

Extensive Study of the Wobbling Properties in ^{163}Lu for the Positive and Negative Parity States

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Abstract

A new interpretation on the wobbling structure in ^{163}Lu is developed, based on the concept of parity symmetry. It is known that four wobbling bands are experimentally observed in this isotope, where three of them are considered as wobbling phonon excitations (namely TSD_2 , TSD_3 , and TSD_4) and the yrast band for the ground state (that is TSD_1). In the present work, the trial function that is used for obtaining the wobbling spectrum is analyzed in terms of its behavior under the rotation operation. Indeed, due to a specific symmetry to rotations with π around the 2-axis of the triaxial system, the parity becomes a good quantum number. As such, the trial function admits solutions with negative parity, which belong to the rotational states in TSD_4 . A unified description of all the triaxial super-deformed bands in ^{163}Lu is achieved with the new formalism.

1 Introduction

Triaxiality in nuclei has become an interesting topic for physicists over the years, mainly due to its great challenge of measure it experimentally, but also for its large number of characteristics that are said to be resulting from these kind of shapes. Moreover, stable triaxial shapes are of rare occurrence across the chart of nuclides [1], since the predominant character of nuclei is either spherical or axially symmetric. Over the last two decades, it has been shown that triaxiality plays a crucial role in measurements of important quantities like proton emission probabilities [2], separation energies of the nucleons [1], and also fission barriers in heavy nuclei [3], however, concrete evidences

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of triaxiality in nuclei were still missing or under investigation. A tremendous work was given in finding a clear signature for non-axially symmetric shapes: effects such as anomalous signature splitting [4], signature inversion [5], and staggering of γ bands [6] were pointed, but only recently two clear fingerprints of nuclear triaxiality have emerged in the literature, based on both experimental and theoretical findings. Indeed, the phenomena of *chiral symmetry breaking* and that of *wobbling motion* (W.M.) are considered as unique characteristics of nuclear triaxiality.

Chirality consists in the existence of a pair of chiral twin bands with an identical structure and almost similar energies. These bands are expected to appear due to the coupling of valence nucleons and the collective mode of rotation that could drive the total spin away from any of the three principal planes, giving rise to both left-handed and right-handed orientation of the angular momentum vectors [7]. A rigorous investigation of all the nuclei with chiral bands is given by Xiong and Wang in [8], where reportedly a total of 59 chiral doublet bands in 47 such nuclei are confirmed. As a matter of fact, 8 of these nuclei have multiple chiral doublets.

On the other hand, the experimental observations regarding wobbling motion have been quite rare, even though this kind of collective motion has been theoretically predicted almost 50 years ago by Bohr and Mottelson [9] when they were investigating the rotational modes of a triaxial nucleus by means of a Triaxial Rotor Model (TRM). Therein, they showed that for a triaxial rotor, the main rotational motion is around the axis with the largest moment of inertia (MOI), as it is energetically the most favorable. This mode is quantum mechanically disturbed by the rotation around the other two axes, since rotation around any of the three principal axes of the system are possible, due to the anisotropy of the three different MOIs (that is $\mathcal{I}_1 \neq \mathcal{I}_2 \neq \mathcal{I}_3$).

W.M. can be viewed as the quantum analogue to the motion of the asymmetric top, whose rotation around the axis with largest MOI is energetically favored and stable. A uniform rotation about this axis will have the lowest energy for a given angular momentum (spin). As the energy increases, this axis will start to precess with a harmonic type of oscillation about the space-fixed angular momentum vector, giving rise to a family of wobbling bands, each characterized by a wobbling phonon number n_w . The resulting quantal spectrum will be a sequence of rotational $\Delta I = 2$ bands, with alternating signature number for each wobbling excitation. According to [9], it is possible to obtain the wobbling spectrum of any triaxial rigid rotor, by using the information related to its angular momentum I , moments of inertia $\mathcal{I}_{1,2,3}$, rotational frequency ω_{rot} , wobbling frequency ω_{wob} as follows:

$$E_{\text{rot}} = \sum_i \left(\frac{\hbar^2}{2\mathcal{I}_i} \right) I_i^2 \approx \frac{\hbar^2}{2\mathcal{I}_1} I(I+1) + \hbar\omega_{\text{wob}} \left(n_w + \frac{1}{2} \right), \quad (1)$$

with ω_{wob} given by the following expression:

$$\hbar\omega_{\text{wob}} = \hbar\omega_{\text{rot}} \sqrt{\frac{(\mathcal{I}_1 - \mathcal{I}_2)(\mathcal{I}_1 - \mathcal{I}_3)}{\mathcal{I}_2\mathcal{I}_3}}, \quad (2)$$

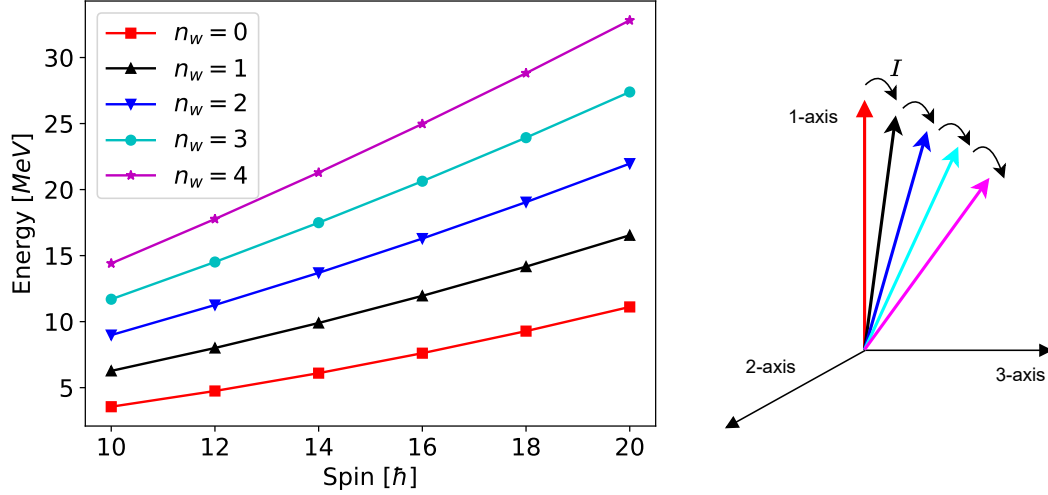


Figure 1: Family of wobbling bands for a simple triaxial rotor (left-side). Tilting of the angular momentum vector away from the rotational axis (right-side). This schematic representation was done for an arbitrary set of MOIs $\mathcal{I}_1 : \mathcal{I}_2 : \mathcal{I}_3 = 25 : 5 : 2$.

where the rotational frequency of the rigid rotor is given by $\hbar\omega_{\text{rot}} = \frac{\hbar I^2}{\mathcal{I}_1}$. In Eq. 1, the approximation of very large MOI along 1-axis is considered (i.e., $\mathcal{I}_1 \gg \mathcal{I}_2, \mathcal{I}_3$), and $I(I+1) = \mathcal{I}_1^2 + \mathcal{I}_2^2 + \mathcal{I}_3^2$. One can see that the wobbling motion is expressed as a 1-dimensional vibration with only one variable, since the energy of the zero-point fluctuation is $\frac{\hbar\omega_{\text{wob}}}{2}$ [10].

Just for an illustrative purpose, Figure 1 shows a theoretical spectrum for the wobbling bands within a triaxial rigid rotor. The family of wobbling bands are obtained from a set of three moments of inertia (along the three principal axes), a given angular momentum, and different wobbling phonon numbers. Moreover, in Figure 1, the tilting of the angular momentum away from the rotational axis is sketched, where the tilt increases with the increase in the wobbling excitation. In a given sequence of wobbling bands, both the intra-band $\Delta I = 2$ as well as inter-band $\Delta I = 1$ transitions have a strong $E2$ collective character.

It is important to mention that the wobbling spectrum described by Eq. 1 and graphically represented in Figure 1 was firstly predicted for an even-even triaxial nucleus [9]. This predicted wobbling mode has not been experimentally confirmed yet. However, the first experimental evidence for wobbling excitations in nuclei was for an even-odd nucleus, namely ^{163}Lu , where a single one-phonon wobbling band was measured initially [11], followed by two additional wobbling bands discovered one year later [12, 13].

After the first discovery of wobbling bands in ^{163}Lu ($Z = 71$), an entire series of even-odd isotopes with $A \approx 160$ were experimentally confirmed as *wobblers*: ^{161}Lu , ^{165}Lu , ^{167}Lu , and ^{167}Ta . In these nuclei, the wobbling mode appears due to the coupling of a valence nucleon (the so-called $\pi(i_{13/2})$ intruder) to a triaxial core, driving the entire nuclear system up to large deformation ($\epsilon \approx 0.4$) [14].

With time, several nuclei in which WM occurs were also reported in regions of smaller A . Indeed, two isotopes with $A \approx 130$: ^{133}La [15] and ^{135}Pr [16, 17] were identified as having wobbling bands, which emerged from the coupling with a triaxial even-even core of another intruder (the $\pi(h_{11/2})$ nucleon) for ^{135}Pr , and an additional pair of positive parity quasi-protons which are making an alignment with the short axis of the triaxial rotor for ^{105}Pr . The resulting coupling in both cases have a deformation $\epsilon = 0.16$ [15, 16], which is obviously smaller than the deformation in the heavier nuclei within the $A \approx 160$ region. A third nucleus that also lies in this mass region was confirmed very recently by Chakraborty et. al. in [18], namely the odd- A ^{127}Xe , where a total of four wobbling bands have been reported by the team (two yrast bands, and two excited phonon bands with $n_w = 1$ and $n_w = 2$).

Some additional progress towards a more comprehensive wobbling spectroscopy was made in the $A \approx 100$ mass region, with an experimental evidence for ^{105}Pd that showed of two such bands that are built on a $\nu(h_{11/2})$ configuration, the first one so far in which a valence neutron couples to the triaxial core [19]. The resulting configuration drives the nuclear system up to deformation $\epsilon \approx 0.26$.

The heaviest nuclei known so far in which WM has been experimentally observed are the isotopes $Z = 79$ with $A = 183$ [20] and $A = 187$ [21], respectively. However, for the case of ^{187}Au , there is an ongoing investigation [22] whether the two wobbling bands ($n_w = 0$ and $n_w = 1$) are bands with wobbling character, or if they are of magnetic nature (which would exclude the wobbling phonon interpretation).

Regarding the wobbling motion for the even-even nuclei (behavior that was described above through the schematic representation from Figure 1), the experimental results are very fragmentary, with unclear evidence on such collective behavior in nuclei. However, some embryos of even-even wobblers have been reported in the recent years. For example, the ^{112}Ru ($Z = 44$) nucleus has three wobbling bands [23], with two of them being excited (one- and two-phonon wobbling bands). Another example is the even-even ^{130}Ba ($Z = 56$) [24–26]. Indeed, for ^{112}Ru , the ground band together with the odd and even spin members of the γ band with were interpreted as zero-(yrast), one-, and two-phonon wobbling bands. Unfortunately, since there are no data concerning the electromagnetic transitions, its wobbling character is still unclear. On the other hand, for the nucleus ^{130}Ba , from its recent study regarding the band structure [24], a pair of bands with even and odd spins were proposed as zero- and one-phonon wobbling bands, respectively. What it is worth noting for this case is the fact that these two bands are built on a configuration in which two aligned protons that emerge from the bottom of $h_{j=11/2}$ shell couple with the triaxial core. One remarks the change in nature of the wobbling motion from a purely collective form, but in the presence of two aligned quasiparticles [25].

Regarding the interpretation of the wobbling motion which occurs in the nuclei that were mentioned above, it is mandatory to discuss some aspects related to its behavior with the increase in total angular momentum (nuclear spin). It is a long lasting debate on whether certain nuclei behave as *longitudinal wobblers* (LW) or *transverse wobblers* (TW). The concepts of LW and TW emerged from an extensive study done by Frauendorf et. al. [27] in which the team discussed the possible coupling schemes that a valence

nucleon can create with the triaxial core, thus giving rise to two possible scenarios. Based on microscopic calculations using the Quasiparticle Triaxial Rotor (QTR) model, they showed that if the odd valance nucleon aligns its angular momentum vector \vec{j} with the axis of largest MOI, the nuclear system is of longitudinal wobbling character. On the other hand, if the odd nucleon aligns its a.m. vector \vec{j} with an axis perpendicular to the one with largest MOI, then the nuclear system has a transverse wobbling character. From the microscopic calculations, it was shown that for LW, the wobbling energy E_{wob} (see Eq. 3) has an *increasing* behavior with an increase in spin, while for TW the energy E_{wob} *decreases* with spin.

Within the nuclei that were mentioned above, most of them are of TW type, with only ^{133}La [15], ^{127}Xe [18], and $^{183,187}\text{Au}$ [20, 21] being nuclei with LW character. The energy that characterizes the type of wobbling in a nuclear system is the energy of the first excited band (the one-phonon $n_w = 1$ wobbling band) relative to the yrast ground band (zero-phonon $n_w = 0$ wobbling band):

$$E_{\text{wob}} = E_1(I) - \left(\frac{E_0(I+1) + E_0(I-1)}{2} \right), \quad (3)$$

with 0 and 1 representing the wobbling phonon number n_w .

The odd nucleons that couple with the rigid triaxial core will influence the appearance of a particular wobbling regime (LW or TW). In all the wobblers, there is a proton from a certain orbital which is coupling with the core, except for the case of ^{105}Pd , where the valence nucleon is a neutron. The nature of the odd quasiparticle (i.e., particle or hole) and its "position" in the deformed j -shell (i.e. bottom or top) will determine whether its angular momentum \vec{j} will align with the *short* (s) or *long* (l) axes of the triaxial rotor, respectively (with the notations short s , long l , and medium m axes of a triaxial ellipsoid). The reasoning behind this has to do with the minimization of the overall energy of the system: in the first case, a maximal overlap of its density distribution with the triaxial core will determine a minimal energy, while in the second case, a minimal overlap of the density distribution of the particle with the core will result in a minimal energy. Moreover, if the quasiparticle emerges from the middle of the j -shell, then it tends to align its angular momentum vector \vec{j} with the *medium* (m) axis of the triaxial core. Figure 2 aims at depicting the type of alignment of a quasiparticle with the triaxial core.

As previously mentioned, for a given angular momentum, uniform rotation around the axis with the largest MOI corresponds to a minimum energy. For a triaxial rotor, this is equivalent to rotation around the m axis. Therefore, Frauendorf [27] classified the LW as the situation when the odd nucleon will align its angular momentum along the m -axis, while TW being the situation where j is aligned perpendicular to the m -axis (with s - or l -axis alignment depending on the j -shell orbital from which the odd nucleon arises). It is worthwhile to mention the fact that the analysis done in Ref. [27] was within a so-called *Frozen Alignment* approximation, where the angular momentum of the odd particle \vec{j} is rigidly aligned with one of the three principal axes of the triaxial ellipsoid (that is s -, l - or m -axis).

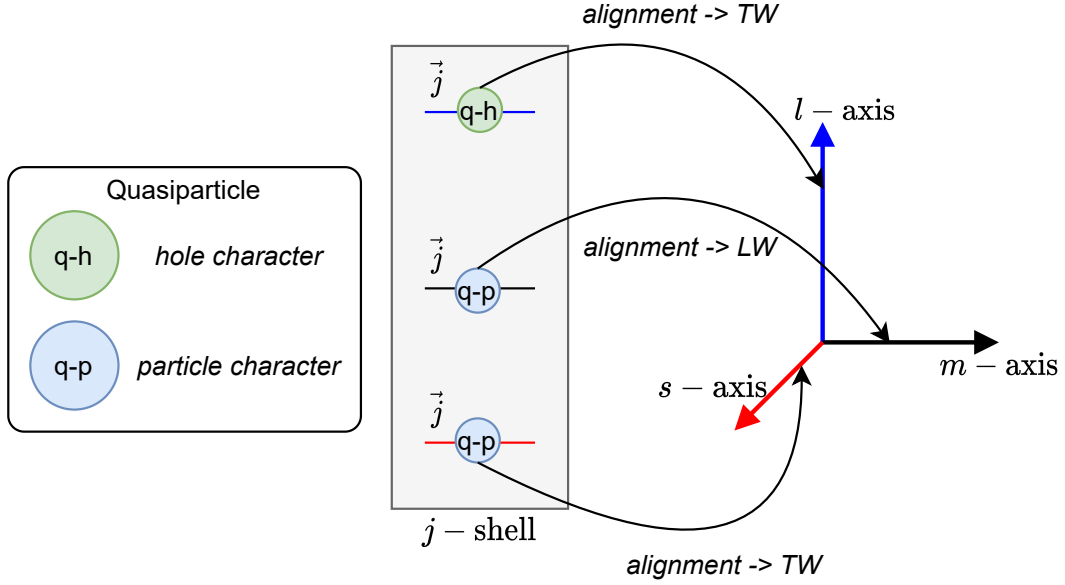


Figure 2: The wobbling regime (LW or TW) based on the type of alignment for an odd quasiparticle with the triaxial core. Figure based on the quantal analysis done in [27].

For a better understanding of the wobbling regimes in terms of angular momentum alignment, the schematic illustration from Figure 3 depicts three particular cases, namely a simple wobbler (the case firstly developed by Bohr and Mottelson [9]) - shown in inset A.0, a longitudinal wobbler - shown in inset A.1, and lastly a transverse wobbler - shown in inset A.1.

In terms of its theoretical analysis, the wobbling motion has been studied using multiple models and interpretations. The Triaxial Particle Rotor Model has been widely used over the recent years [9, 27–30], these being quantal models that can be exactly solved in the laboratory frame. TRM was however, firstly introduced for the motion of a rotating nuclear system by Davydov and Filippov in [31], where they obtained a complete quantal description for the motion of a triaxial nucleus (considering the fact that the nucleus must have a well-defined potential minimum at a non-zero value for the triaxiality parameter γ). Starting from the framework of Cranking Mean Field Theory (CMFT), there were attempts at extending the cranking model for the study of WM. however, using the mean field approximations, CMFT only help at describing the yrast sequence for a given configuration. In order to improve that, the framework was extended with proper quantum correlations by incorporating the Random Phase Approximation (RPA) theory (see Refs. [32–39] for more details). The method of Collective Hamiltonian [40, 41] was used for the investigation of wobbling spectra in nuclei with the help of deformed potentials which were calculated from the Tilted Axis Cranking (TAC) model. TAC single j -shell model is also used for the description of the chiral vibrations and rotational motion in deformed nuclei [42, 43]. Mean field approxima-

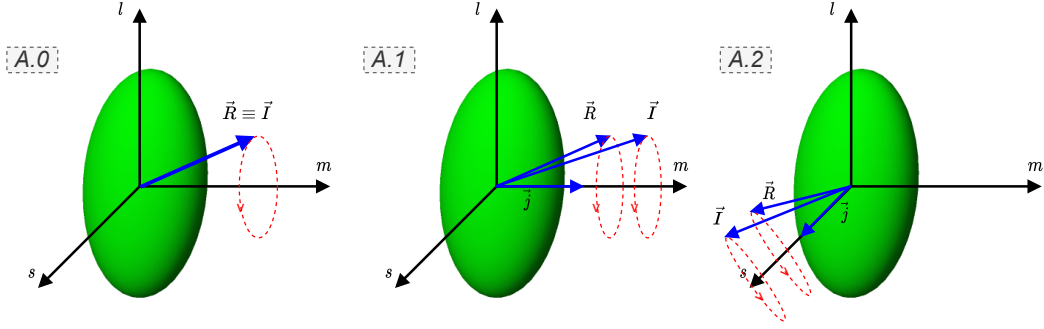


Figure 3: The geometry for the angular momentum for a simple wobbler: A.0, a longitudinal wobbler: A.1, and a transverse wobbler: A.2. The short, long, and medium axes are defined in the body-fixed frame. The vectors \vec{R} , \vec{j} , and \vec{I} represent the set of angular momenta of the core, odd particle, and the total nuclear system, respectively.

tions were also developed by the so-called *generator coordinate method after angular momentum projection* (GCM+AMP for short), with calculations that emerged from intrinsic cranking states [44]. Some analytical solutions were also developed (based on certain approximations), such as the harmonic approximation (HA) [9, 27, 40, 45], Dyson boson expansion [45, 46], and Holstein-Primakoff (HP) formula [29, 45–48]. The angular momentum projections were also incorporated into the mean field framework, with the recent development of a completely microscopic description of the wobbling motion by Shimada et. al. [49]. A Projected Shell Model (PSM) [50] which starts from the shell-model configuration mixing that is based on a Nilsson deformed mean field was also used for the theoretical study concerning WM. There are alternative developments based on the PSM approach, based on Density Functional Theories (DFT) that can be both non-relativistic [51] as well as relativistic [52].

Other tools that proved to be very efficient for the analysis of the wobbling nuclei are the semi-classical approaches, through which one can obtain equations of motion that describe the nuclear system quite well, starting from quantal Hamiltonians and further applying some de-quantization procedures. The semi-classical approach applied to generalized rotor Hamiltonians has the *advantage* of keeping close contact with the classical picture embedded in the dynamic of the systems. Recently, there has been quite an impressive progress towards realistic description of the wobbling motion [27, 45, 53–57]. As a matter of fact, the present team was able to describe (with a very good agreement) the experimental data concerning the wobbling energies and electromagnetic (e.m.) transition probabilities for the Lu isotopes with $A = 161, 163, 165, 167$ (see the work done in Refs. [45, 54]), and more recently the odd- A ^{135}Pr isotope [46]. Indeed, starting from a quantal Hamiltonian specific to a triaxial rotor model (that is a triaxial core coupled with an odd valence nucleon) and applying the Time Dependent Variational Equation (TDVE), with a trial function that was carefully chosen, the complete wobbling spectrum of the mentioned isotopes was reproduced, together with the e.m. (intra-band and inter-band) transitions.

Concluding this section, the importance of nuclear triaxiality and the challenges of identifying it experimentally were contoured in the beginning, serving as the starting point of the current work. Furthermore, all the known nuclei in which wobbling motion appears were mentioned (with important observation for some of them). Additionally, the mechanism behind the simple wobblers (the one developed by Bohr and Mottelson [9]) was sketched (starting with Eq. 1), and a family of wobbling bands were schematically represented for a given set MOIs associated to the triaxial rotor (see Figure 1). Lastly, a brief overview with most of the theoretical *tools* that were/are used for describing this elusive phenomenon was realized. Having this in mind, one can say that a detailed outlook for the topic of wobbling motion was properly shown.

Going further, the remaining structure of this current work must be pointed out. In Section 2, an overview with regards to the team's reinterpretation of the wobbling band structure in ^{163}Lu will be illustrated. This will be the *core-idea* that serves as foundation of this newly developed model. The theoretical formalisms and analytical formulas will be properly presented in Section 3. Experimental results concerning the wobbling spectrum of this isotope will be compared with the newly obtained data in Section 4. Overall conclusions and discussions are reserved to Section 5.

2 Re-interpretation of the wobbling bands structure for ^{163}Lu

Now that a complete overview of the recent experimental and theoretical results regarding wobbling motion has been made, together with the description of its two regimes (namely, longitudinal wobbling and transverse wobbling), it is worth mentioning the latest progress made by the present team towards the actual interpretation of the wobbling structure of ^{163}Lu . Considered the *best wobbler* to date, ^{163}Lu has a rich wobbling spectrum [11,12], with no less than four such wobbling bands: one yrast - TSD_1 , which has a zero-phonon wobbling number $n_w = 0$), and three excited wobbling bands - $TSD_{2,3,4}$ with their corresponding wobbling phonon numbers $n_w = 1, 2, 3$, respectively. The name TSD comes from Triaxial Strongly Deformed bands. The triaxial bands emerge due to the coupling of an odd- \vec{j} nucleon with an even-even triaxial core. Thus, for ^{163}Lu , it is the intruder $\pi(i_{13/2})$ that couples to the triaxial core [11,13,28], driving the nuclear system up to large deformation, and stabilizing the deformed structure. Indeed, a triaxial shape with deformation parameters $(\epsilon_2, \gamma) \approx (0.38, +20^\circ)$ is assumed to be in agreement with the observed data, based on calculations using the Ultimate Cranker Code [58] for the potential energy surface (PES).

In terms of experimental evidence which should be pointing out wobbling nature for the four TSD bands belonging to ^{163}Lu , the large transition quadrupole moment $Q_t \approx 10\text{ b}$ [59] (which is substantially larger than it is the case for normal-deformed bands), the predominantly $E2$ character of the transitions linking adjacent bands ($I \rightarrow I - 1$), a large $E2/M1$ mixing ratio $\delta > 1$ for the transitions linking the yrare ($n_w = 1$) and yrast ($n_w = 0$) bands (which obviously should result in smaller transitions with magnetic character, or in other words, there is no $M1$ dominance) are clear fingerprints

of wobbling nature. For a set of results concerning these quantities (both theoretical and experimental), see Ref. [45], and the references cited therein. Another quantity that indicate strong deformation with wobbling character is the relative rigid rotor energy, and for this isotope calculations show that all four bands have a similar behavior with respect to this value (see Figures 3 and 4 from Ref. [60]).

Considering the experimental evidence which was indicated above and calculations based on particle rotor models, it can be summarized that the *generally accepted* formalism for the band structure in ^{163}Lu is the following:

- There are three excited wobbling bands (w.b.): TSD_2 , TSD_3 , and TSD_4 .
- The three excited w.b. have wobbling-phonon numbers $n_{w_2} = 1$, $n_{w_3} = 2$, and $n_{w_4} = 3$, respectively.
- All three bands are built on top of the yrast state (the ground state band) with zero-wobbling-phonon number $n_{w_1} = 0$.
- Stable triaxial super-deformation is achieved due to the alignment of the odd $\pi(i_{13/2})$ nucleon which couples to a triaxially deformed core \vec{R} .
- $TSD_{1,2,3}$ have all positive parity $\pi_1 = \pi_2 = \pi_3 = +1$, while the spin states belonging to TSD_4 have negative parity $\pi_4 = -1$. All states within the four bands have half-integer spin, obviously.

In accordance with the band structure which was just formulated, a full description of the wobbling spectrum of ^{163}Lu was done within a semi-classical formalism by Raduta et. al. [45]. Therein, with the TDVE applied on the PRM Hamiltonian with a trial wave-function that encapsulates both the states of the deformed nucleus I and the single-particle states j , a set of analytical expressions for the excitation energies of all four bands was obtained. The energies belonging to the excited wobbling phonons were populated by the action of a phonon operator Γ^\dagger on the ground state. Indeed, by acting with the operator one on the ground state with the spin $I = R + j$ and $R = 0, 2, 4, \dots$, the states from TSD_2 ($n_w = 1$) can be obtained. By applying twice ($n_w = 2$) the phonon operator, the rotational states from TSD_3 will be created. Lastly, the states from TSD_4 are obtained with the action on the ground state with three ($n_w = 3$) phonon operators: two of positive parity and one of negative parity (due to the overall negative parity $\pi_4 = -1$ of TSD_4). One has to remark the fact that for TSD_4 , the model assumes an odd-particle-rotor with a different intruder: the $\pi(h_{9/2})$ nucleon. This was suggested by the negative parity orbital which might be occupied by this proton, in the spherical shell model. Several calculations in the literature point out that this nucleon might be causing the third excited wobbling band to have negative parity [61]. It is worthwhile mentioning that for the work described in [45], the variational principle was only applied for the states in TSD_1 , since the other three wobbling bands are obtained through phononic excitations with the corresponding operator.

In what follows, it is useful to introduce some notations that will refer to the formalisms developed by the team in describing the wobbling motion in ^{163}Lu . As such,

the work developed in [56, 57] will be denoted to W1, while the current work (which is in fact an extension of W1) will be shortly denoted by W2. For the sake of a self-consistent presentation, in the following subsection 2.1 a brief overview of the recently published work W1 will be made, with further development that has W1 as a starting ground being presented in the second subsection 2.2 - representing the *core concept* of the current team's investigation.

2.1 W1 - Signature Partner Bands

Working with a semi-classical approach that is based on the triaxial particle rotor model, a full description of the wobbling bands for ^{163}Lu was achieved, but with a slightly modified band structure. Indeed, rather than applying a TDVE just for the yrast TSD_1 band, the states from TSD_2 were also obtained variationally. This was possible due to the different coupling schemes that emerged for TSD_1 and TSD_2 , respectively. More precisely, in [57] and [56] there are three different coupling schemes ($\vec{R} + \vec{j}$): states from TSD_1 arise from the odd $\pi(i_{13/2})$ intruder coupling with a core with angular momentum sequence $R_1 = 0, 2, 4, \dots$; states from TSD_2 arise from the same odd proton but coupling with a different triaxial core with angular momentum sequence $R_2 = 1, 3, 5, \dots$. The band TSD_3 is obtained as a set of states which are built on top of TSD_2 , with the action of a wobbling frequency with wobbling phonon number $n_w = 1$; this being different than the band structure previously mentioned where the third band was a two-phonon excitation of the yrast TSD_1 . Lastly, the fourth band TSD_4 is a ground state band which results from the coupling of the same core as for TSD_2 (that is defined with the angular momentum sequence $R_2 = 1, 3, 5, \dots$) but with a different odd nucleon: $\pi(h_{9/2})$. Consequently, TSD_2 and TSD_4 are yrast states, alongside TSD_1 . Each band represents a collection of energy levels describing ground states that correspond to distinct sets of angular momenta.

For the first three bands, the MOIs are the same, and they are considered to be free parameters within the numerical calculations. However, this is not true for the fourth band, where a different set of MOIs had to be introduced, since for TSD_4 the core polarization effects are changed by coupling scheme.

Using W1, the final results pointed out to a largest MOI corresponding to the 1-axis (with \mathcal{I}_1 being the largest MOI obtained through the fitting procedure), making the system rotate around the 1-axis (that is the short axis). Moreover, the odd proton is aligned to the short axis as well, suggesting that the nucleus has a LW character. By representing the experimental wobbling energies according to Eq. 3, it was obtained that both theoretical as well as the experimental values were increasing functions with respect to an increase in angular momentum (keep in mind that the first wobbling band $n_w = 1$ within the W1 model is TSD_3). The agreement between the two sets of data (see Figure 6 from [56]) indicate that the condition for LW/TW character of the wobbling bands stated by Frauendorf et. al. in [27] is not strictly related to the increasing/decreasing wobbling energy E_{wob} . In fact, Guo et. al. [22] also point out that the approximation of Frozen Alignment (FA) from [27] neglects the Coriolis interaction of the single particle. There is an ongoing debate whether the behavior of an LW or TW triaxial nucleus is

strictly related to the change in E_{wob} with total a.m. [62–64].

A final aspect that needs to be mentioned regarding W1 has to do with the interpretation of TSD_1 and TSD_2 as being Signature Partner Bands (SPB). Signature [9] is a quantum property which appears in deformed systems. It is strictly related to the invariance of a system with quadrupole deformation (its nuclear wave-function) to a rotation by an angle π around a principal axis. For example, a rotation around the x -axis will be defined as an operator:

$$\hat{R}_x = e^{i\pi\hat{I}_x} . \quad (4)$$

As for the framework used in [56, 57], due to the wave-function describing the system being written as a product between the $|I\rangle$ basis state corresponding to the total angular momentum and the single-particle basis state $|j\rangle$, the rotation operator used in W1 achieves the following form:

$$\hat{R}_x(\pi) = e^{-i\pi\hat{I}_x} \otimes e^{-i\pi\hat{j}_x} . \quad (5)$$

If the system has an axial symmetry, only the rotation around any of the principal axes that are perpendicular to the symmetry one can define the signature quantum number. Consequently, signature is a property specific to a deformed system and it translates to a so-called *deformation invariance* with respect to the space and time reflection properties [9]. For an even-even nucleus, the signature operator \hat{R}_x has two eigenvalues, -1 and 1. For the even-odd case, the eigenvalues are $-i$ and $+i$, and depending on its total spin, the signatures can have to values, given by the following assignment:

$$\alpha_I = \frac{1}{2} (-1)^{I-1/2} . \quad (6)$$

Indeed, Eq. 6 describes the signature quantum number for a state of angular momentum I belonging to an odd mass nucleus. Such a rotational band with a sequence of states differing in spin by $\Delta I = 1$ will be divided into two branches, each branch consisting of levels differing in spin by $\Delta I = 2$, being related by the signature number $\alpha_I = \pm 1/2$. In [56] the signature concept is brought to the classical picture associated to a triaxial nucleus by means of rotation operators which act on the trial function (this function is a product of two coherent states, one that is associated to the core and one to the valence nucleon). Eqs. 27-29 from [56] will extract two signatures for TSD_1 and TSD_2 , namely the *favoured* signature $\alpha_{1f} = +1/2$ for the first band, and *un-favoured* signature $\alpha_{2u} = -1/2$ for the second band, respectively. A justification for the possibility of TSD_1 and TSD_2 of being SPB was based on the calculation of the triaxial potential (which was systematically performed in [45] and [54]), concluding that the minimum is very deep, preventing in this way the states from TSD_2 to share other minima through tunneling effects. Other experimental and theoretical results [65–68] for deformed nuclei around this mass region suggest that the calculations performed in W1 regarding the connection between TSD_1 and TSD_2 as belonging to a signature splitting phenomenon are valid and consistent with already existing interpretations.

It is instructive to mention a few key-points which arise based on the above discussion regarding W1 can be summarized as follows:

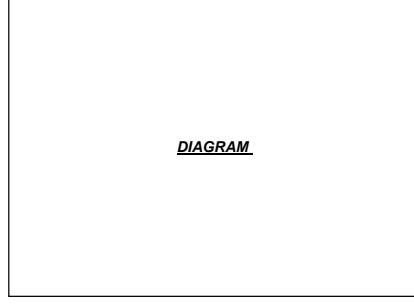


Figure 4: The general workflow involving W1 model.

- (a) The wobbling band structure in ^{163}Lu was re-interpreted: three bands are now yrast ground states, and only TSD_3 is one-phonon excited wobbling band (built on top of TSD_2)
- (b) Both TSD_1 and TSD_2 are obtained variationally, by solving the time dependent variational equation associated to the initial quantal Hamiltonian
- (c) There are three different $R + j$ coupling schemes that will produce the entire spectra:
 - (i) Coupling C_1 : The odd proton $j_1 = 13/2$ is coupled to a core sequence with a.m. $R_1 = 0, 2, 4, \dots$ (even spin states for the triaxial rotor).
 - (ii) Coupling C_2 : The same odd proton $j_1 = 13/2$ as in C_1 is coupled to a core sequence with a.m. $R_2 = 1, 3, 5, \dots$ (odd spin states for the triaxial rotor).
 - (iii) Coupling C_3 : A different odd proton $j_2 = 9/2$ is coupled to the same core as in C_2 .
- (d) Two different sets of MOIs corresponding to the triaxial nucleus (that is the rotor coupled with the odd proton) are obtained as fitting parameters throughout the numerical calculations: one for the set $TSD_{1,2,3}$ and one for TSD_4 .
- (e) TSD_1 and TSD_2 are Signature Partner Bands: with TSD_1 (TSD_2) being the favored (un-favored) partner. Their corresponding signature quantum numbers are $\alpha_{1f} = +1/2$ and $\alpha_{2u} = -1/2$.
- (f) As a side-by-side comparison with regards to the overall agreement with the experimental data, W1 yielded better results than compared to the previous work depicted in Ref. [53], although it must be mentioned that both models are based on semi-classical approaches.

A diagram which shows the workflow involved in W1 can be seen in the Figure 4. Also, a comparison with previous calculations can be seen in Figure 21 from Ref. [56].

2.2 W2 - Signature Partner Bands + Parity Partner Bands

The main question which can be asked regarding the formalism W1 that was described in 2.1 is whether it is possible to obtain a *unified* description for all four bands in ^{163}Lu in relation to the coupling scheme. In other words, it is worth investigating the possibility of having a unique single-particle state j that is coupled to a core of positive parity for the bands $TSD_{1,2,3}$ and a core of negative parity for TSD_4 .

Fortunately, the answer is positive: starting from the semi-classical formalism of W1, one can properly adjust the coupling scheme, making sure that the entire numerical recipe used for obtaining the energy spectrum of ^{163}Lu remains consistent with the experimental results.

Regarding the unique single-particle that couples to the triaxial core, it is natural to pick the $i_{13/2}$ proton (that is j_1 from W1). Reasoning behind this choice has to do with the microscopic calculations [10, 13, 61] that showed stable triaxial structures in the ^{163}Lu potential energy surface when the triaxial core couples with a highly aligned j -shell particle, strongly indicating the $\pi(i_{13/2})$ proton. Keep in mind that a highly aligned j -nucleon will *prefer* to keep a certain triaxial deformation when coupled to a core [69–71] (in the sense that the triaxiality parameter γ will have a certain value based on the orbital of the odd nucleon), and using microscopic calculations following the Ultimate Cranker code, it has been shown that a value of $\gamma \approx +20^\circ$ is preferred by the odd $\pi(i_{13/2})$ nucleon.

By taking j_1 as the sole intruder that couples to a positive core and also a negative core, the sequences with even/odd integer spins for the core do not change. In fact, the coupling schemes can be readily obtained, making sure that the final (experimental) spin sequences for each TSD band is intact.

- (a) Coupling C'_1 : the odd j_1 proton aligns with the core of even-integer spin sequence $R_1 = 0, 2, 4, \dots$, with a parity of the R_1 core that is positive $\pi(R_1) = +1$.
- (b) Coupling C'_2 : the odd j_1 proton aligns with the core of even-integer spin sequence $R_2^+ = 1, 3, 5, \dots$, with a parity of the R_2^+ core that is positive $\pi(R_2^+) = +1$.
- (c) Coupling C'_3 : the odd j_1 proton aligns with the core with an odd-integer spin sequence $R_2^- = 1, 3, 5, \dots$, which has negative parity $\pi(R_2^-) = -1$.

From the three schemes defined above, it is clear that C'_1 corresponds to the yrast TSD_1 , C'_2 to the ground state TSD_2 , and finally C'_3 to the ground state TSD_4 . Obviously, the odd valence nucleon j_1 has a positive parity $\pi_{j_1} = +1$. There has not been attributed a coupling scheme for TSD_3 , since this band still remains as the one-wobbling phonon excitation that is built on top of TSD_2 with the action of a phonon operator which will be characterized later on. The three couplings are schematically represented in Figure 5. One should keep in mind the fact that \vec{j}_1 is aligned with the axis with largest MOI does not necessarily mean the fact that the model works within the Frozen Alignment approximation - it is just an illustration.

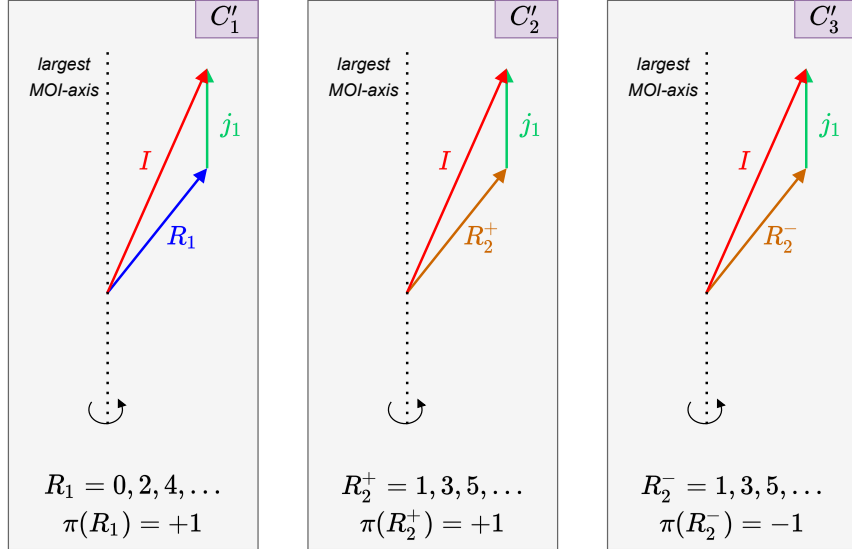


Figure 5: A schematic representation with the three coupling schemes that characterize the W2 model. The same odd particle (the $j_1 = i_{13/2}$ proton) is coupled with two positive cores with even/odd integer spin sequences for TSD_1/TSD_2 , and one negative core in the case of TSD_4 with odd integer spin sequence. The total spin of the system precesses around the axis with the largest MOI, as it is the case for a triaxial rotor.

It is expected that the Hamiltonian of the system will keep a similar form, since there are no new interactions or modified particle-core states added in the problem. The argument holds because the coupling scheme which involves the fourth triaxial band TSD_4 only the single-particle's angular momentum state changes, but the core has negative parity. As a result, the treatment of the problem will follow the same manner as it did in the W1 case.

The last step in searching for a unified coupling scheme in ^{163}Lu is to establish a possible relationships between the four bands. As per the calculations involved in W1, it was proven that signature quantum number is a good quantum number and indeed, a sign that TSD_1 and TSD_2 are signature partners emerged. Their overall similar properties and spin difference enforce this argument. Furthermore, in this new W2 approach, the difference in parity between the TSD_2 and TSD_4 but the same angular momentum sequence of their corresponding triaxial core R_2^+ and R_2^- strongly suggest that the two bands are *Parity Partner Bands*: two rotational sequences with energy states characterized by opposite parity, increasing energy that follows a trend $\propto I(I+1)$, and a spin difference $\Delta I = 2$ between states belonging to the same band. In the following section, calculations which will show that parity is indeed a good quantum number for the triaxial rotor + odd-particle system will be provided. For what it is worth mentioning now is that the concept of parity partners between TSD_2 and TSD_4 emerge from the idea that a stable strongly deformed structure is achieved from a single quasiparticle which moves in a quadrupole mean-field generated by a triaxial even-even core. However, there is a

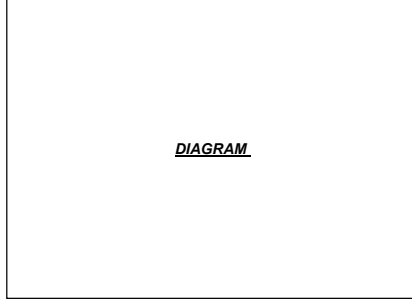


Figure 6: The general workflow involving W2 model.

splitting in two different cases of coupling mechanisms, namely C'_2/C'_3 depending on the alignment of the high- j -shell particle with a core of positive/negative parity.

Similar structures with alternating positive-negative parity bands have been also reported in other nuclei such as ^{40}Ca [72], or some heavier isotopes like ^{218}Fr [73]. In fact, one of the authors of this present work developed a unified description of states with positive and negative parity in odd-mass nuclei [74, 75], although therein, a quadrupole-octupole term was introduced within the particle-core Hamiltonian to describe this feature. Concluding the current subsection, a diagram which shows the workflow involved in W2 can be seen in the Figure 6.

3 Theoretical Background

In this section, a description of the framework used for obtaining the wobbling spectrum of ^{163}Lu is made. As stated in the previous section, the system is described with a similar Hamiltonian used in W1, namely the Hamiltonian for a triaxial PRM model.

$$H = H_{\text{core}} + H_{\text{s.p.}} \quad . \quad (7)$$

The Hamiltonian in Eq. 7 describes a system in which an odd j quasiparticle interacts with a triaxial even-even core, where the odd nucleon is moving in a quadrupole deformed mean field that is generated by the core itself. As such, the first H_{core} term in the Hamiltonian corresponds describes the motion triaxial core, while the second term $H_{\text{s.p.}}$ represents the aligned valence proton (j_1 in this case).

Indeed, the core Hamiltonian is given by:

$$H_{\text{core}} = \sum_{i=1,2,3} \frac{1}{2\mathcal{I}_i} (I_i - j_i)^2 \quad , \quad (8)$$

where the core angular momentum is $\vec{R} = \vec{I} - \vec{j}$ and the terms \mathcal{I}_i represent the moments of inertia for triaxial ellipsoid along the principal axes. These three moments of inertia will be considered as free parameters in the present calculations, but, compared to the

work W1, a unique set of MOIs will be attributed to the four bands, since the triaxial core will create an alignment with a unique single particle, that is j_1 . Because of this, there is not option for their nature (i.e., rigid or hydrodynamic).

The single-particle Hamiltonian from Eq. 7 is derived from the well-known Nilsson potential [76, 77]:

$$h(\beta_2, \gamma) = C \left\{ \cos \gamma Y_{20}(\theta, \varphi) + \frac{\sin \gamma}{\sqrt{2}} [Y_{22}(\theta, \varphi) + Y_{2-2}(\theta, \varphi)] \right\} , \quad (9)$$

where the coupling parameter C causes the level splitting in the deformed field and it is directly proportional to the quadrupole deformation β_2 . The potential h from Eq. 9 is written in terms of the quadrupole deformation and triaxiality parameter that play the role of deformation parameters within a triaxial system (β_2, γ) . Its expression using the coupling parameter C is widely used when working with a particle-rotor-model [78–80]. In our case, by applying a Wigner-Eckart theorem for the single- j particle, the following expression for $H_{s.p.}$ will be obtained:

$$H_{s.p.} = \frac{V}{j(j+1)} \left[\cos \gamma (3j_3^2 - \vec{j}^2) - \sqrt{3} \sin \gamma (j_1^2 - j_2^2) \right] + \epsilon_j . \quad (10)$$

This term describes the motion of an odd particle with angular momentum j in a mean field generated by a triaxial core, with a potential strength V characterized by the quadrupole deformation ($V \propto \beta_2$). In fact, the single-particle potential strength V will be considered as fourth free parameter within the calculations and its behavior will dictate the coupling of the j_1 particle with all four TSD bands. The term ϵ_j from Eq. 10 represents the single particle energy that corresponds to the odd j_1 proton from the i -orbital.

Regarding the triaxial deformation γ which enters in Eq. 10, its value will be considered as another free parameter of the current problem. In other words, having V and γ as free parameters means that the system will be described by its deformation parameters which will be obtained through a fitting procedure, keeping an agreement with the experimental data regarding the excitation energies of the rotational states belonging to $TSD_{1,2,3,4}$.

From Eqs. 8 and 10, the free parameter set can be obtained, hereafter denoted by \mathcal{P} . It comprises three moments of inertia, the single particle potential strength, and the triaxial deformation. As such, \mathcal{P} can be written as:

$$\mathcal{P} = [\mathcal{I}_1, \mathcal{I}_2, \mathcal{I}_3, V, \gamma] . \quad (11)$$

Solving the problem of W2 is equivalent to finding the eigenvalues of H given in Eq. 7. In a similar approach as in W1, the eigenvalues of interest are obtained on the base of a semi-classical approach. Thus, the first step is to perform a de-quantization procedure on H through a TDVE [45, 53, 55]:

$$\delta \int_0^t \langle \Psi_{IjM} | H - i \frac{\partial}{\partial t'} | \Psi_{IjM} \rangle dt' = 0 . \quad (12)$$

Working within a semi-classical approach allows one to keep a close contact with the system's dynamics in terms of *friendly* and *familiar* equations of motion and generalized coordinates. The trial function from Eq. 12 is carefully chosen as a product of two basis states comprising the states with total angular momentum I and j , respectively:

$$|\Psi_{IjM}\rangle = \mathbf{N} e^{z\hat{I}_-} e^{s\hat{j}_-} |IMI\rangle |jj\rangle , \quad (13)$$

where the operators \hat{I}_- and \hat{j}_- denote the lowering operators for the intrinsic angular momenta \vec{I} and \vec{j} , respectively, and \mathbf{N} plays the role of the normalization constant. One must remark the fact that the states from Eq. 13 are extremal states for the operators (\hat{I}^2, \hat{I}_3) and (\hat{j}^2, \hat{j}_3) , respectively, and they correspond to the maximally allowed states for a given set of angular momenta I and j . As an observation, the trial function is an admixture of components of definite K , which is consistent with the fact that for a triaxial nucleus, K is not a good quantum number.

The variables z and s from Eq. 13 are complex functions that depend on time, and they play the role of classical coordinates of the phase spaces that describe the motion of the core and the odd particle:

$$z = \rho e^{i\varphi} , \quad s = f e^{i\psi} . \quad (14)$$

In order to obtain a set of classical equations which describe the motion of the system, a new pair of variables are introduced from the ones defined in Eq. 14. Indeed, defining the variables (φ, r) and (ψ, t) with r and t given by the relations:

$$r = \frac{2I}{1 + \rho^2} , \quad t = \frac{2j}{1 + f^2} , \quad (15)$$

where $r \in [0, 2I]$ and $t \in [0, 2j]$, will bring the classical equations of motion (provided by the TDVE), to a straightforward canonical form:

$$\begin{aligned} \frac{\partial \mathcal{H}}{\partial r} &= \dot{\varphi} ; \quad \frac{\partial \mathcal{H}}{\partial \varphi} = -\dot{r} , \\ \frac{\partial \mathcal{H}}{\partial t} &= \dot{\psi} ; \quad \frac{\partial \mathcal{H}}{\partial \psi} = -\dot{t} . \end{aligned} \quad (16)$$

The function \mathcal{H} denotes the average of the Hamiltonian operator H (Eq. 7) with the trial function $|\Psi_{IjM}\rangle$ given in Eq. 13, and it plays the role of classical energy. It represents a crucial quantity within the present analysis, as it will provide an insight with regards to the stability of the wobbling motion in ^{163}Lu , and can help to pinpoint several regions where the wobbling mode is stable/unstable. This analysis will be performed in the following sections.

$$\mathcal{H} \equiv \mathcal{H}(\bar{q}_{\text{core}}; \bar{q}_{\text{s.p.}}) = \langle \Psi_{IjM} | H | \Psi_{IjM} \rangle , \quad (17)$$

where the set of coordinates (φ, r) and (ψ, t) were conveniently denoted with \bar{q}_{core} and $\bar{q}_{\text{s.p.}}$, respectively, since they describe the motion of the core and that of the odd single-particle.

Starting from the equations of motion given in Eq. 16, one can observe that the function \mathcal{H} is a constant of motion, that is $\dot{\mathcal{H}} \equiv 0$. This equation will define a surface, a so-called equi-energy surface $\mathcal{H} = \text{const.}$ It is worth mentioning the fact that such an equality holds since the entire set of equations of motion emerged from a variational principle. The sign of the Hessian associated to this classical function will indicate its stationary points. Among them, some are minima. One such example of minimum points for \mathcal{H} is the point $p_0 \equiv (q_{\text{core}}; q_{\text{s.p.}})|_{\text{min}} = (\varphi, r; \psi, t)|_{\text{min}} = (0, I; 0, j)$. The energy function is minimum $\mathcal{H}_{\text{min}}^{(I,j)} = H(p_0)$ as long as the condition on the MOIs is: $\mathcal{I}_1 > \mathcal{I}_2 > \mathcal{I}_3$. There is no restriction on γ .

With a linearization procedure for the equations of motion around the minimum point p_0 of \mathcal{H} , a dispersion equation will be obtained:

$$\Omega^4 + B\Omega^2 + C = 0 . \quad (18)$$

The above equation describes a harmonic type of motion for the nuclear system, with the solutions to this algebraic equation as the *wobbling frequencies* Ω . The terms B and C are complex functions of total angular momentum I , single particle a.m. j , inertial parameters $A_k = 1/(2\mathcal{I}_k)$, $k = 1, 2, 3$, single particle potential strength V , and triaxiality parameter γ . In fact, one can conclude that B and C are functions which depend on the parameter set defined in Eq. 11: $B = f_B(\mathcal{P}, I, j)$ and $C = f_C(\mathcal{P}, I, j)$, respectively. The B term from Eq. 18 has the expression [56]:

$$-B = [(2I - 1)(A_3 - A_1) + 2jA_1] [(2I - 1)(A_2 - A_1) + 2jA_1] + 8A_2A_3Ij + (\mathbf{T}_B^1 \times \mathbf{T}_B^2) , \quad (19)$$

where the terms \mathbf{T}_B^1 and \mathbf{T}_B^2 are defined as:

$$\begin{aligned} \mathbf{T}_B^1 &= \left[(2j - 1)(A_3 - A_1) + 2IA_1 + V \frac{2j - 1}{j(j + 1)} \sqrt{3}(\sqrt{3} \cos \gamma + \sin \gamma) \right] , \\ \mathbf{T}_B^2 &= \left[(2j - 1)(A_2 - A_1) + 2IA_1 + V \frac{2j - 1}{j(j + 1)} 2\sqrt{3} \sin \gamma \right] . \end{aligned} \quad (20)$$

Accordingly, the C term from Eq. 18 has the expression [56]:

$$\begin{aligned} C &= \{[(2I - 1)(A_3 - A_1) + 2jA_1] \times \mathbf{T}_C^1 - 4IjA_3^2\} \times \\ &\quad \times \{[(2I - 1)(A_2 - A_1) + 2jA_1] \times \mathbf{T}_C^2 - 4IjA_2^2\} , \end{aligned} \quad (21)$$

where the terms \mathbf{T}_C^1 and \mathbf{T}_C^2 are defined as:

$$\begin{aligned} \mathbf{T}_C^1 &= \left[(2j - 1)(A_3 - A_1) + 2IA_1 + V \frac{2j - 1}{j(j + 1)} \sqrt{3}(\sqrt{3} \cos \gamma + \sin \gamma) \right] , \\ \mathbf{T}_C^2 &= \left[(2j - 1)(A_2 - A_1) + 2IA_1 + V \frac{2j - 1}{j(j + 1)} 2\sqrt{3} \sin \gamma \right] . \end{aligned} \quad (22)$$

It can be seen that the terms which enter in B and C , namely $(\mathbf{T}_B^1, \mathbf{T}_B^2)$ from Eq. 20 and $(\mathbf{T}_C^1, \mathbf{T}_C^2)$ from Eq. 22 that enter in B and C correspond to the quadrupole

deformation that causes the single particle to move in the mean-field of the triaxial core. The terms also define the triaxiality that the nucleus achieves once the odd proton couples to the triaxial core, driving the system up to a large (and stable) deformation.

Going back to Eq. 18, under the restrictions for the MOIs defined above, the dispersion equation admits two real and positive solutions (hereafter denoted with Ω_1^I and Ω_2^I , where $\Omega_1^I < \Omega_2^I$) defined for $j_1 = i_{13/2}$, given by:

$$\Omega_{1,2}^I = \sqrt{\frac{1}{2} (-B \mp (B^2 - 4C)^{1/2})} . \quad (23)$$

These two solutions are interpreted as *wobbling frequencies* associated to the motion of the core, and the motion of the odd-particle. As such, each wobbling frequency has an associated wobbling-phonon number:

$$\Omega_1^I \rightarrow n_{w_1} ; \Omega_2^I \rightarrow n_{w_2} . \quad (24)$$

It is worth noting that they take part in the energy states belonging to all the bands, including the yrast ones, however, in that case, the zero-point motion is considered (meaning that the two wobbling quanta (Ω_1^I, Ω_2^I) are halved).

Now that the energy function \mathcal{H} was defined in terms of classical coordinates \bar{q}_{core} and $\bar{q}_{\text{s.p.}}$, and the wobbling frequencies were obtained, it is possible to define the analytical expressions for the four TSD bands of ^{163}Lu . Indeed, taking care of the coupling scheme for each band, they can be formulated as follows:

$$\begin{aligned} E_{\text{TSD1}}^I &= \epsilon_j + \mathcal{H}_{\text{min}}^{(I,j)} + \mathcal{F}_{00}^I , \quad I = 13/2, 17/2, 21/2 \dots \\ E_{\text{TSD2}}^I &= \epsilon_j^1 + \mathcal{H}_{\text{min}}^{(I,j)} + \mathcal{F}_{00}^I , \quad I = 27/2, 31/2, 35/2 \dots \\ E_{\text{TSD3}}^I &= \epsilon_j + \mathcal{H}_{\text{min}}^{(I-1,j)} + \mathcal{F}_{10}^I , \quad I = 33/2, 37/2, 41/2 \dots \\ E_{\text{TSD4}}^I &= \epsilon_j^2 + \mathcal{H}_{\text{min}}^{(I,j)} + \mathcal{F}_{00}^I , \quad I = 47/2, 51/2, 55/2 \dots , \end{aligned} \quad (25)$$

where $\mathcal{F}_{n_{w_1} n_{w_2}}^I$ is a function of the wobbling frequencies, defined accordingly to the wobbling phonon numbers n_{w_1}, n_{w_2} :

$$\mathcal{F}_{n_{w_1} n_{w_2}}^I = \Omega_1^I \left(n_{w_1} + \frac{1}{2} \right) + \Omega_2^I \left(n_{w_2} + \frac{1}{2} \right) , \quad (26)$$

and $\mathcal{H}_{\text{min}}^{(I,j)}$ is the classical energy evaluated in its minimal point p_0 .

A few aspects regarding the energy spectrum defined in Eq. 25 are worth mentioning. To each band there is a specific energy ϵ_j associated to the single-particle state from the orbital that it belongs to. In this case, the odd-proton j_1 with $j = 13/2$ from the i -orbital is the one that couples to the triaxial core. However, for the bands TSD_2 and TSD_4 , a different re-normalization is considered, since the former is the unfavored partner within the structure, and the latter is the negative parity partner within the band structure. These quantities will shift the overall energy states belonging to the two bands, each by a different amount. As a result, both ϵ_j^1 and ϵ_j^2 will be adjusted throughout the

Band	n_{w_1}	n_{w_2}	π	α	Coupling scheme
TSD_1	0	0	+1	+1/2	C'_1
TSD_2	0	0	+1	-1/2	C'_2
TSD_3	1	0	+1	+1/2	Built on top of TSD_2
TSD_4	0	0	-1	-1/2	C'_3

Table 1: The wobbling phonon numbers, parities, signatures, and coupling schemes assigned to each triaxial band in ^{163}Lu , within the W2 model. The three coupling schemes were defined in Section 2.2.

the numerical calculations such that the energy spectrum is best reproduced. Another aspect concerns the band TSD_3 ; since this is the only excited wobbling band within the family, its configuration is built on top of TSD_2 , with the action of a single phonon ($n_{w_1} = 0$) operator, that is the wobbling frequency Ω_1^I . Consequently, an energy state I belonging to TSD_3 is obtained from a state $I - 1$ from TSD_2 . In Table 1, the rest of the wobbling phonon numbers are mentioned, with the parity, signature, and coupling scheme for each band in particular.

3.1 Parity quantum number for the wave-function

In W1 it was shown that signature emerges from the calculations on the total wave-function as a good quantum number for this triaxial system. This is why in [56] the bands TSD_1 and TSD_2 appeared as Signature Partner Bands (SPB). In W2, this property still stands.

Since the backbone of the current work started from the need for a single odd-particle that couples to a triaxial core in ^{163}Lu , one has to look at the band TSD_4 (which was interpreted as having a different nucleon: j_2 with $j = 9/2$ from the h -orbital), and see if its differentiating properties can be linked to *main group* of bands (namely $TSD_{1,2,3}$). Indeed, from the experimental measurements regarding spin and parity assignment [61], it turns out that the parity of the rotational states are negative. Therefore, a forensic analysis on this quantum property should be considered as the necessary ingredient in a unified description of all four bands.

Choosing the rotation axis as the 2-axis, one can define the parity operator as a product of the complex conjugation operation and a rotation of angle π around this axis: $P = e^{-i\pi\hat{I}_2}C$. Obviously, \hat{I}_2 represents the rotation operator and C the operator performs a complex conjugation when applied on a wave-function. The total parity operator is defined as the product of an operator corresponding to the core and one corresponding to the single particle:

$$\mathcal{P}_T = P_{\text{core}}P_{\text{s.p.}} . \quad (27)$$

Acting with the total parity operator defined above, on the trial function $\Psi(\bar{q}_{\text{core}}; \bar{q}_{\text{s.p.}})$ associated to the triaxial system (which was defined in Eq. 13), the following is obtained:

$$\mathcal{P}_T\Psi(r, \varphi; t, \psi) = \Psi(r, \varphi + \pi; t, \psi + \pi) \stackrel{\text{not.}}{=} \bar{\Psi}. \quad (28)$$

Keep in mind that the set of classical phase space which describe the individual dynamics of the core and the single-particle, namely \bar{q}_{core} and $\bar{q}_{\text{s,p}}$, were defined as (r, φ) and (t, ψ) .

The classical energy function \mathcal{H} has an invariance property at changing the angles with π :

$$\mathcal{H}(r, \varphi; t, \psi) = \mathcal{H}(r, \varphi + \pi; t, \psi + \pi) . \quad (29)$$

From Eqs. 28 and 29, it can be concluded that the wave-function describing the triaxial system Ψ and its image through \mathcal{P}_T , $\bar{\Psi}$, are two linearly dependent functions which differ only by a multiplicative constant μ , with $|\mu| = 1$. Thus, μ can either be -1 or +1, such that:

$$\Psi(r, \varphi + \pi; t, \psi + \pi) = \pm \Psi(r, \varphi; t, \psi) . \quad (30)$$

The above result concludes the parity analysis for the wave-function, showing that the triaxial rotor admits eigenfunctions (i.e., deformed rotational states) of negative parity. Therefore, a single wave-function characterized by the coupling of a triaxial core to the odd proton $i_{13/2}$ is describing both positive parity states ($\in TSD_{1,2,3}$) as well as negative parity states ($\in TSD_4$). This analysis, together with the fact that TSD_2 and TSD_4 have the same a.m. sequences (although TSD_2 has more states with low spin than TSD_4) suggest the fact that these two bands might be Parity Partners.

3.2 Energy function - geometrical interpretation

The analytical expression for the average of H with the trial function describing the system was previously calculated in W1. Indeed, the energy function \mathcal{H} was given in terms of the phase space coordinates as follows [56]:

$$\begin{aligned} \mathcal{H} = & \frac{I}{2}(A_1 + A_2) + A_3 I^2 + \frac{2I-1}{2I} r(2I-r) \mathcal{A}_\varphi + \frac{j}{2}(A_1 + A_2) + A_3 j^2 + \frac{2j-1}{2j} t(2j-t) \mathcal{A}_\psi \\ & - 2\sqrt{r(2I-r)t(2j-t)} \mathcal{A}_{\varphi\psi} + A_3 [r(2j-t) + t(2I-r)] - 2A_3 Ij + V \frac{2j-1}{j+1} \mathcal{A}_\gamma , \end{aligned} \quad (31)$$

with:

$$\begin{aligned} \mathcal{A}_\varphi &= (A_1 \cos^2 \varphi + A_2 \sin^2 \varphi - A_3) , \\ \mathcal{A}_{\varphi\psi} &= (A_1 \cos \varphi \cos \psi + A_2 \sin \varphi \sin \psi) , \\ \mathcal{A}_\psi &= (A_1 \cos^2 \psi + A_2 \sin^2 \psi - A_3) , \\ \mathcal{A}_\gamma &= \left[\cos \gamma - \frac{t(2j-t)}{2j^2} \sqrt{3} (\sqrt{3} \cos \gamma + \sin \gamma \cos 2\psi) \right] \end{aligned} \quad (32)$$

It is instructive to check the dependence of the energy function on the angular momentum components, e.g. the coordinates $x_k \stackrel{\text{not.}}{=} I_k$, $k = 1, 2, 3$, where the quantization

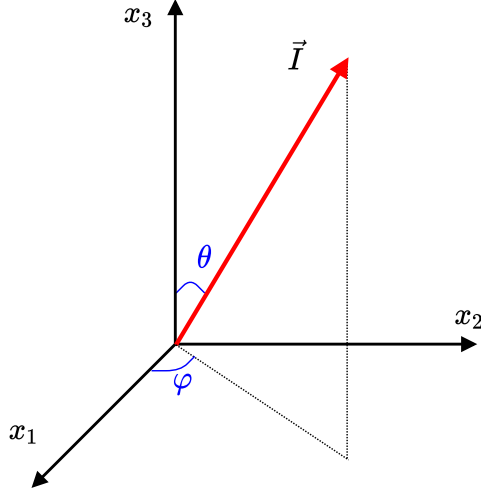


Figure 7: The angular momentum vector of the triaxial nucleus, represented in the angular momentum space. The polar angles (θ, φ) that the vector makes with the x_3 axis and x_1x_2 -plane, respectively, are also shown.

axis is chosen as the 3-axis. By expressing the angular momentum coordinates $x_{1,2,3}$ in terms of the polar angles (θ, φ) and a radius r given by the value of the total angular momentum I , one obtains:

$$x_1 = I \sin \theta \cos \varphi, \quad x_2 = I \sin \theta \sin \varphi, \quad x_3 = I \cos \theta. \quad (33)$$

The total angular momentum \vec{I} of the nucleus can be visualized as a vector within the angular momentum space, generated by its three components. Figure 7 shows such a representation.

Within this spherical coordinates, and evaluating the energy function around its minimum point p_0 , the following expression for \mathcal{H} arises:

$$\mathcal{H} |_{p_0} = I \left(I - \frac{1}{2} \right) \sin^2 \theta (A_1 \cos^2 \varphi + A_2 \sin^2 \varphi - A_3) - 2A_1 I j \sin \theta + T_{\text{core}} + T_{\text{s.p.}}. \quad (34)$$

The last two terms in this equation are independent on the polar angles (θ, φ) , and they characterize the minimal energy for a triaxial rotor with given spin I and inertial parameters $A_{1,2,3}$ (i.e., T_{core}), and the minimal energy which arises from the coupling j nucleon that is moving in the mean-field of strength V and triaxial deformation γ (i.e., $T_{\text{s.p.}}$). The two terms have the following form:

$$\begin{aligned} T_{\text{core}} &= \frac{I}{2} (A_1 + A_2) + A_3 I^2, \\ T_{\text{s.p.}} &= \frac{j}{2} (A_2 + A_3) + A_1 j^2 - V \frac{2j-1}{j+1} \sin \left(\gamma + \frac{\pi}{6} \right). \end{aligned} \quad (35)$$

The classical equation of motion admits two constants of motion: the total system energy (E) and the total angular momentum (I). In fact, this is consistent with the quantum theory, where the angular momentum vector of a rotating system is conserved. A 3-dimensional rotation can be visualized in a semi-classical description as a precession of the total angular momentum vector around the rotational axis (that is the axis with largest MOI). Consequently, by finding the intersection line(s) between the surface of the energy ellipsoid E and the surface of the angular momentum sphere generated by the spin I , then one can assert that a trajectory which belongs to a defined rotational state (with given spin and energy) is found. It is possible to construct a set with all the allowed trajectories for ^{163}Lu if the energy ellipsoid E for a given state $I \in TSD_k$, $k = 1, 2, 3, 4$ is evaluated, then intersected with its corresponding angular momentum I . Such representations will be made in the following section.

The expression of the energy ellipsoid within Cartesian coordinates can be obtained by expressing the classical energy using the coordinate representation which was fixed in Eq. 33. As a result, its expression becomes:

$$E = \left(1 - \frac{1}{2I}\right) A_1 x_1^2 + \left(1 - \frac{1}{2I}\right) A_2 x_2^2 + \left[\left(1 - \frac{1}{2I}\right) A_3 + A_1 \frac{j}{I}\right] x_3^2 - I \left(I - \frac{1}{2}\right) A_3 - 2A_1 I j + T_{\text{rot}} + T_{\text{sp}} . \quad (36)$$

The above expression describes a deformed ellipsoid characterized by a total angular momentum I and single particle a.m. j . For a total angular momentum \vec{I} , the vector generates a sphere of radius $r = I$ within the space generated by the angular momentum components denoted by (x_1, x_2, x_3) . The sphere's equation is given by:

$$I^2 = x_1^2 + x_2^2 + x_3^2 . \quad (37)$$

As a quick side-note, the radius of the angular momentum sphere generated by a vector \vec{I} is in fact $r \approx \sqrt{I(I+1)}$, since $\hat{I}^2 = I(I+1)$, however, in the high spin limit $I \gg 1$, an approximation $r \approx I$ holds without issues. The trajectories obtained through the intersection of Eqs. 36 and 37 will aim to show a classical visualization of the wobbling character for a triaxial nucleus.

4 Numerical results

Now that the theoretical formalism was established, with the W2 approach described in Section 2.2, it is natural to proceed to the numerical analysis. As a first step, this will consist in the calculation of the energies from the four TSD bands in ^{163}Lu . If the formulated theory regarding the wave-function of the triaxial system is indeed consistent, then an agreement with the experimental data should be achieved. In fact, this will be best verified after the negative parity states from TSD_4 will be obtained.

Regarding the wobbling spectrum of ^{163}Lu , its theoretical interpretation was given in Eq. 25. As mentioned, those energy formulas are parametrized in terms of \mathcal{P} , which

Band	n_s	\vec{j}	\vec{R} - Sequence	\vec{I} - Sequence	Coupling scheme
TSD_1	21	j_1	$R_1 = 0, 2, 4, \dots$	$13/2, 17/2, 21/2, \dots$	C'_1
TSD_2	17	j_1	$R_2^+ = 1^+, 3^+, 5^+, \dots$	$27/2, 31/2, 35/2, \dots$	C'_2
TSD_3	14	j_1	1-phonon excitation	$33/2, 37/2, 41/2, \dots$	1-phonon excitation
TSD_4	11	j_1	$R_2^- = 1^-, 3^-, 5^-, \dots$	$47/2, 51/2, 55/2, \dots$	C'_3

Table 2: The number of energy states n_s within each wobbling band, the coupling proton a.m. \vec{j} , the core a.m. \vec{R} , the nucleus's a.m. \vec{I} , and the corresponding coupling scheme that was established according to the W2 model.

is the set of free parameters established in Eq. 11. Determining \mathcal{P} , results in obtaining the theoretical energies. Consequently, a fitting procedure can be applied within the present formalism, finding an arbitrary parameter set \mathcal{P} such that the deviations from the experimental results are minimal. Working with a χ^2 fit test:

$$\chi^2 = \frac{1}{N_T} \sum_i^{N_T} \frac{(E_{\text{exp}} - E_{\text{th}})^2}{E_{\text{exp}}}, \quad (38)$$

where N_T represents the total number of states (comprising each rotation state I from all four bands $TSD_{1,2,3,4}$) it was possible to develop a numerical searching procedure for *the best* parameter set. It should be retained that each energy state E_{th} has its own given formula (function of spin I and single particle a.m. j), depending on the band where that state is located. Table 2 contains the number of states within each band, with the corresponding spin sequences for the core a.m. R and total a.m. I , in accordance with the coupling schemes specific to W2 formalism.

Performing the minimization procedure on all four bands, the parameter set was determined, its values being illustrated in Table 3. This W2 method contrasts the approach in W1, where a second minimization process was needed separately for TSD_4 . The root mean square error provided by the obtained parameter set \mathcal{P} has a value of $E_{\text{rms}} \approx 79 \text{ keV}$. This result is much better than the one obtained with previous formalism W1 (where an E_{rms} of approximately 240 keV was obtained [56]). As a matter of fact, this is the first semi-classical formalism in the literature which achieves agreement with the experimental data with less than 100 keV for the entire wobbling spectrum of ^{163}Lu . It is worth mentioning that the fitting procedure was done not for the absolute wobbling energies E_{TSDk}^I , $k = 1, 2, 3, 4$, but for the *excitation energies* which are relative to the hand-head $I = 13/2^+$ from the first yrast state TSD_1 . Comparison between the theoretical values obtained within the current formalism and the experimental data is shown in Figures 8 and 9. For the sake of completeness, the wobbling frequencies which enter in the expression of the $\mathcal{F}_{n_{w_1}n_{w_2}}^I$ given by Eq. 26 are graphically represented as function of total angular momentum I for the same parameter set in Figure 10. It is remarkable the fact that the wobbling frequency Ω_2^I is much larger than its partner, suggesting the fact the coupling effects caused by the highly aligned proton have a stronger influence in achieving a wobbling character for ^{163}Lu , which is in line with the characteristics of

\mathcal{I}_1 [\hbar^2/MeV]	\mathcal{I}_2 [\hbar^2/MeV]	\mathcal{I}_3 [\hbar^2/MeV]	γ [deg.]	V [MeV]
72	15	7	22	2.1

Table 3: The parameter set \mathcal{P} that was determined by a fitting procedure of the excitation energies for ^{163}Lu .

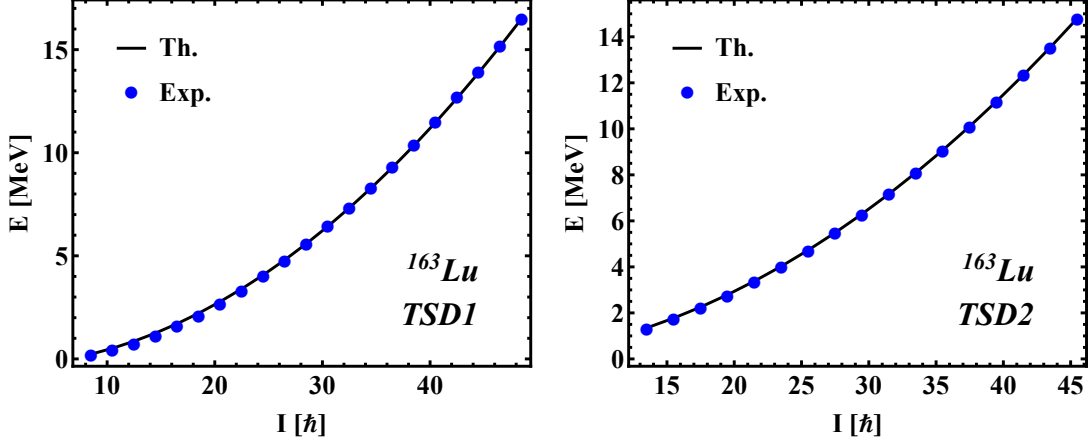


Figure 8: Comparison between theoretical and experimental excitation energies for the first two wobbling bands in ^{163}Lu . The theoretical results are obtained with the parameters listed in Table 3. Experimental data is taken from [81].

a particle-rotor coupling. Another feature of these wobbling frequencies is their linear behavior with respect to the nuclear spin.

Concerning the single-particle energies from Eq. 25, namely ϵ_j^1 and ϵ_j^2 that emerge from the un-favored signature of TSD_2 and negative parity of TSD_4 , respectively, they induce a correction for the mean-field with the quantities $\epsilon_j^1 - \epsilon_j = 0.3 \text{ MeV}$ and $\epsilon_j^1 - \epsilon_j = 0.6 \text{ MeV}$. Note that since the energy state $I_{13/2} \in TSD_1$ (the band-head of TSD_1) was subtracted from all bands, the single particle energies for band 2 and 4 are adjusted accordingly. One remarks that in the case of ϵ_j^1 , which is related to the shift due to the un-favored partner of the yrast TSD_1 band, the obtained numerical value is three times smaller than the signature splitting predicted by Jensen et. al. in [13]. Moreover, a lower single-particle shift allows for TSD_2 states to lie close to the yrast region from TSD_1 , making the current interpretation for the band structure valid, since a larger shift might correlate with cranking effects and the current Hamiltonian does not take into account such constraints.

Another noteworthy aspect of this current formalism is the fact that the theoretical difference $\delta_{42} = E_{TSD_4}^I - E_{TSD_2}^I \approx 0.3 \text{ MeV}$. This suggests that the states of the same a.m. from TSD_2 and TSD_4 bands might emerge through the parity projection of a sole wave-function which does not have reflection symmetry. In the present case, this is caused by the fact that the wobbling frequency is parity independent. Consequently, these two bands indeed behave as a pair of parity partners, as defined in [82–84].

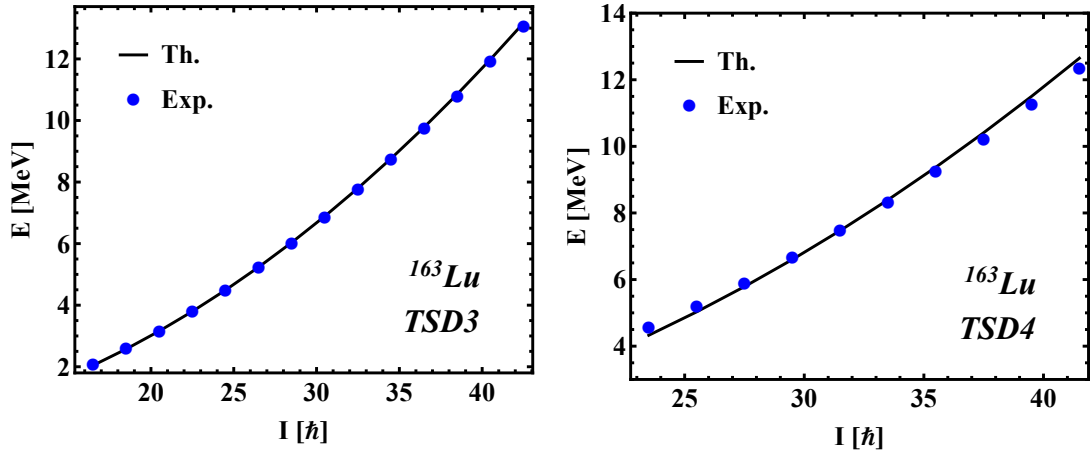


Figure 9: Comparison between theoretical and experimental excitation energies for third and fourth wobbling bands in ^{163}Lu . The theoretical results are obtained with the parameters listed in Table 3. Experimental data is taken from [81].

4.1 Interpretation of the parameter set \mathcal{P}

Performing the fitting procedure for the excitation energies of ^{163}Lu results in the moments of inertia \mathcal{I}_k given in Table 3, together with the single-particle potential strength V , and triaxiality parameter γ . Interpretation of their numerical values is mandatory in order to check whether the current formalism is valid or not.

Regarding the moments of inertia, it is clear that the axis of rotation for the energy ellipsoid is the 1-axis, as the largest MOI is \mathcal{I}_1 , causing a maximal density distribution across this axis [27]. The MOI ordering is $\mathcal{I}_1 \gg (\mathcal{I}_2 > \mathcal{I}_3)$, and compared with the results of the previous work W1, the current 1-axis MOI is bigger than both $\mathcal{I}_1^{\text{TSD1,2,3}} = 63.2$ and $\mathcal{I}_1^{\text{TSD4}} = 67$ (data taken from Table 1 in Ref. [57]). This is expected, since here, TSD_4 band is obtained by the coupling of a higher aligned j particle, driving the system to an even larger deformation. One must remember that these are the *effective* MOIs of the entire system, that is the triaxial-rotor + odd-particle. No spin dependence has been inferred for the MOIs, so a possible change in the MOIs ordering with the increase in spin I cannot be studied within the current description. Furthermore, this formalism does not contain microscopic terms, so no presumptions on what causes the obtained MOI ordering can be stated. Although, by working with a quadrupole deformed mean-field, the moments of inertia of the triaxial core should be indeed consistent with the hydrodynamic model. For the sake of completeness, Figure 10 shows the evolution of a hydro-dynamical set of MOIs with respect to the triaxiality parameter γ .

Concerning the triaxiality parameter γ , it has a positive value $\gamma = 22^\circ$. This is consistent with the microscopic calculations based on cranking mechanism for the potential energy surface (PES) of ^{163}Lu (discussion on PES was done in the previous sections). In fact, the agreement is quite good with the predicted deformed minima of $(\beta_2, \gamma) \approx (0.38, 20^\circ)$ [13, 61]. Comparing the current W2 model with already existing de-

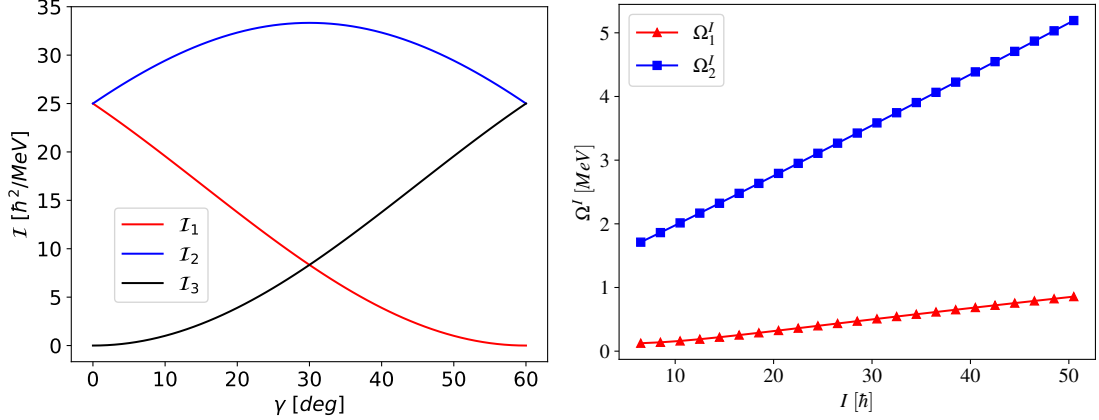


Figure 10: Left-side: The hydrodynamic moments of inertia [29] as a function of the triaxiality parameter γ , for the positive interval $\gamma \in [0^\circ, 60^\circ]$, evaluated for a scale factor $\mathcal{I}_0 = 25 \text{ MeV}^{-1}$. Right-side: The wobbling frequencies defined in Eq. 23 as function of total angular momentum, evaluated with the parameter set \mathcal{P} which was obtained through the fitting procedure.

scriptions which take γ to be fixed a-priori (e.g. [29, 62]), here γ is obtained through the fitting process in a self-consistent manner. Moreover, its value is slightly larger than the one obtained in W1 formalism ($\gamma = 17^\circ$). This might be due to the larger ratios $\mathcal{I}_1/\mathcal{I}_{2,3}$, which in the present case they appear to be bigger ($\mathcal{I}_1/\mathcal{I}_2 \approx 4.8$ for W2, compared to ≈ 3.2 in the previous approach).

Finally, the single-particle potential strength, which causes the odd-proton to move in the quadrupole deformed mean-field, has a value of $V = 2.1 \text{ MeV}$. In W1, this parameter was $V^{\text{TSD1,2,3}} = 3.1 \text{ MeV}$ and $V^{\text{TSD4}} = 0.7 \text{ MeV}$. An explanation for its decrease in the present case might be due to the upward shift in the energy caused by the un-favored partner, or due to the energetic shift of the parity partner, indicating a quenching effect on the quadrupole deformation of the triaxial system. Nevertheless, the obtained value seems to be consistent with the previous calculations. Other interpretations [62] that were developed using a similar single-particle potential term in the Hamiltonian adopted values of around $V = 1.6 \text{ MeV}$, however, that was for an isotope with smaller quadrupole deformation $\beta_2 = 0.18$. An interesting research devoted to the particle deformation in a single- j shell model aimed at obtaining a realistic expression for the deformation parameter was done in [85]. Therein, results for the potential strength of odd nuclei with similar mass, but different quasiparticle configuration were numerically obtained. As a future work, it is worth extending the current W2 formalism with the a more realistic term for the quadrupole deformation. Concluding this subsection, the obtained values of \mathcal{P} seem to not only describe the wobbling spectrum of ^{163}Lu very well (see spectrum in Figures 8 and 9), but they are also consistent with the previous formalism developed by the same team, or even with other interpretations from the literature.

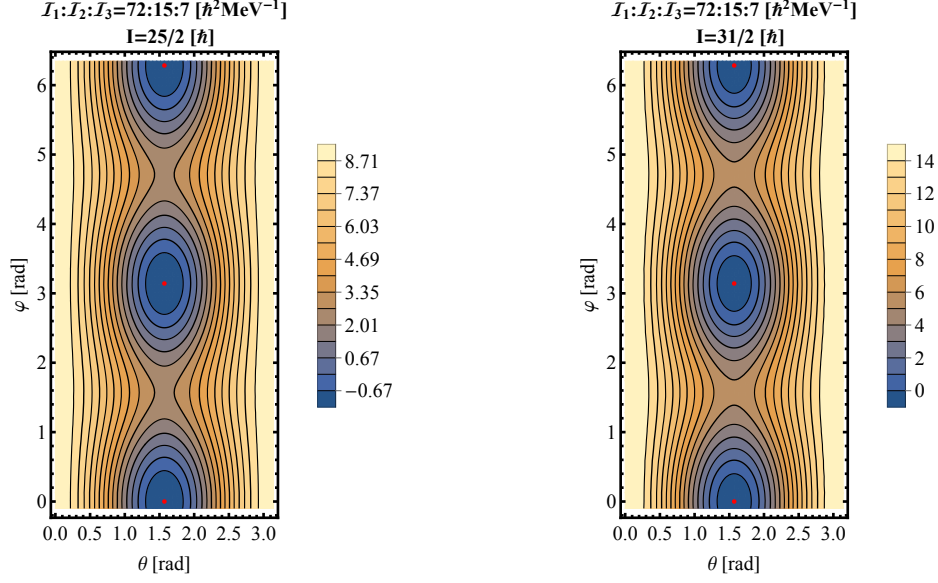


Figure 11: Contour plots with the energy function \mathcal{H} for a state in TSD_1 (left) and a state from TSD_2 (right). Calculations were performed with the numerical parameters obtained from the fitting procedure. The minimum points for \mathcal{H} are marked by red dots, and they represent the regions in space where the nucleus has a stable wobbling character. The blue *islands* also indicate a stable motion of the triaxial nucleus.

4.2 Stability of the wobbling region

The expression for the classical energy function, which plays a crucial role in analyzing the nucleus' stability for a given rotational state, was formulated in the previous section, with Eq. 34. This will be used within the present numerical calculations in order to pinpoint the regions in the space where the minimal points of \mathcal{H} arise. A special interest is devoted to the low-lying states from each of the four bands. Namely, for each band, a spin-state close to the band-head is chosen, then using the parameter set \mathcal{P} , a graphical representation in the (θ, φ) -coordinate space is realized, and in each case the extremal points with minimum character are identified. These graphical representations are shown in Figures 11 and 12.

The four contour plots shown in Figures 11 and 12 have many similarities, suggesting common collective properties, but also differences caused by the fact that the minima have different depths. A common feature consists in that the equi-energy curves surround a sole minimum for low values in energy, but as the energy increases, the trajectories go around all minima, the lack of localization indicating unstable wobbling motion. The unstable regions might also relate to phase transitions, where the nucleus can undergo a major change in its rotational character. This aspect will be also discussed in the next subsection for the case of a 3-dimensional representation of the energy ellipsoid and the classical trajectories of the triaxial system.

Regarding the minimum points (marked by red dots on the contour plots), their

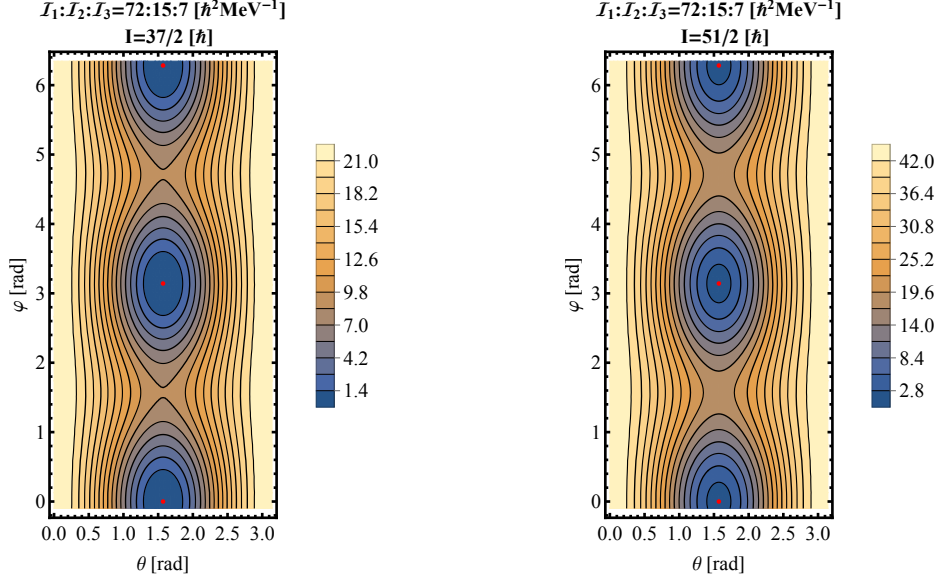


Figure 12: Contour plots with the energy function \mathcal{H} for a state in TSD_3 (left) and a state from TSD_4 (right). Calculations were performed with the numerical parameters obtained from the fitting procedure. The minimum points for \mathcal{H} are marked by red dots, and they represent the regions in space where the nucleus has a stable wobbling character. The blue *islands* also indicate a stable motion of the triaxial nucleus.

position remain unchanged for all four bands and for any rotational state I , as long as the MOI's ordering stays the same. Remarkable is the fact that only with the obtained set of parameters (the current MOI ordering) it was possible to define contours with stable motion (marked by the blue regions). Indeed, if the two ratios $\mathcal{I}_1/\mathcal{I}_2$ and $\mathcal{I}_2/\mathcal{I}_3$ would have been smaller, a larger unstable region would prevail (with islands of maximal character), constraining thus the stable wobbling motion. This could indicate the fact that the single-particle term $T_{s.p.}$ from \mathcal{H} actually *prefers* a larger triaxiality in order to achieve a stable motion characterized by large deformation (see Eq. 35).

An additional step consists in the analysis of the energy function, or more precisely to see its evolution in one of the minimal points with respect to the angular momentum I . As it was already observed from the contour plots, the depth of the minima differ from one spin state to another, so it would be useful to have a quantitative view on that change. By fixing \mathcal{H} in one of its critical points (e.g., the minimum $p_{\min}(\theta, \varphi) = (\frac{\pi}{2}, 0)$), the angular momentum I was varied within a large interval, and the evolution of \mathcal{H} was evaluated. Graphical representation is shown in Figure 13.

As it can be seen from Figure 13, the classical energy \mathcal{H} is an increasing function of angular momentum, which is to be expected, since the wobbling energies of the four bands increase with respect to the increase in spin. The negative values of \mathcal{H} for low-lying wobbling states do not indicate that the nucleus has negative energy states, since the rest of nucleus' energy is also given by the single-particle energy ϵ_j terms and the

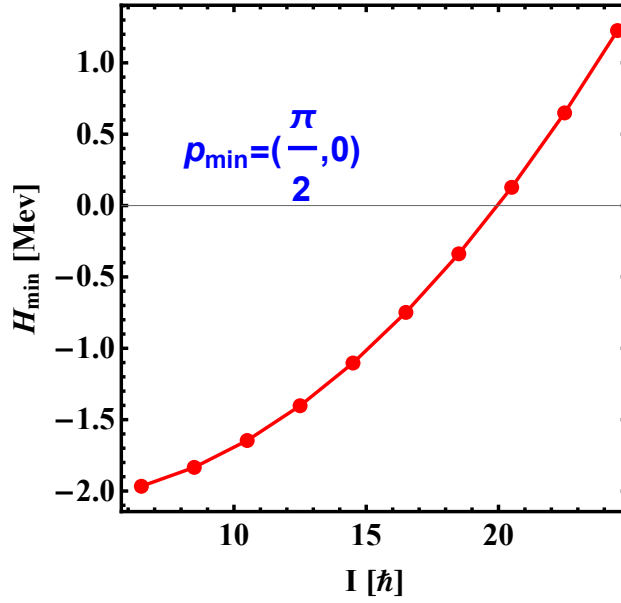


Figure 13: The change in the minimum depth of \mathcal{H} , for the given parameter set \mathcal{P} , evaluated in the point $(\theta, \varphi) = (\frac{\pi}{2}, 0)$.

phononic $\mathcal{F}_{n_{w1}n_{w2}}^I$ terms.

Another useful insight would be the study of the classical energy function \mathcal{H} for the obtained parameter set, as a function of the polar angles (θ, φ) . This can be achieved by choosing a minimum point, then keeping one of the polar coordinates fixed, and let the other one vary across its corresponding interval. For ^{163}Lu , such a graphical representation was done for the point $p_{\min} = (\frac{\pi}{2}, 0)$ (that is the bottom-most red dot from each of the four contour plots depicted in Figures 11 and 12). Results can be seen in Figure 14.

4.3 Comment on the wobbling nature of ^{163}Lu

It is worthwhile discussing the results obtained regarding the wobbling spectrum of ^{163}Lu . Indeed, using a fitting procedure which minimized the χ^2 function, it was possible to find a parameter set \mathcal{P} that when applied in the calculations for the excitation energies, an agreement with the experimental data is observed. However, in the current state, there is no clear evidence on whether the formalism W2 predicts a TW or a LW behavior for the nucleus. According to [27], the wobbling character is given by the coupling of the odd particle which aligns parallel (LW) or perpendicular (TW) to the axis with largest MOI. But in order to see this within the measured data, an interpretation of the wobbling energy as it was defined in Eq. 3 must be performed. Indeed, according to the definition, one has to subtract an energy state within the first excited wobbling band (the one-wobbling-phonon band) from the average of its adjacent energies that belong to the yrast partner. In the present calculations, the first excited state is TSD_3 , with its

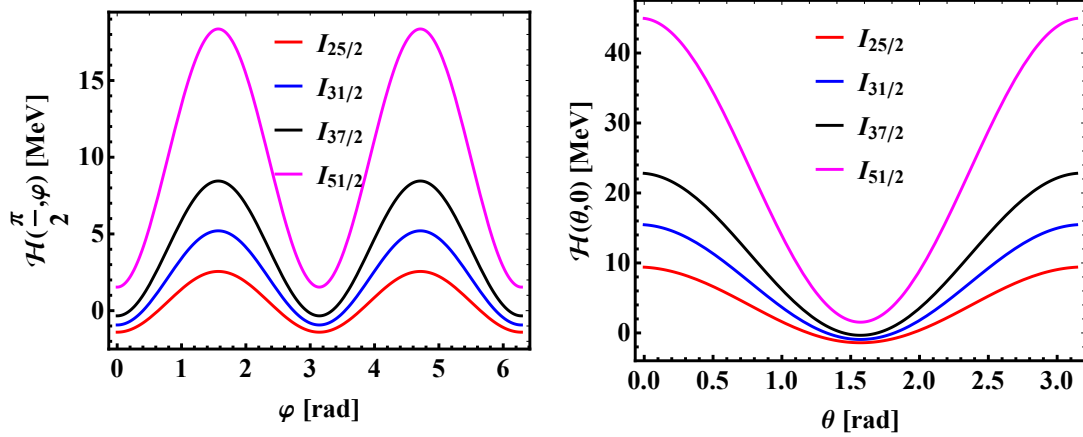


Figure 14: The energy function \mathcal{H} , evaluated in one of its minimum points, as a function of the polar coordinates. One coordinate is fixed while the other one is varied within its interval of existence. For $\theta \in [0, \pi]$ and $\varphi \in [0, 2\pi]$.

yrast partner being the band TSD_2 . Following this procedure, both the experimental wobbling energies, as well as the theoretical ones were calculated according to the Eq. 3. The obtained results are plotted in Figure 15.

From the behavior of E_{wob} from Figure 15, it can be seen that the theoretical wobbling spectrum is clearly an increasing function of angular momentum I , suggesting that ^{163}Lu would have a LW character. This contrasts the current interpretation on which the wobbling energies are decreasing functions with the increase in spin. However, within those formalisms [27, 63], the wobbling energies are calculated with TSD_2 and TSD_1 instead, since the one-wobbling-phonon band is TSD_2 , which is inconsistent with the W2 approach.

By contrast, analyzing the experimental data from Figure 15, a slight increase with spin can also be observed, also suggesting that the coupling in ^{163}Lu achieves a LW character. Indeed, from the lower limit of around $11/2 \hbar$ and up to a spin of about $39/2 \hbar$, the energy is increasing, then starts to decrease once $I \geq 39/2 \hbar$. The increasing behavior of the theoretical data also appears to be quenched in the high-spin limit, indicating that indeed, once the nucleus reaches high rotational states, a change in the wobbling regime might emerge, and the nucleus can transition from a wobbling regime (LW) to another (TW).

In a recent work [20], an observation of two wobbling bands was made for ^{183}Au (as it was already mentioned throughout the introduction). The bands are based on a positive $\pi(i_{13/2})$ and negative $\pi(h_{9/2})$ parity configuration. The remarking aspect of this research is that within a PRM model amended with the HFA (harmonic frozen alignment) approximation, it results that both bands have a transverse character (TW). However, they have different behavior in respect to the wobbling energies. Namely, this quantity increases (decreases) with spin for the positive (negative) parity configuration (see Figure 3 from [20]). According to the same study, the increasing trend of the positive

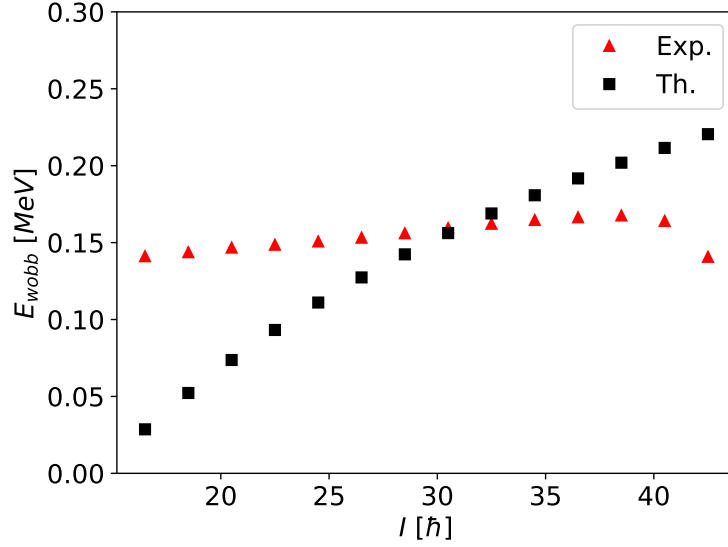


Figure 15: The wobbling energies for ^{163}Lu given as $E_{\text{wob}} = E_1(I) - \frac{1}{2}(E_0(I+1) + E_0(I-1))$. According to the current formalism, the sets of energies E_1 belong to TSD_3 , while E_0 correspond to TSD_2 . Experimental data is taken from [81]. Theoretical values were calculated with the parameter set \mathcal{P} .

parity band with spin has indeed a TW character, but it might be the *initial part* of a TW band predicted by Frauendorf [27] for ^{163}Lu , but never observed. This indicate that a increasing/decreasing behavior for E_{wob} is not enough evidence for asserting a wobbling character on a triaxial nucleus (at least, not without some strong constraints on the coupling scheme between the core and the odd particle).

Concluding this comment on the wobbling nature for ^{163}Lu , if the behavior of the wobbling energies with spin is the sole player in determining the wobbling character of a nucleus, then one could argue that indeed, based on the current results, ^{163}Lu behaves as a longitudinal wobblers. On the other hand, considering the newly obtained results discussed in the previous paragraph, the evidence is not enough for making a clear assumption on the type of wobbling motion which occurs.

4.4 Classical trajectories - 3-dimensional representation

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