Distribution of Low-Lying Quadrupole Phonon Strength in Nuclei

N. Pietralla,¹ P. von Brentano,¹ R. F. Casten,^{1,2} T. Otsuka,³ and N. V. Zamfir^{2,4,5}

¹Institut für Kernphysik, Universitat zu Köln, D-50937 Köln, Germany

²Brookhaven National Laboratory, Upton, New York 11973

³Department of Physics, University of Tokyo, Hongo, Bunkyo-ku, Tokyo 113, Japan

⁴Clark University, Worcester, Massachusetts 01610

⁵Institute of Atomic Physics, Bucharest-Magurele, Romania

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The distribution of quadrupole phonon strength to low-lying states in nuclei is examined. It is found that the 2_1^+ and 4_1^+ states of all nuclei from Z = 30 to 100 display almost perfect quadrupole phonon character even though they encompass widely ranging structures. The interacting boson approximation reproduces both the very small deviations from phonon purity and the distribution of those deviations as a function of mean field structure in a natural fashion. Simple physical mechanisms explain these results.

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There has recently been renewed interest in the applicability of phonon models [1] in nuclear structure. Examples of phonon and multiphonon behavior have been found [2-10] in a variety of contexts ranging from spherical to transitional and deformed nuclei and from yrast states to intrinsic collective modes. The growing recognition that phonon modes are widespread, and that multiphonon excitations often survive intact in the low degeneracy nuclear environment, motivates a further study of the empirical and theoretical phonon structure and interrelationships of the low-lying yrast levels of collective nuclei. In particular, it was shown in Ref. [4] that an anharmonic vibrator, with constant anharmonicity, can describe the yrast energies of all nonrotational collective nuclei from Z = 30 to 82. It is therefore pertinent to ask if it is possible to directly test this phonon picture, that is, to assess whether or not the wave functions of the low-lying levels of collective even-even nuclei can be accurately described in terms of phonon excitations.

It is therefore the purpose of this Letter to use B(E2) values to show that the 2_1^+ and 4_1^+ levels for all nuclei from Z = 30 to 100 display excellent quadrupole phonon purity and that this result is a natural property of the interacting boson approximation (IBA).

A quadrupole phonon excitation $|2_1^+2\rangle$ of the ground state is defined by

$$|2_{1}^{+}2\rangle_{\rm ph} \equiv NQ_{22}|0_{1}^{+}\rangle, \tag{1}$$

where Q_{22} is the quadrupole operator, N is a normalization factor, and the labels refer to spin, parity, and magnetic substate. Multiple "Q-phonon" states are created by repeated application of Q. Note that such states can be called phonons even for rotational excitations in that the phonon accelerates or decelerates the rotation.

This Q-phonon picture can be tested empirically or theoretically. If

$$B(E2:0_1^+ \to 2_i^+) = 0 \text{ for all } i > 1,$$
 (2)

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then Eq. (1) is satisfied for the 2_1^+ state. However, if the phonon strength is distributed over several 2^+ states, that is, if

$$|2^+2\rangle_{\rm ph} = \beta_1 |2_1^+\rangle + \sum_{i>1} \beta_i |2_i^+\rangle$$
, where $\sum_i \beta_i^2 = 1$. (3)

then Eq. (2) does not hold. A measure of the phonon distribution is given by

$$R^{(2)} \equiv \frac{\sum_{i>1} B(E2:0_1^+ \to 2_i^+)}{\sum_{i\geq 1} B(E2:0_1^+ \to 2_i^+)}.$$
 (4)

where *i* labels the 2^+ states. From Eqs. (3) and (4), we have

$$R^{(2)} = \sum_{i>1} \beta_i^2 = 1 - \beta_1^2.$$
 (5)

Thus, $R^{(2)}$ directly measures the fraction of the Qphonon strength to 2_i^+ levels *other* than 2_1^+ and indicates the 1-phonon purity of the 2_1^+ level. Of course, that collectivity may ultimately have origins in higher states such as the giant quadrupole resonance (GQR), but our purpose here is to focus on the *nature* of the resultant collectivity, not its microscopic *origins*, and on the phonon interrelationships of the low-lying yrast levels. Hence, in the sums in Eq. (4), it is necessary to explicitly exclude the GQR and to focus on the low-lying collective 2^+ states. In effect, we view Eqs. (1)–(5) as referring to the $0\hbar\omega$ space, regardless of whether or not the microscopic origin of this collectivity resides in the GQR.

The same prescription for Eq. (4)' arises from a complementary source. As the aim of the present study is to confront the data with calculations of collective nuclei, we need to choose a model that can offer predictions for a full variety of collective structures in a unified framework with a minimum of parameters. In practice, the only available model is the IBA [11], in which the full panopoly of observed collective structures can be reproduced with only two parameters. (Ultimately, of course,

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a shell model understanding of phonon structure would be desirable, but practical calculations for all mass regions and structural varieties are currently not feasible.) Inasmuch as the IBA is a valence space model for low-lying collective states, it is essential that the data set used in Eq. (4) do likewise.

First, we consider the experimental situation. We have evaluated empirical $R^{(2)}$ values for all collective nuclei [nuclei with $R_{4/2} = E(4_1^+)/E(2_1^+) > 1.9$] from Zn to Fm, where absolute B(E2) values to more than one 2^+ state are known. All known $B(E2:0_1^+ \rightarrow 2_i^+)$ values to discrete 2^+ states were included. The results for 101 nuclei are summarized in Fig. 1. Though it is, of course, commonly recognized [12] that the B(E2) from the ground state to the 2_1^+ level is usually dominant, the quantitative degree of this dominance has not been discussed and is striking indeed. $R^{(2)}$ is always small: it is <0.09 in 96% of these nuclei (moreover three of the four exceptions have large errors). $R^{(2)}$ is <0.06 in 85% of the cases and its average value is about 0.03: Note that this overwhelming (97%) dominance and phonon purity is independent of the detailed structure of these nuclei, which comprise all varieties of structures including near harmonic and anharmonic vibrators, γ -soft nuclei, transitional nuclei, and rotors. The data in Fig. 1 also show interesting features in the distribution of $R^{(2)}$ values, in particular, the peak near 8%-9%. This is not a statistical fluctuation but reflects a specific group of nuclei (Os) and type of structure: To see this, we show the $R^{(2)}$ values for Hf-Pt in the inset to Fig. 1.

We now turn to the IBA. The 2_1^+ (and 4_1^+) states are pure phonon excitations in all three symmetries. Though long known for U(5), the phonon description of O(6) was only recently given [8]. $R^{(2)}$ vanishes for all three dynamical symmetries U(5), O(6), and SU(3), because the selection rules ($\Delta n_d = 1$), ($\Delta \tau = 1, \Delta \sigma = 0$), and



FIG. 1. Histogram of experimental $R^{(2)}$ values for 101 nuclei from Zn to Fm obtained from all known $B(E2:0_1^+ \rightarrow 2_i^+)$ values tabulated in the Nuclear Data Sheets. The inset shows detailed $R^{(2)}$ values for Hf-Pt nuclei plotted against $E(4_1^+)/E(2_1^+)$.

 $[\Delta(\lambda, \mu) = 0]$ forbid all E2 transitions from the ground state to any state other than 2_1^+ . Outside the symmetries we study the behavior of $R^{(2)}$ by systematically exploring the symmetry triangle (see Fig. 2) with the consistent Qformalism (CQF) [13], by writing

$$H = -\kappa Q(\chi) \bullet Q(\chi) + \varepsilon n_d, \qquad (6)$$

where

$$Q_{2\mu}(\chi) = s^{\dagger} d_{\mu} + d_{\mu}^{\dagger} s + \chi (d^{\dagger} d)_{\mu}^{(2)}.$$
 (7)

The E2 operator is $T(E2) = e_B Q(\chi)$. [The effective charge e_B cancels out in Eq. (4).] With no loss of generality, we take κ constant throughout at -0.025 MeV.

We first study the U(5) \rightarrow O(6) transition leg which is the easiest to understand. In both limits we take $\chi = 0$ and N = 6. The transition region is traversed by varying ε (or, equivalently, ε/κ) from infinity [U(5)] to 0 [O(6)]. The results for $R^{(2)}$ are shown in Fig. 2(a). The most striking feature is the extreme smallness of $R^{(2)}$



FIG. 2. Top right, the IBA symmetry triangle. The other panels give calculated values of $R^{(2)}$ [or $R^{(4)}$ for panels (f) and (g)] giving the absolute magnitude of phonon mixing in the 2_1^+ and 4_1^+ states in the IBA for various trajectories along the symmetry triangle. (See text for details.)

everywhere along the U(5)-O(6) transition: $R_{\text{max}}^{(2)}$ is only 0.002.

This result may be understood by expanding the IBA wave functions in an O(6) basis instead of the usual U(5) basis. The O(5) symmetry is common to both U(5) and O(6), and hence [14,15] τ is a good quantum number throughout the transition from $U(5) \rightarrow O(6)$, and the $\Delta \tau = \pm 1$ selection rule for the E2 operator also persists. Hence also the mixing of the 0^+ and 2^+ O(6) basis states is restricted to states with the same τ and different σ . Except very close to U(5), then, the low-lying physical states are composed mostly of two O(6) components, with $\sigma = N$ and N - 2 [see Fig. 3(a)]. Contributions from O(6) basis states with $\sigma < N - 4$ are only significant close to U(5). From the wave functions, the contribution of various E2 matrix elements to the physical $B(E2:0_1^+ \rightarrow 2_i^+)$ values can be quantitatively determined. There are only two significant nonvanishing contributions to $B(E2:0_1^+ \rightarrow 2_i^+)$ values. For the $B(E2:0_1^+ \rightarrow 2_1^+)$ value, the two O(6) amplitudes $0^+(\sigma = N, \tau = 0) \rightarrow 2^+(\sigma = N, \tau = 1)$ and $0^+(\sigma = N - 2, \tau = 0) \rightarrow 2^+(\sigma = N - 2, \tau = 1)$ add coherently while for the other [usually the $B(E2:0_1^+ \rightarrow 2_3^+)$] value], by orthogonality, they nearly cancel. [The component $0^+(\sigma = N - 4, \tau = 0) \rightarrow 2^+(\sigma = N - 4, \tau = 1)$ further enhances the cancellation as U(5) is approached.] Hence, the $B(E2:0_1^+ \rightarrow 2_1^+)$ nearly exhausts the strength.

For $O(6) \rightarrow SU(3)$, $\varepsilon = 0$ and χ varies from 0 to -1.32. Calculations were performed for N = 6 and with N varying linearly with χ from 4 to 16. The results in Figs. 2(b) and 2(c) show that $R^{(2)}$ is again very small, with a maximum of ~0.05. A phonon description is thus again preserved to a very high degree. The U(5) \rightarrow SU(3) transitional region calculations were carried out by decreasing ε/κ as χ varies in the range $0 \rightarrow -1.32$. As Figs. 2(d) and 2(e) show, $R^{(2)} < 0.05$.

These results can also be easily understood. It is again useful to use a different basis: We expand the wave functions in SU(3) [see Fig. 3(b)]. The primary mixing matrix elements, in this basis, have $\Delta K = 0$ [15]. Thus



FIG. 3. Partial level schemes of O(6) and SU(3) showing the most relevant basis states.

the principal mixing for low-lying states is between the SU(3) K = 0 ground band [(2N, 0) representation] and the K = 0 (" β ") band of the [2N - 4, 2] representation. However, the mixing of 0^+ and of 2^+ states is comparable, as are the intraband SU(3) E2 matrix elements, and, hence, by orthogonality, the $B(E2:0_1^+ \rightarrow 2_1^+)$ value is little affected by the mixing, and the $B(E2:0^+_1 \rightarrow 2^+_\beta)$ value nearly vanishes by cancellation. However, there is also weak $\Delta K = 2$ mixing of the SU(3) 2^+_{γ} state (quasi- γ bandhead) with the 2^+ states of both the ground and the [2N - 4, 2] K = 0 excitations. Since there is no " 0_{ν}^{+} " level available to provide cancellation, both of these admixed 2⁺ states contribute (the latter via the intermediary of the aforementioned $\Delta K = 0$ mixing) to yield a small $B(E2:0_1^+ \rightarrow 2_{\gamma}^+)$ value which is the main contributor to the numerator of $R^{(2)}$.

The results in Fig. 2 encourage us to pursue the validity of this Q-phonon picture in a more general context that comprises all the manifestations of collectivity in the IBA. We have therefore carried out a thorough set of calculations with a rather fine mesh of internal paths in the symmetry triangle. The results for N = 6 are shown in the contour plot in Fig. 4. (Similar results are obtained for other N values.) Remarkably, $R^{(2)}$ is always small. Its maximum values are along the $O(6) \rightarrow SU(3)$ transition leg and, for any $N \leq 16$, the maximum value, independent of the IBA parameters (i.e., for any $\kappa, \varepsilon, \chi$ in the symmetry triangle), is $R^{(2)} < 0.07$. The average value is ~ 0.02 . As we saw in Fig. 1, the experimental values are equally small, averaging ~ 0.03 and nearly always <0.09. Thus, the IBA predicts that the 2^+_1 level is a nearly pure one-quadrupole-phonon excitation of the ground state induced by the $0\hbar\omega$ operator Q. Note that this prediction is inherent to the model. Were the data different, the model predictions could not be altered accordingly.

The peak in $R^{(2)}$ in Figs. 2 and 3 between O(6) and SU(3) is also interesting since it mirrors the empirical $R^{(2)}$ distribution in the inset in Fig. 1. The maximum $R^{(2)}$ values occur for precisely the nuclei ¹⁸⁶⁻¹⁹²Os that are the archetypical representatives of the O(6) \rightarrow rotor transition [16]. Thus, not only does the IBA reproduce the



FIG. 4. Contour plot of $R^{(2)}$ throughout the symmetry triangle. N = 6.

empirical fact of very small $R^{(2)}$ values, but it reproduces the detailed distribution of $R^{(2)}$ values across transition regions, as exemplified in Fig. 1 (inset).

It is useful to put these results in a more general context. The usual fermion E2 operator consists of proton and neutron parts, with separate effective charges. This feature is conveyed to the boson situation in the IBM-2 (Ref. [17]), where

$$T(E2) = e_{p}^{B}Q_{p}^{B} + e_{n}^{B}Q_{n}^{B}.$$
 (8)

In low-lying states, the proton and neutron motions are coherent, and such states are connected to the ground state predominantly by the isoscalar component of T. There is an overwhelming concentration of strength due to this operator from the ground state, leading to the Q-phonon picture. In this sense, the Q phonon is largely of isoscalar nature (F scalar for bosons), and the quadrupole-quadrupole interaction naturally favors this excitation at lower energy. At higher energies, there should be states excited by the isovector E2 operator. In the IBM-2, these are mixed-symmetry states but are excited very weakly by E2 transitions [18] [with $B(E2:0_1^+ \rightarrow 2) \sim 10^{-1}-10^{-2}$ of $B(E2:0_1^+ \rightarrow 2_1^+)$].

The above IBA analysis can be extended to the 4_1^+ state. As for the 2_1^+ state, we ask the extent to which $|4_1^+4\rangle = NQ_{22}|2_1^+2\rangle$: that is, we write

$$NQ_{22}|2_1^+2\rangle = \alpha_1|4_1^+4\rangle + \sum_{i>1}\alpha_i|4_i^+4\rangle.$$
(9)

We define $R^{(4)}$ analogously to Eq. (4) in terms of $B(E2: 2_1^+ \rightarrow 4_i^+)$ values. Examples of the results are shown in Figs. 2(f) and 2(g): $R^{(4)}$ is again very small. The maximum value for any N, κ, ε , or χ is only 0.01, which is nearly an order of magnitude less than $R^{(2)}_{max}$. The reason is simple and interesting. As with $R^{(2)}$, the largest $R^{(4)}$ values occur between O(6) and SU(3) and mixing arguments given for $R^{(2)}$ apply here *except* that even the $\gamma \rightarrow g E2$ contribution to $R^{(4)}$ is now largely cancelled since there are *both* 2_{γ}^+ and 4_{γ}^+ levels to mix with 2_g^+ and 4_g^+ . Note that the high purity of the 4⁺ state holds for the $Q_{22}|2_1^+2\rangle$ configuration: $Q_{22}Q_{22}|0\rangle$ is less pure.

The above discussions highlight an important point. The small values of $R^{(2)}$ and $R^{(4)}$ are not somehow a trivial result but rather stem from basic properties of the symmetry limits, from the characteristic ways in which these symmetries are broken, and from subtle but pervasive and strong cancellation mechanisms.

To conclude, we have shown quantitatively that the empirical 2_1^+ states of all collective even nuclei from Z = 30 to 100 nearly exhaust the known $0\hbar\omega E2$ strength from the ground state. Strength to all other known 2^+ states averages only about 3% and exceeds a 10% fraction in

only 1%-2% of nuclei (and most of these cases have large uncertainties). A thorough mesh of IBA calculations, comprising essentially all types of collective structures, shows that the model naturally and unavoidably produces exactly the empirical result: $R^{(2)}$ is always less than 0.07 independent of the CQF-IBA Hamiltonian parameters. The agreement of calculated with empirical $R^{(2)}$ values is not limited to global averages. For example, the large $R^{(2)}$ values between Hf and Pt (Fig. 1, inset) are reproduced by the IBA for nuclei spanning an O(6) \rightarrow rotor region [Fig. 2(c), Fig. 4]. Finally, an even higher Q-phonon purity is calculated for the 4_1^+ state as a single Q-phonon excitation of the 2_1^+ state.

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