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December 17, 2018

Physical Review C

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Triaxiality in selenium-76

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(Dated: January 23, 2019)

Background: Selenium and germanium nuclei are associated with both triaxiality and shape co-existence. The relative influence of these deformation effects on the low-lying nuclear structure remains the subject of much discussion, with additional attention drawn to ⁷⁶Se and ⁷⁶Ge due to the potential for the observation of neutrinoless double-beta decay.

Purpose: Experimental observables related to the deformation of ⁷⁶Se are lacking in precision. The purpose of the present work is to provide electric quadrupole matrix elements with improved precision in order to determine the deformation of low-lying states in a model-independent manner.

Methods: Sub-barrier Coulomb excitation was employed at the reaccelerated beam facility of the National Superconducting Cyclotron Laboratory using the JANUS setup. Using this method nineteen *E2* matrix elements were extracted.

Results: Extracted matrix elements agree within uncertainties with those in the literature but with improved precision. Through both a comparison with geometric models and a model-independent evaluation of *E2* matrix elements using rotational invariants the ground state of ⁷⁶Se is best described as having a significant triaxial component, while not being maximally triaxially deformed.

Conclusions: Selenium-76 exhibits a significant degree of triaxiality in its ground state. A detailed comparison with configuration interaction calculations indicates that this can be well reproduced theoretically.

I. INTRODUCTION

Atomic nuclei exhibit the bulk property of deformation, commonly assuming non-spherical shapes far from nuclear magic numbers. It is becoming increasingly clear that axially symmetric prolate and oblate extremes of quadrupole deformation are not sufficient to describe many nuclei, with triaxial components to the deformation often found to be significant. One area of the nuclear landscape where the role of axially asymmetric deformation has long been inferred to be important is the selenium-germanium region. The low-lying γ band in these isotopes - associated with a triaxial vibrational mode - is indicative of this, while the staggering of the band members has been used to infer the rigidity of this triaxial deformation (see e.g. Ref. [1]). More recently, comprehensive Coulomb excitation analyses using state-of-the-art γ -ray detection arrays have cemented this picture of triaxial deformation in, for example, ⁷²Ge [2].

Understanding the role of deformation in these isotopes is made a more pressing issue by its influence on the calculation of the nuclear matrix element for the hypothesised neutrinoless double-beta decay process, with ⁷⁶Ge being an excellent candidate for observation [3, 4]. Differing ground-state deformations in the parent (⁷⁶Ge) and daughter (⁷⁶Se) nuclei result in a reduced spatial overlap

between the initial and final state configurations. Any reduction in overlap will inhibit the process, as inferred more explicitly through theoretical calculations (see e.g. Ref. [5]).

Reproducing the germanium-selenium region microscopically provides a challenge for nuclear theory. The *fpg* shell itself is large, while small-amplitude excitations into the $2d_{5/2}$ orbital are known to play an important role in driving quadrupole deformation but are typically too computationally expensive to include in calculations. Over the last decade, however, Hamiltonians have been developed to tackle this region of the nuclear landscape with examples pertinent to this work being the JUN45 [6] and the *jj44b* (see appendix A in Ref. [7]) interactions.

In order to provide a clearer picture of the role of deformation in the selenium isotopes, more extensive and precise experimental data are required. In particular, transition matrix elements connecting low-lying states are known to be sensitive to deformation effects and might be reproduced theoretically by, for example, shell-model calculations with the aforementioned interactions. Here, we present results from a sub-barrier Coulomb excitation measurement of ⁷⁶Se performed on a high-*Z* target to allow for multi-step excitation. This methodology allows us access to multiple *E2* matrix elements connecting low-lying states, improving precision over literature values. The new, more precise results are then used to construct a series of rotationally invariant quantities related to the quadrupole deformation parameters of the nucleus, both with respect to the absolute magnitude of the ground-

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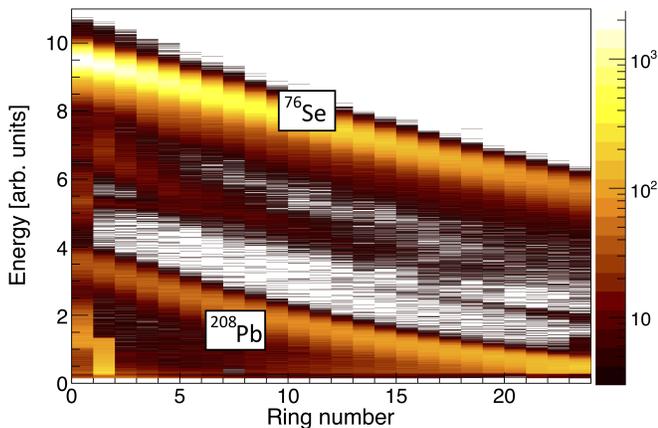


FIG. 1. Pixel energy in the JANUS silicon detectors against ring number with smaller ring numbers corresponding to shallower angles. Dominant in the spectrum are the kinematic lines originating from the scatter of ^{76}Se and ^{208}Pb . Also seen is a lighter contaminant at approximately a 1% level. Note that bins with two or fewer counts are excluded from this figure to aid clarity.

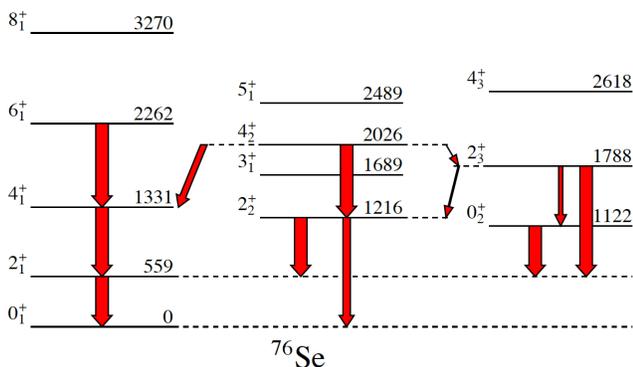


FIG. 2. Level scheme indicating the observed transitions. Note that the experiment is sensitive to unobserved transitions by including data from the literature such as branching ratios. Arrow widths correspond to relative branching ratios.

state deformation and its degree of triaxiality.

II. EXPERIMENTAL DETAILS

The present measurements were performed at the reaccelerated beam facility (ReA3 [8]) of the National Superconducting Cyclotron Laboratory (NSCL [9]). Stable ^{76}Se ions were injected into the NSCL's electron beam ion trap [10] and charge-bred prior to injection into the ReA accelerator chain. To enhance sensitivity to the nuclear matrix elements, ^{76}Se was delivered at two energies: 4.07 MeV/u and 3.55 MeV/u, maintained for 24 hours and 72 hours, respectively and impinged upon an enriched, 0.92-mg/cm² thick, ^{208}Pb target. Mean on-target intensities were maintained at approximately

$2 \cdot 10^5$ pps for the duration of the experiment. The beam energies corresponded to minimum impact parameters of 5 fm and 7.5 fm, 72% and 62% of the Coulomb barrier for the higher and lower energies respectively. Scattered beam- and target-like nuclei were detected in the JANUS setup for Coulomb excitation [11], consisting of a pair of S3-type annular silicon detectors located upstream and downstream of the target. The downstream and upstream target-silicon detector separations were measured to be 26 ± 1 mm and 34 ± 1 mm, respectively. Emitted γ rays were detected in the Segmented Germanium Array (SeGA) [12], consisting of sixteen 32-fold segmented HPGe detectors. Data were extracted using a digital data-acquisition system, made up of 33 100-MHz (SeGA) and 8 250-MHz (silicon detectors) XIA Pixie-16 modules in a triggerless, continuous-running mode. Events were constructed on the basis of the master clock and analyzed using the GRUTinizer [13] software package, built in a ROOT framework [14]. The experimental setup was identical to that described in Ref. [15].

III. ANALYSIS

Silicon pixels were constructed from time- and energy-coincident ring and segment hits in the S3 detectors. Beam-like (^{76}Se) and target-like (^{208}Pb) recoil loci were identifiable in the observed kinematics. Gamma rays detected in SeGA were Doppler corrected for ^{76}Se velocities event-by-event on the basis of the reaction kinematics and the beam- or target-like nature of the detected particle. Target-like events were excluded from the analysis at angles greater than 40° to avoid biasing of the data due to dead-layer effects in the silicon detectors. Detections in the upstream detector were similarly excluded for the lower beam energy. The observed particle energies are shown in Fig. 1 plotted against the downstream ring number. Clearly, beam- and target-like loci are well separated. Low-level contamination from a lighter ion is seen at roughly the 1% level. This contamination does not impact the analysis of the experimental data presented below.

The beam-like locus in the downstream S3 detector was subdivided into six angular bins, each composed of four annular rings. In the case of the ^{208}Pb -gated kinematics, the data were subdivided into two angular bins containing six rings, corresponding to the twelve shallowest angle rings. Wider angle target-like events were excluded because of the dead-layer effects discussed previously. The upstream detector was considered as a whole in the higher-energy experimental setting. Doppler-corrected γ -ray spectra were constructed from SeGA coincidences with these angular bins for both the high- and low-energy portions of the experiment. The experimental data were thereby separated into seventeen groups for Coulomb-excitation analysis. Example Doppler-corrected γ -ray spectra, corresponding to beam- and target-like detection at the higher beam energy are shown in Fig. 3. The

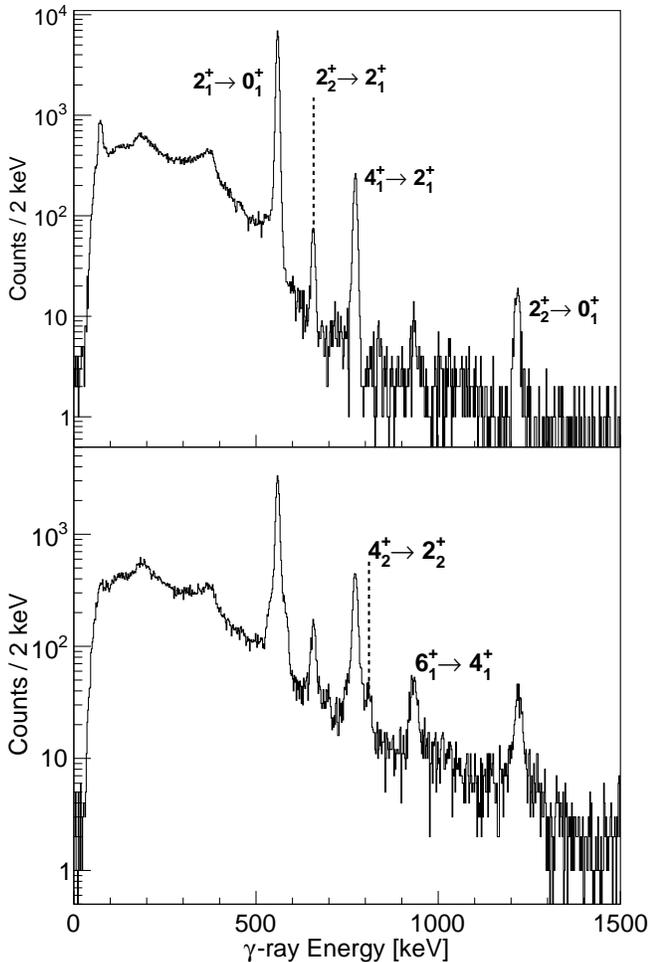


FIG. 3. Doppler-corrected γ rays coincident with the detection of a beam-like (top) and target-like (bottom) particle in the downstream silicon detector at a beam energy of 4.07 MeV/u with the strongest transitions highlighted. The non-Gaussian shapes in coincidence with the detection of target-like particles arise from the wide scattering angles of the beam-like particles in coincidence, resulting in them traversing a larger distance in the target material and inducing a line-shape.

Doppler correction is superior for beam-like particle detection. In the case of target-like detection, the beam-like particle is scattered at angles approaching 90° and therefore travels further in the target material, causing a broadening of the lineshape. Figure 4 shows example Doppler-corrected γ -rays detected in coincidence with a particle in the upstream detector. This detection condition corresponds to a best-case scenario for Doppler correction. In the general case, multiplets such as those shown in the bottom two panels of Fig. 4 were unresolvable.

The $0_2^+ \rightarrow 2_1^+$ and $2_1^+ \rightarrow 0_1^+$ transitions (bottom panel, Fig. 4) are separated by only 4 keV and, as such, cannot be resolved in coincidence with downstream detection

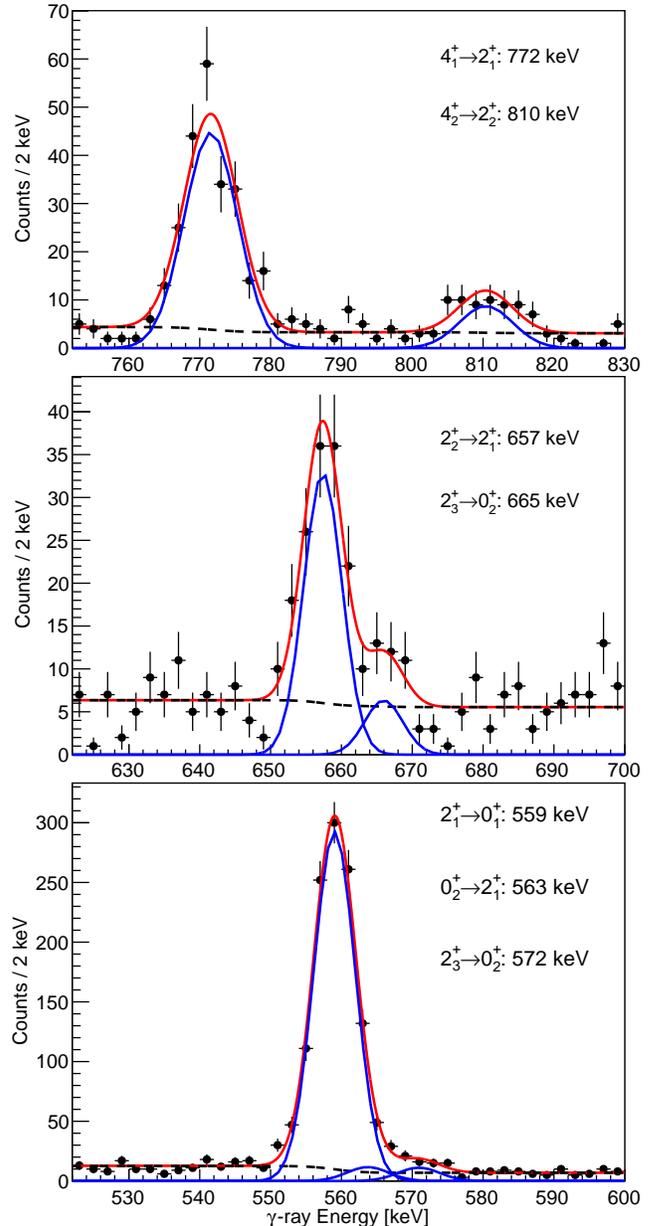


FIG. 4. (Color online) Doppler-corrected γ rays coincident with the detection of a beam-like particle in the upstream silicon detector. Shown are three cases in which doublets were observed and fitted, with the total fit (red), individual peaks (blue) and background (dashed black) shown.

due to the larger velocities. In order to enhance sensitivity both to the multi-step excitation: $0_1^+ \rightarrow 2_1^+ \rightarrow 0_2^+$, and to the quadrupole moment of the 2_1^+ state it is desirable to quantify the relative contributions to the combined peak around 560 keV. Utilizing the high efficiency of the SeGA array this was achieved using $\gamma - \gamma$ coincidences. In order to use $\gamma - \gamma$ coincidences it is necessary to account for changes in detection efficiency as a result of the $\gamma - \gamma$ angular distribution. This efficiency correc-

tion was determined empirically through comparison to cascades both following the β decay of ^{152}Eu and through comparison with the γ -ray spectrum at backward angles where, as shown in the bottom panel of Fig. 4, the two transitions were resolvable. The correction factor was then taken as a weighted average of that determined from the two empirical methods. The uncertainty on the values extracted in this method remain dominated by the statistical component.

Coulomb-excitation yields were evaluated using the GOSIA code [16] to determine transition matrix elements. Literature experimental values pertaining to the transitions observed in the present work were additionally used to constrain the fit and are summarized in Table I. The $E0$ decay of the 0_2^+ state was incorporated through decay to a fictional 1^+ state via an $M1$ decay, constrained by the measured branching ratio. Consistency between the high- and low-energy data was confirmed by performing an independent GOSIA analysis for each. The final results presented below are from a simultaneous GOSIA analysis, with matrix elements minimized to best reproduce both the high- and low-energy data.

TABLE I. Experimental values as constraints for the GOSIA calculations presented here. Values taken from Ref. [17] except where indicated.

State	Lifetime [ps]	Transition	Mixing ratio (δ)
2_1^+	17.75(29)	$2_2^+ \rightarrow 2_1^+$	5.2(2)
2_2^+	4.01(29)	$2_3^+ \rightarrow 2_2^+$	0.80(55) [†]
4_1^+	2.19(8)	$2_3^+ \rightarrow 2_1^+$	-0.51(5)
2_3^+	1.86(46) [†]	$4_2^+ \rightarrow 4_1^+$	1.7(4)
6_1^+	0.89(10)		

Initial (i)	Final (f_1)	Final (f_2)	BR [$\frac{I(i \rightarrow f_1)}{I(i \rightarrow f_2)}$]
0_2^+	0_1^+	2_1^+	$2.3(4) \cdot 10^{-4}$
2_2^+	0_1^+	2_1^+	0.59(3)
2_3^+	2_1^+	0_1^+	4.00(23)
2_3^+	2_1^+	0_2^+	3.13(23)
2_3^+	2_1^+	2_2^+	10(2)
2_3^+	2_1^+	4_1^+	33(3)
4_2^+	2_2^+	4_1^+	2.04(23)
4_2^+	2_2^+	2_1^+	20(12)

[†] - from Ref. [18]

IV. DISCUSSION

The matrix elements resulting from the present analysis are given in Table II. Note that as previously mentioned, literature data were used to constrain a number of matrix elements and so the present values should not be considered as independent of those presented in Table I. $E2$ matrix elements and transition strengths can be used to infer the collective behaviour of the nucleus

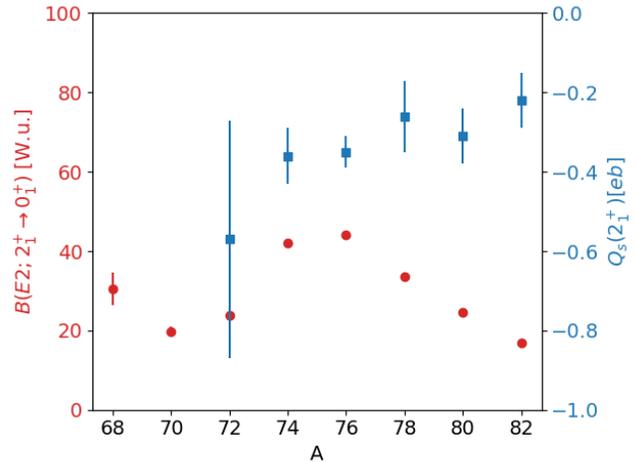


FIG. 5. $B(E2; 2_1^+ \rightarrow 0_1^+)$ and $Q_s(2_1^+)$ values for selenium isotopes. A shallow maximum in absolute $B(E2)$ and $Q_s(2_1^+)$ is seen to occur around ^{76}Se . Data taken from Refs. [15, 17].

through a number of relations. ^{76}Se has the largest $B(E2; 2_1^+ \rightarrow 0_1^+)$ value, indicating a maximum of collectivity, along with the lowest 2_1^+ excitation energy, another indicator of a large degree of collectivity. This is also borne out by the systematics of the spectroscopic quadrupole moments, $Q_s(2_1^+)$, with maximum absolute values approximately centered on ^{76}Se , as shown in Fig. 5, however this picture is far from complete. Under the assumption of a rigid axial rotor, one can relate $|Q_s(2_1^+)|$ to the $B(E2; 2_1^+ \rightarrow 0_1^+)$ by:

$$|Q_s(2_1^+)| = \frac{2}{7} \sqrt{16\pi \cdot B(E2; 2_1^+ \rightarrow 0_1^+)}. \quad (1)$$

The value of $|Q_s(2_1^+)|$ determined in the present work is only approximately 60% of that determined from the above relation. Coupled with the existence of a low-lying 2_2^+ state which can be best described as the band-head of the so-called γ -band, this indicates that a simple symmetric axial rotor description is insufficient to explain the low-lying structure and triaxiality likely plays a significant role in ^{76}Se .

One can compare the level energies and transition strengths of the low-lying states to, for example, the Davydov-Filippov model [22] to estimate the degree of triaxiality in the nuclear system from the assumption of an axially asymmetric rotor. Such an analysis is presented in Fig. 6 for ^{76}Se . Clearly, the level energies and transition strengths are best described by a nucleus approaching maximum triaxiality ($\gamma \sim 25 \rightarrow 30^\circ$). Note that the behaviour of level energies and transition strengths in the model is symmetric about $\gamma = 30^\circ$. Using the triaxial rotor model presented in Ref. [23], in which the inertia tensor is treated independently of the electric quadrupole tensor, one can calculate a value for the triaxiality parameter, $\gamma \approx 25^\circ$ - in good qualitative agreement with the value expected from Fig. 6.

TABLE II. Reduced matrix elements, reduced transition strengths and quadrupole moments determined in the present work. Where available literature values are provided for comparison. Note that due to the use of experimental observables in the GOSIA minimization the present results cannot be considered as independent of the literature. Note that $B(E2) \uparrow$ and $B(M1) \uparrow$ values are given.

This Work			Literature		
	$\langle J_i^\pi E2 J_f^\pi \rangle$ [eb]	$B(E2; J_i^\pi \rightarrow J_f^\pi)$ [$e^2\text{fm}^4$]	$\langle J_i^\pi E2 J_f^\pi \rangle$ [eb]	$B(E2; J_i^\pi \rightarrow J_f^\pi)$ [$e^2\text{fm}^4$]	Reference
$0_1^+ \rightarrow 2_1^+$	0.647 (5)	4190 ($\frac{60}{70}$)	0.649 (7)	4210 (68)	[17]
$0_1^+ \rightarrow 2_2^{\dagger}$	0.110 (2)	119 (5)	0.108 ($\frac{4}{5}$)	116 (10)	[17]
$0_1^+ \rightarrow 2_3^+$	0.040 (1)	16 (1)	0.044 ($\frac{10}{14}$)	19 (10)	[18]
$2_1^+ \rightarrow 0_2^+$	0.285 ($\frac{26}{27}$)	163 ($\frac{30}{29}$)	0.300 ($\frac{64}{81}$)	180 (84)	[17]
			0.47 ($\frac{11}{10}$)	440 ($\frac{230}{168}$)	[19]
$2_1^+ \rightarrow 2_2^+$	0.640 (11)	820 (30)	0.641 ($\frac{22}{23}$)	822 (57)	[17]
$2_1^+ \rightarrow 4_1^+$	1.108 ($\frac{12}{11}$)	2450 (50)	1.105 ($\frac{15}{16}$)	2444 (69)	[17]
$2_1^+ \rightarrow 2_3^+$	-0.093 ($\frac{7}{6}$)	17 (2)	0.102 ($\frac{14}{16}$)	21 (6)	[18]
$2_1^+ \rightarrow 4_2^+$	0.039 ($\frac{35}{7}$)	3 ($\frac{8}{1}$)	0.04 (1)	2.4 (17)	[17]
$0_2^+ \rightarrow 2_2^+$	0.182 ($\frac{33}{47}$)	330 ($\frac{130}{150}$)	0.15 ($\frac{8}{18}$)	225 ($\frac{304}{225}$)	[19]
$0_2^+ \rightarrow 2_3^+$	0.532 ($\frac{21}{18}$)	2830 ($\frac{220}{190}$)	0.595 ($\frac{7}{9}$)	3540 ($\frac{1910}{960}$)	[18]
			0.59 ($\frac{30}{74}$)	3481 ($\frac{4440}{3481}$)	[19]
$2_2^+ \rightarrow 4_1^+$	0.047 ($\frac{44}{29}$)	4.5 ($\frac{122}{38}$)	0.12 ($\frac{16}{12}$)	31 ($\frac{131}{31}$)	[19]
$2_2^+ \rightarrow 2_3^+$	0.262 ($\frac{67}{43}$)	140 ($\frac{80}{40}$)	0.19 ($\frac{6}{4}$)	73 ($\frac{57}{27}$)	[18]
			0.30 ($\frac{10}{6}$)	880 ($\frac{650}{325}$)	[19]
			< 0.022	< 4.8	[17]
$2_2^+ \rightarrow 4_2^+$	0.768 ($\frac{36}{37}$)	1180 (110)	0.706 ($\frac{81}{91}$)	998 (241)	[17]
$4_1^+ \rightarrow 2_3^+$	0.418 ($\frac{18}{15}$)	190 ($\frac{20}{10}$)	0.489 ($\frac{6}{7}$)	266 (74)	[18]
$4_1^+ \rightarrow 4_2^+$	0.733 ($\frac{51}{38}$)	600 ($\frac{90}{60}$)	0.62 ($\frac{9}{11}$)	421 (134)	[17]
$4_1^+ \rightarrow 6_1^+$	1.390 ($\frac{64}{58}$)	2050 ($\frac{200}{180}$)	1.3 ($\frac{8}{7}$)	1880 (220)	[17]
	$\langle J_i^\pi M1 J_f^\pi \rangle$ [μ_n]	$B(M1; J_i^\pi \rightarrow J_f^\pi)$ [μ_N^2]	$\langle J_i^\pi M1 J_f^\pi \rangle$ [μ_n]	$B(M1; J_i^\pi \rightarrow J_f^\pi)$ [μ_N^2]	Reference
$2_1^+ \rightarrow 2_2^+$	0.067 (3)	$9.1 (\frac{0.8}{0.7}) \cdot 10^{-4}$	0.068 (3)	$9.2 (9) \cdot 10^{-4}$	[17]
$2_1^+ \rightarrow 2_3^+$	0.186 ($\frac{4}{5}$)	$7.0 (3) \cdot 10^{-3}$	0.16 (2)	$5 (1) \cdot 10^{-3}$	[18]
$2_2^+ \rightarrow 2_3^+$	0.168 ($\frac{4}{2}$)	$6 (\frac{3}{1}) \cdot 10^{-3}$	0.14 ($\frac{4}{3}$)	$4 (2) \cdot 10^{-3}$	[18]
$4_1^+ \rightarrow 4_2^+$	0.158 ($\frac{5}{3}$)	$3 (\frac{2}{1}) \cdot 10^{-3}$	0.13 ($\frac{4}{3}$)	$2 (1) \cdot 10^{-3}$	[18]
	$\langle J_i^\pi E2 J_i^\pi \rangle$ [eb]	$Q_s(J^\pi)$ [eb]	$\langle J_i^\pi E2 J_i^\pi \rangle$ [eb]	$Q_s(J^\pi)$ [eb]	Reference
$2_1^+ \rightarrow 2_1^+$	-0.463 ($\frac{52}{53}$)	-0.35 (4)	-0.449 (92)	-0.34 (7)	[20]
			-0.396 (106)	-0.30 (5)	[21] via [17]
			-0.45 (7)	-0.34 (5)	[19]
$2_2^+ \rightarrow 2_2^+$	0.245 ($\frac{57}{60}$)	0.19 (4)	0.24 ($\frac{6}{8}$)	0.18 ($\frac{5}{8}$)	[19]
$4_1^+ \rightarrow 4_1^+$	-0.387 ($\frac{55}{53}$)	-0.29 (4)	-0.36 ($\frac{24}{14}$)	-0.27 ($\frac{18}{11}$)	[19]

While comparison with geometric models therefore indicates that triaxiality plays a role in the low-lying structure of ^{76}Se , the quality of the present data combined with that in the literature allows for a more rigorous analysis with the presented set of $E2$ matrix elements in Table II sufficient to perform a model-independent analysis of the nuclear shape.

The electric multipole transition operator is a spherical tensor allowing for rotationally invariant, zero-coupled products to be constructed which can themselves be linked to the intrinsic deformation of the nucleus [24, 25]. The two invariant products discussed here relate to the quadrupole deformation parameters of the nucleus. The first quadrupole invariant to be constructed, Q , corresponds to the overall deformation of the nucleus, in analogy to the β parameter of the Bohr Hamiltonian, but relating to the distribution of charge rather than mass. By

expanding over all intermediate states using the Wigner $6j$ symbol, one can write:

$$\frac{\langle Q^2 \rangle}{\sqrt{5}} = \frac{(-1)^{2I_i}}{\sqrt{(2I_i + 1)}} \sum_t \langle i || E2 || t \rangle \langle t || E2 || i \rangle \begin{Bmatrix} 2 & 2 & 0 \\ I_i & I_i & I_t \end{Bmatrix}. \quad (2)$$

The second invariant to be considered, δ , relates to the degree of triaxiality in the intrinsic frame, in analogy to Bohr's γ parameter, but again relating to the charge distribution. This is a higher-order invariant which, expanding again over all intermediate states, yields:

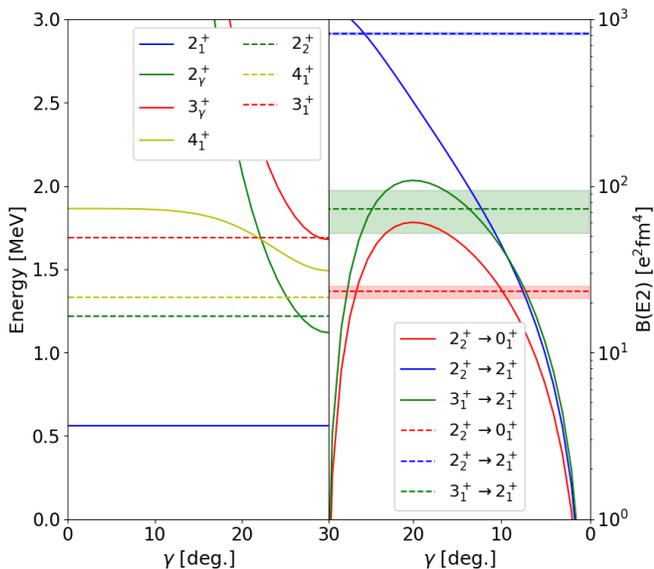


FIG. 6. (Color online) Level energies (left) and transition strengths (right) expected for the asymmetric rotor, Davydov-Filippov model (solid lines) as a function of the triaxiality parameter, γ . Experimental values (dashed lines) for ^{76}Se are shown, with shaded regions indicating the uncertainties. The experimental data are best described by a triaxiality parameter, $\gamma \approx 25^\circ$. Note that the behaviour of both energies and transition strengths is symmetric about $\gamma = 30^\circ$. Level energies and transition strengths were normalized to reproduce the 2_1^+ excitation energy and $2_1^+ \rightarrow 0_1^+$ transition strength, respectively.

$$\sqrt{\frac{2}{35}} \langle Q^3 \cos(3\delta) \rangle = \mp \frac{(-1)^{2I_i}}{(2I_i + 1)} \sum_{t,u} \langle i || E2 || u \rangle \langle u || E2 || t \rangle \langle t || E2 || i \rangle \begin{Bmatrix} 2 & 2 & 2 \\ I_i & I_t & I_u \end{Bmatrix}. \quad (3)$$

Where a negative sign corresponds to integer spins and a positive sign to half-integer. Here, we further assume that:

$$\langle \cos(3\delta) \rangle \approx \frac{\langle Q^3 \cos(3\delta) \rangle}{(\langle Q^2 \rangle)^{3/2}}. \quad (4)$$

Invariants were thereby determined for stable selenium isotopes and the $\langle \delta \rangle$ and $\langle Q \rangle$ parameters are presented in Fig. 7 in $(\langle Q \rangle, \langle \delta \rangle)$ space, in analogy to the (β, γ) space often used to represent the Bohr deformation parameters. Note that $\langle 2_3^+ | E2 | 2_3^+ \rangle$ has not been determined in any stable selenium isotope. The contribution of this matrix element to the ground-state $\langle \cos(3\delta) \rangle$ values is expected to be small, due to $\langle 0_1^+ | E2 | 2_3^+ \rangle$ being small in all cases where it has been measured. In the case of the 0_2^+ state however, the contribution can be significant due to the larger magnitude of the $\langle 0_2^+ | E2 | 2_3^+ \rangle$ matrix ele-

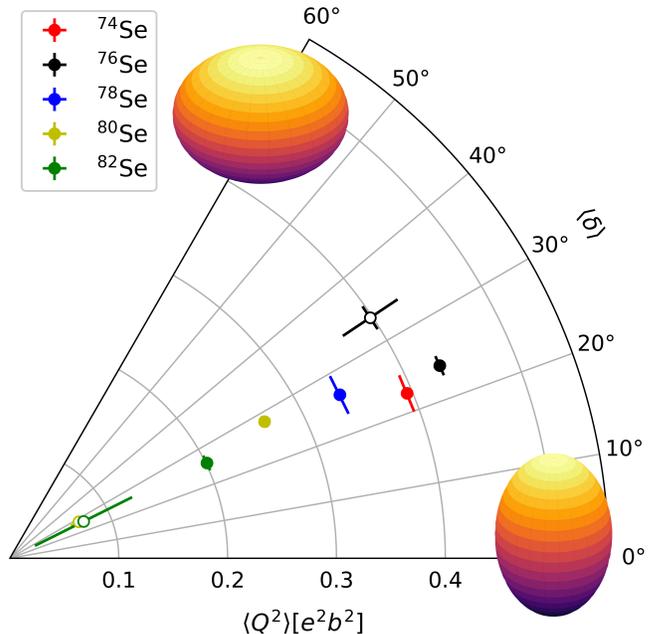


FIG. 7. (Color online) $\langle Q^2 \rangle$ and $\langle \delta \rangle$ calculated using the invariant values described in equations 2 and 3 for the 0_1^+ (solid symbols) and 0_2^+ (open symbols) states in stable selenium isotopes. ^{76}Se values were deduced from the matrix elements determined in the present work, while $^{74,78,80,82}\text{Se}$ were determined from data presented in Refs. [17, 19, 26–28]. Note that, as discussed in the text, uncertainties correspond to matrix-element uncertainties only and do not incorporate any uncertainties in convergence.

ment. The $\langle \delta \rangle$ values for the 0_2^+ states presented in Fig. 7 (open symbols) might therefore not be converged.

The $\langle \cos(3\delta) \rangle$ parameter, as previously stated, relates to the degree of triaxiality of the state in question - in this case the 0_1^+ and 0_2^+ states. A value of $\delta = 60^\circ$ ($\langle \cos(3\delta) \rangle = -1$) corresponds to an axially symmetric oblate shape, and a value of $\delta = 0^\circ$ ($\langle \cos(3\delta) \rangle = 1$) to an axially symmetric prolate shape. The $\langle Q \rangle$ parameter, meanwhile, relates to the absolute magnitude of deformation. From Fig. 7, therefore, it is clear that while the absolute degree of deformation in selenium isotopes increases from $A = 82 \rightarrow 74$, the nuclear shapes never stabilize into an axially symmetric rotor. Instead, the nuclei exhibit a significant - albeit not maximal - degree of triaxiality. Note that the presented invariant values are only sensitive to the average degree of triaxiality and not to the degree of rigidity of the triaxial deformation. Higher order invariant quantities than those expressed in Eq. 2 and 3 can also be constructed to quantify the rigidity of the Q and δ values, but require an even larger collection of matrix elements than presented here.

TABLE III. Experimental $B(E2)$ transition strengths and spectroscopic quadrupole moments for ^{76}Se compared to those calculated using the configuration interaction in the $jj44$ model space with two different interactions using an isoscalar effective charge: $e^\pi = 1.8, e^\nu = 0.8$. See text for details.

$B(E2; I \rightarrow F)$ [$e^2 fm^4$]	Expt.	$jj44b$	JUN45
$2_1^+ \rightarrow 0_1^+$	842 (14)	788	678
$2_2^+ \rightarrow 0_1^+$	23.2 (2)	2.4	44.2
$2_2^+ \rightarrow 2_1^+$	822 (57)	1073	273
<hr/>			
$Q_s(I)$ [eb]			
2_1^+	-0.34 (4)	+0.08	+0.49
2_2^+	+0.19 (4)	-0.09	-0.42
4_1^+	-0.29 (4)	+0.27	+0.27

A. Shell-Model Calculations

A microscopic comparison to the present data can be made through comparison with configuration interaction (CI) calculations in the so-called $jj44$ model space, which is made up of the $0f_{5/2}, 1p_{3/2}, 1p_{1/2}$ and $0g_{9/2}$ orbitals for both protons and neutrons. Calculations were performed using the shell-model code NuShellX [29] using both the JUN45 and $jj44b$ Hamiltonians with an isoscalar effective charge of $e^\pi + e^\nu = 2.6$. These calculations were previously presented more broadly for low-lying states in ^{76}Se [18] and ^{76}Ge [7], with a detailed description of the calculations presented in Appendix A of the latter reference. Here, we will focus on the transition strengths most pertinent to the present results.

Table III shows transition strengths for the lowest lying 2^+ states and quadrupole moments for the $2_1^+, 2_2^+$ and 4_1^+ states determined both experimentally and in the CI calculations. The $2_1^+ \rightarrow 0_1^+$ transition strength is relatively well reproduced, with the $jj44b$ Hamiltonian providing the best agreement. Similarly, the $jj44b$ interaction provides the best reproduction of the $2_2^+ \rightarrow 2_1^+$ transition strength, albeit overestimating the relative strength as compared to the $2_1^+ \rightarrow 0_1^+$ transition. The $2_2^+ \rightarrow 0_1^+$ strength is underpredicted by the $jj44b$ interaction, however, with the JUN45 providing the better agreement. If one were to interpret these results in the context of the Davydov-Filippov model (Fig. 6), the $jj44b$ calculations would appear to predict near maximal triaxiality, while the calculations using the JUN45 Hamiltonian predict a rather less triaxial structure - albeit not axially symmetric as evidenced by non-negligible $2_2^+ \rightarrow 2_1^+$ strength. This interpretation is strengthened by comparison of the quadrupole moments, with the JUN45 calculations resulting in larger absolute values, while the $jj44b$ values are considerably smaller. Investigation of the quadrupole moments appears to indicate a discrepancy between experiment and both calculations however, with predictions of a positive quadrupole moment - indicative of an oblate deformation - being in conflict with the experimentally determined negative value (prolate).

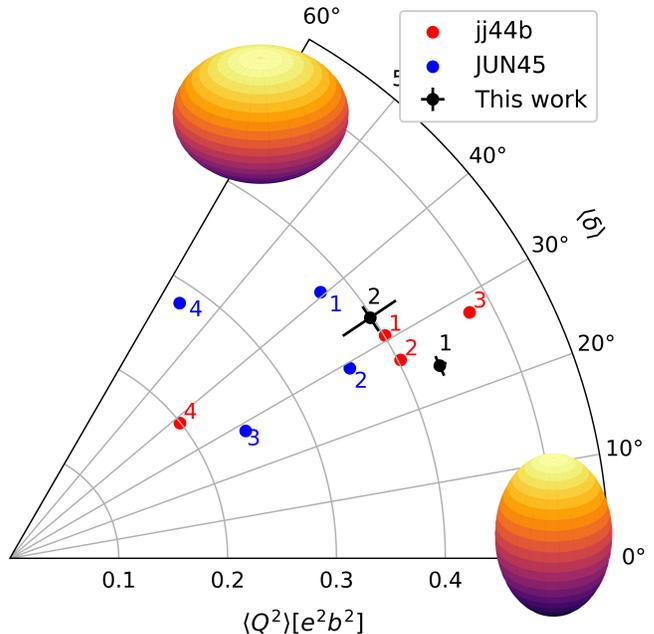


FIG. 8. (Color online) $\langle Q^2 \rangle$ and $\langle \delta \rangle$ calculated using the invariant values described in equations 2 and 3. Present experimental results for ^{76}Se (black points) along with values determined from shell-model calculations performed using the $jj44b$ and JUN45 interactions. Numbered points indicate the energy ordering of the 0^+ states (i.e. 1 corresponds to 0_1^+ , 2 to 0_2^+ , etc.). Uncertainties correspond to matrix-element uncertainties only and do not incorporate any uncertainties in convergence. See text for details.

To investigate further, the rotational invariants in equations 2 and 3 were constructed from the calculated matrix elements. 2^+ states up to 2_5^+ were included in the determination of the invariants, increasing confidence in convergence. As before, we present these calculated values in Fig. 8 in $(\langle Q^2 \rangle, \langle \delta \rangle)$ space. $\langle Q \rangle$ and $\langle \delta \rangle$ values were calculated for 0_{1-4}^+ and are compared to those calculated from experimentally determined matrix elements for the 0_1^+ and 0_2^+ states in ^{76}Se . The $jj44b$ calculations appear to provide the best description of the two lowest-lying 0^+ states, albeit with the ground state exhibiting a modest dominance by oblate deformation as compared to prolate deformation in the experimental data. Nonetheless, the agreement is considerably better than might be assumed if one were to merely compare the signs and magnitudes of the $Q_s(2_1^+)$ values. Due to the more comprehensive collection of $E2$ matrix elements determined from the CI calculations, higher order invariants can be constructed to assess the degree of rigidity in $\langle Q^2 \rangle$ and $\cos(3\delta)$. It is found that the $jj44b$ calculations correspond to a very δ soft structure, while the JUN45 calculations result in a more rigid configuration.

V. CONCLUSIONS

A low-energy Coulomb excitation measurement of ^{76}Se was performed using the JANUS setup at the ReA3 facility of the NSCL. A number of electric quadrupole matrix elements were extracted using the GOSIA code, with good agreement between the present results and those in the literature, where available. The spectroscopic quadrupole moment of the first 2^+ state was measured at improved precision, remaining consistent with a prolate deformation. An analysis of rotational invariants was performed to probe the intrinsic structure of the bands built upon the 0_1^+ and 0_2^+ states. The ground-state is found to exhibit significant triaxiality, albeit with a dominant prolate component. The 0_2^+ state is also found to be consistent with a triaxial deformation; however the present measurement was unable to determine the diagonal matrix element of the 2_3^+ state which, due to the large magnitude of $\langle 0_2^+ | E2 | 2_3^+ \rangle$, will likely contribute significantly to the triaxial invariant of the 0_2^+ state.

Comparison with configuration interaction calculations were performed. The present result highlights the need for a detailed analysis of theoretical results before comparison with experiment. A first-order analysis might imply that the CI calculations failed dramatically to reproduce the experimental data, due to the differing signs of the spectroscopic quadrupole moments. The

construction of invariants using the calculated $E2$ matrix elements, however, demonstrate that the calculations in fact reproduce the structure relatively well, as shown in Table III. A more comprehensive set of $E2$ matrix elements for the even-even selenium isotopes might allow further investigation of the role of triaxiality in the excited 0^+ states, which cannot be conclusively determined in the present work.

Previous studies into the effects of deformation on $0\nu 2\beta$ matrix elements have focussed on axial shapes (e.g. [5]). The present result, coupled to the low-lying structure of ^{76}Ge which also points towards triaxiality indicates that future investigations should take non-axial shapes into account.

ACKNOWLEDGMENTS

The authors thank the ReA3 beam delivery group at the NSCL for the high-quality beam of ^{76}Se . The authors acknowledge discussions with S. W. Yates. Work at LLNL was performed under contract DE-AC52-07NA27344. This work was supported in part by the National Science Foundation under Contract No. PHY-1565546 (NSCL), by the US Department of Energy, Office of Nuclear Physics, under Grant No. DE-FG02-08ER41556 (NSCL) and by the DOE National Nuclear Security Administration through the Nuclear Science and Security Consortium, under Award No. DE-NA000318.

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