p, p\alpha)$ reaction were extracted for the coplanar angle pairs $(\theta_p, \theta_\alpha)$ chosen to include the zero recoil momentum condition of the residual nucleus. The experimental TDX is given in the unit of $\mu b/(sr^2 \cdot MeV)$ in the laboratory system. For a given angle pair $(\theta_p, \theta_\alpha)$, the TDX for a given transition can be written as

$$\frac{\mathrm{d}^{3}\sigma^{\mathrm{exp}}}{\mathrm{d}T_{p}\mathrm{d}\Omega_{p}\mathrm{d}\Omega_{a}} = \frac{\Delta N(T_{p})}{\varepsilon_{pa}(\theta_{p})N_{t}N_{b}\Delta T_{p}\cdot PV(T_{p})},\qquad(1)$$

where index p and α stand for the outgoing proton and α particle, respectively; T_p is the proton kinetic energy; ΔN is the number of counts in an energy bin ΔT_p ; $\varepsilon_{p\alpha}(\theta_p)$ is the overall efficiency of detecting a p- α pair at θ_p and θ_α obtained from the NPTool simulation [34]; N_t is the number of protons per unit area of the SHT; N_b is the number of incident beam particles; the phase volume term $PV(T_p)$ corresponding to the portion of the $\Delta \Omega_p \times \Delta \Omega_\alpha$ volume kinematically allowed can be defined in discrete form by

$$PV(T_p) = \sum_{\theta_p} \sum_{\varphi_p} \sum_{\theta_a} \sum_{\varphi_a} \sin \theta_p \Delta \theta_p \Delta \varphi_p \sin \theta_a \Delta \theta_a \Delta \varphi_a,$$
(2)

where the summation ranges over θ_p , θ_α , φ_p , and φ_α are restricted to satisfy the energy-momentum conservation.

Figure 3(a) shows the extracted experimental TDXs as a function of the recoil proton kinetic energy T_p for the ${}^{10}\text{Be}(p, p\alpha){}^{6}\text{He}(\text{g.s.})$ reaction at the angle pair $(\theta_p/\theta_\alpha = 65^\circ/ - 7.7^\circ)$ compared with the reaction calculations discussed below. The arrow in the plot indicates the value of T_p^{RL} corresponding to the recoilless condition. The error bars correspond to the square root of the quadratic sum of statistical uncertainties and those on the PV induced by the error on the scattering angles. We note that the shape of the TDX distribution in inverse kinematics is heavily influenced by the PV term shown in Fig. 3(b), which appears in the denominator of Eq. (1). Unlike the TDX



FIG. 3. (a) TDX distribution of ${}^{10}\text{Be}(p, p\alpha){}^{6}\text{He}(\text{g.s.})$ reaction extracted at the coplanar angle pair $\theta_p/\theta_\alpha = 65^{\circ}/-7.7^{\circ}$, which is chosen to include recoilless condition of the residual nucleus. The arrow indicates the proton kinetic energy T_p at the recoilless condition. The red and blue solid lines are the DWIA predictions using the THSR and AMD structural models, respectively. The blue dashed line (dot-dashed line) is the TDX for an artificial state describing the compact shell-model-like (loosely bound gaslike) structure of ${}^{10}\text{Be}$ nucleus [19]. (b) The corresponding phase volume distribution. The horizontal dashed line indicates the product of solid angles $\Delta\Omega_p \cdot \Delta\Omega_{\alpha}$. (c) and (d) are the density distribution of the protons and valence neutrons in the ground state of ${}^{10}\text{Be}$ predicted by the THSR model, adapted from Refs. [19,36].

distribution in forward kinematics for an orbital angular momentum transfer L = 0, it is no longer a peak distribution centered at the value of T_p^{RL} [35].

In Fig. 3(a) the experimental TDX is compared to the result of the distorted-wave impulse approximation (DWIA) calculations using microscopic reduced width amplitude (RWA) of the ground state of ¹⁰Be obtained through the THSR and AMD models. Explicit formulae of $(p, p\alpha)$ TDX calculated within the DWIA framework can be found in Eqs. (4)–(7) of Ref. [37]. See Refs. [38–40] for detailed DWIA description. TDX calculations have been performed for an incident energy of 150 MeV/nucleon used in the experiment. The proton optical potentials were deduced from the democratic parametrization of Dirac phenomenology [41]. The global optical potential of Ref. [42] is applied to the emitted α . As for the *p*- α elementary process, p- α differential cross section obtained by the folding model potential [43] using the Melbourne G-matrix interaction [44] is adopted. For the structure

TABLE I. Comparison of the experimental and theoretical double-differential cross sections for the ground state and 2_1^+ excited-state transitions at quasifree conditions. Both σ_{exp} and σ_{th} are integrated over the angle bin size of $\pm 1^\circ$ for the quasifree angle pairs at $\theta_p = 65^\circ$.

Final state	$\begin{array}{c} [\theta_p/\theta_\alpha] \\ (\mathrm{deg}) \end{array}$	DDX _{exp} (mb/sr ²)	DDX _{THSR} (mb/sr ²)	$\begin{array}{c} DDX_{AMD} \\ (mb/sr^2) \end{array}$
⁶ He(g.s.)	65°/7.7°	23.6(28)	22.7	25.9
⁶ He(2 ⁺ ₁)	65°/7.5°	17.6(30)	5.2	7.9

calculation of the ¹⁰Be ground-state, the RWA used in the TDX calculation has been obtained from the approximation method described in [45] for both THSR and AMD cases. The ¹⁰Be ground-state calculated within the THSR model was presented in Ref. [19]. It corresponds to a molecular configuration of the 2α core with two valence neutrons occupying ring-shape π orbitals. The intrinsic proton and valence density distribution are shown in Figs. 3(c) and 3(d), respectively. The ¹⁰Be ground-state from AMD model [46] exhibits very similar features.

The shape of the experimental TDX distribution of ${}^{10}\text{Be}(p, p\alpha){}^{6}\text{He}(\text{g.s.})$ is very well reproduced by both calculations, as can be seen in Fig. 3. Furthermore, the normalization of the calculated distributions to the experimental one by a fitting procedure leads to normalization factors of 1.04(7) and 0.90(6) for the THSR and AMD, respectively, very close to unity. Given the strong dependence of the TDX on the α -cluster wave function reported in Ref. [19], one can conclude that the present microscopic descriptions of the ${}^{10}\text{Be}(p, p\alpha){}^{6}\text{He}(\text{g.s.})$ reaction allow an accurate reproduction of the data.

The population of excited states in the ⁶He residues through the $(p, p\alpha)$ reaction measures the contribution of ⁶He core-excited states in the ground state of ¹⁰Be. Extracting the corresponding cross sections in addition to the one for the ground-state transition provides a stringent test to the structure models of ¹⁰Be. Table I gives the extracted double-differential cross sections (DDX) for the ground-state and 2_1^+ excited-state transitions at the quasifree condition obtained by integrating the TDX over the proton kinetic energy T_p . Because of the slight variation of the residue mass, there is a corresponding change in the quasifree angle pairs, while $\theta_p = 65^\circ$ remains unchanged for both cases. The number of counts for the g.s. transition is extracted straightforwardly as only the g.s. of ⁶He is bound, while that for the 2^+_1 excited state transition is obtained through a decomposition of the excitation energy spectrum, as shown in Fig. 2(b). The spectrum was fitted by two resonances modeled as the convolution of a Breit-Wigner distribution with a Gaussian function, taking into account the experimental resolution. The amplitudes and energies (E) of the two resonances are treated as free parameters, while keeping the width of the 2_1^+ state (0.113 keV) fixed. Two approaches for handling the width of the additional resonance were explored: energy-dependent $\Gamma(E)$ from the *R*-matrix method [47,48] and the approach described in Ref. [49] [shown in Fig. 2(b)], yielding consistent results. The present spectrum is well reproduced by a fit with a resonance identified to the well-known 2_1^+ state and an additional resonance at 5.0(2) MeV, $\Gamma = 2.3(5)$ MeV, consistent with Refs. [27,31].

The experimental and calculated DDXs are compared in Table I. A very good agreement is obtained for the groundstate transition as expected from the TDX comparison in Fig. 3(a). The DDXs of the transition to the 2_1^+ state are smaller than for the ground-state transition, in particular for the theoretical predictions. The discrepancy between the experimental and theoretical values may arise from the inadequate treatment of time-dependent resonances in the structure models and the reaction models. The present THSR and AMD calculations described the 2_1^+ resonant state with boundary conditions as those of the bound states, resulting in lower predictions for cross sections due to the different matrix elements between initial and final states.

Discussion.—The α -cluster molecular structure of the ¹⁰Be ground-state within the THSR-based framework shown above $(2\alpha \text{ cores with the two valence neutrons})$ occupying ring-shape π orbitals) allows the reproduction of our experimental cross sections very well. This configuration is spatially extended, although to a lower extent than for other beryllium isotopes because of the attractive effect of the π neutrons compared to, e.g., neutrons in σ orbits. Consistently, the corresponding root-mean-square charge radius of ¹⁰Be is 2.31 fm, very close to the experimental value. To estimate the uncertainty from potential choices on cross sections, we performed further TDX calculations using the Koning-Delaroche potential for $p - {}^{10}\text{Be}$ and $p - {}^{6}\text{He}$ [50] and the $\alpha - {}^{6}\text{He}$ optical potential was obtained by the nucleon-nucleus folding calculation using the Koning-Delaroche potential and the phenomenological alpha density [51]. Both sets of potentials reasonably reproduced the experimental TDX, with the later calculation showing scaling factors of approximately 80% compared to the original calculation.

The sensitivity of the TDX of the ${}^{10}\text{Be}(p, p\alpha){}^{6}\text{He}(\text{g.s.})$ reaction to the spatial extension of the α -cluster wave function, quantified by the intercluster distance has been clearly demonstrated in [19], and the present results thus provide a direct evidence of the above molecular structure. To confirm that these conclusions are valid in our present experimental conditions (incident energy, finite angular acceptances), we performed the TDX calculation using the two "extreme" cases of a compact shell-model-like and a loosely bound gaslike configuration of ${}^{10}\text{Be}$ nucleus described in Ref. [19]. Although unphysical, these states test the impact of α clusters with different spatial distribution on the TDX magnitude. As can be seen in Fig. 3(a),

very large normalization factors are needed to match the magnitude of the data, far beyond the cross-section calculation uncertainty. It can be also noted that the TDX calculation corresponding to the gaslike configuration has a different shape compared to the experimental data. Altogether, our measurement validates the microscopic physical THSR α -cluster wave function mentioned above.

The AMD approach has been successfully used to describe and establish low-lying molecular structures in light nuclei. This framework provides a microscopic description of both single-nucleon properties as well as cluster structure, without assuming the cluster *a priori*. Within this framework, the ¹⁰Be ground-state is found to have a structure similar to the one extracted from the THSR model wave function in the region of interest. Namely, the RWA has similar behavior at the surface region that contributes to the cross section. Consistently, we find that the calculated TDX using AMD RWA shows an agreement of similar quality with the data as when using the THSR approach.

Conclusion.—The α -cluster structure of an unstable neutron-rich nucleus, ¹⁰Be, has been investigated by measuring for the first time the TDX of the $(p, p\alpha)$ reaction in inverse kinematics with a setup allowing inclusion of the recoilless condition. Double-differential cross sections to the ground and 2_1^+ states of the ⁶He residue have been extracted independently. Obtained data have been compared with cross-section calculations performed within a microscopic DWIA framework involving up-to-date α -cluster wave functions, describing the ground-state in terms of a dumbbell-shape 2α core (with moderate extension) surrounded by the two valence neutrons occupying the π orbit. There is a remarkable agreement in both shape and magnitude between the experimental and calculated TDX for the ${}^{10}\text{Be}(p, p\alpha){}^{6}\text{He}(\text{g.s.})$. Because of the previously established sensitivity of the TDX to the extension of the α wave function in the ground-state of ¹⁰Be, our results provide direct experimental evidence of the above molecular structure of ¹⁰Be implemented in the THSR approach, and validated by the general AMD framework. Concerning the doubledifferential cross sections to the g.s. and 2^+_1 states of the residue, a good agreement between calculations and experimental results is found. A consistent picture is then obtained. In the near future, based on the present work, further studies will lead to the understanding of the evolution of the α -cluster structure in the ground-state of neutron-rich nuclei with increasing number of valence neutrons. Besides, it is also planned to apply knockout reactions in inverse kinematics to investigate the formation of other types of clusters in the ground-state of nuclei away from the stability valley.

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