

Fine structure of the isoscalar giant quadrupole resonance in ^{28}Si and ^{27}Al

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The Isoscalar Giant Quadrupole Resonance (ISGQR) in ^{27}Al and ^{28}Si has been investigated with high energy-resolution proton inelastic scattering at $E_p = 200$ MeV and at scattering angles close to the maximum of $\Delta L = 2$ angular distributions with the K600 magnetic spectrometer of iThemba LABS, South Africa. Characteristic scales are extracted from the observed fine structure with a wavelet analysis and compared for ^{28}Si with Random Phase Approximation (RPA) and Second-RPA (SRPA) calculations with an interaction derived from the Argonne V18 potential by a unitary transformation. A recent extension of the method to deformed nuclei provides the best description of the data suggesting the significance of Landau damping.

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I. INTRODUCTION

High energy-resolution inelastic proton scattering data obtained at beam energies of a few hundred MeV with magnetic spectrometers in combination with dispersion matching techniques provide a suitable approach to study fine structure of the ISGQR in nuclei. The necessary experimental techniques to perform (p, p') experiments with an energy resolution better than the energy spread of the incident projectile beam were discussed in Ref. [1] and for the experiments discussed here using the K600 spectrometer at iThemba LABS in Ref. [2].

For a quantitative analysis of the observed fine structure wavelet techniques have been shown to be most useful [3]. A systematic study of the ISGQR in medium-mass to heavy nuclei [4, 5] showed that scales characterizing the fine structure originate from a collective damping mechanism induced by the coupling of elementary one-particle one-hole ($1p1h$) states to low-lying surface vibrations [6]. In contrast, Landau damping (i.e. fragmentation on the $1p1h$ level) significantly contributes to the fine structure of the isovector giant dipole resonance [7]. Such conclusions were based on the comparison of RPA calculations with more refined approaches allowing for a coupling of the $1p1h$ excitations to two particle-two hole ($2p2h$) states.

Recently, a high energy-resolution measurement of the Isoscalar Giant Quadrupole Resonance (ISGQR) in ^{40}Ca was performed [8] motivated by the question whether

collective damping remains the most important decay mechanism in lighter nuclei or if other mechanisms might play a leading role [9]. From a comparison of RPA and second-RPA (SRPA) calculations with a realistic interaction derived from the Unitary Correlation Operator Method (UCOM), it was concluded that Landau damping produces all but the lowest scale derived from the wavelet analysis [8]. The importance of Landau damping obtained with the UCOM interaction is in contrast to many RPA results for the ISGQR in ^{40}Ca with phenomenological interactions (see references in [8]) which all predict a single collective state and correspondingly no fine structure.

The present work extends these investigations to an even lighter nucleus, ^{28}Si . The fine structure properties are compared with those of RPA and SRPA calculations in the framework discussed in Ref. [8]. Since ^{28}Si is an open-shell nucleus in the middle of the sd shell with a deformed ground state (g.s.), the SRPA approach was modified to define a g.s. wave function using shell-model occupation numbers for the valence orbits. Alternatively, a new approach based on a deformed Hartree-Fock g.s. [10] is tested. The impact of coupling a single particle (or hole) to an even-even core on the fine structure of the ISGQR is studied by an experimental comparison with the isotonic nucleus ^{27}Al . In the sequel we discuss experiments, model calculations and extractions of scales always first and in greater detail for ^{28}Si and then for ^{27}Al .

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II. EXPERIMENTAL RESULTS

A. Experiment and spectra

The experiments were carried out with a 200 MeV proton beam produced by the Separated Sector Cyclotron (SSC) of the iThemba Laboratory for Accelerator Based Sciences (iThemba LABS), South Africa. The protons were scattered inelastically from self-supporting targets of $^{\text{nat}}\text{Si}$ (92.2% ^{28}Si) and ^{27}Al with areal densities of 0.232 mg/cm² and 0.819 mg/cm², respectively, and then momentum analysed with the K600 magnetic spectrometer. Dispersion matching techniques were used in order to exploit the high energy-resolution capability of the spectrometer. A scattering angle $\theta_{\text{Lab}} = 12^\circ$ close to the maximum of the cross section for $\Delta L = 2$ transitions populating the ISGQR was selected. Details of the data analysis procedures are described in Ref. [11].

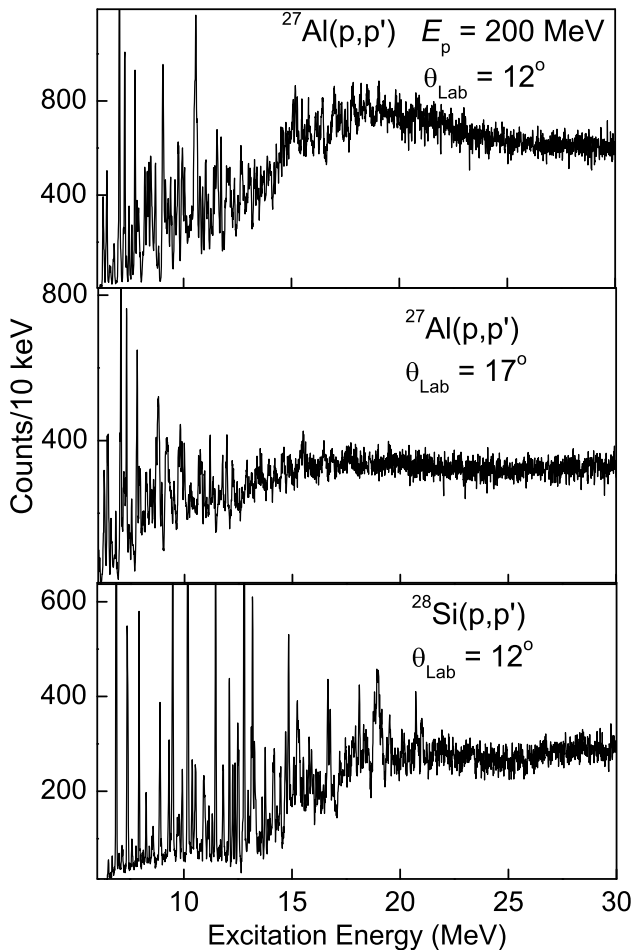


FIG. 1: Excitation energy spectra for $^{27}\text{Al}(p,p')$ reactions between 6 and 30 MeV at $E_p = 200$ MeV and $\theta_{\text{Lab}} = 12^\circ$ (top), $\theta_{\text{Lab}} = 17^\circ$ (middle), and the $^{28}\text{Si}(p,p')$ reaction at $\theta_{\text{Lab}} = 12^\circ$ (bottom) covering low-lying states and the energy region of the ISGQR.

The excitation energy spectrum for the nucleus ^{28}Si is

shown in the bottom part of Fig. 1 for the full energy-bite of the K600 magnetic spectrometer, $6 \leq E_x \leq 30$ MeV. An energy resolution $\Delta E = 38$ keV (Full Width at Half Maximum, FWHM) was achieved. A broad peak about 5 MeV wide between about 14 and 25 MeV centered at $E_x \approx 18$ MeV resulting from excitation of the ISGQR can be identified. This is consistent with features of the ISGQR in ^{28}Si observed in (α, α') scattering [12]. Many discrete states below the respective proton (11.584 MeV) and neutron (17.178 MeV) thresholds in ^{28}Si are resolved, and above these thresholds, fine structure is clearly visible in the energy region of the ISGQR. For ^{27}Al at the same scattering angle (upper part of Fig. 1) a very similar behavior is found. In the middle part of Fig. 1, a spectrum of ^{27}Al measured at $\theta_{\text{Lab}} = 17^\circ$ is displayed. Here, the ISGQR bump is significantly reduced with respect to the background originating from other multipoles and quasi-free reactions. However, in analogy to the case of ^{40}Ca a cross correlation analysis [8] of the two spectra demonstrates that the visible fine structure originates from the ISGQR.

B. Comparison of different probes exciting the ISGQR in ^{28}Si

For sd -shell nuclei with $A \leq 32$, extensive experimental investigations have been made on the fragmentation of the ISGQR. Results from proton (present work), α [12] and electron scattering [14] for the case of ^{28}Si are compared in Fig. 2.

Van der Borg *et al.* [12] investigated the properties of the ISGQR for a variety of sd -shell nuclei using (α, α') scattering at $E_\alpha = 120$ MeV with a resolution $\Delta E = 125$ keV (FWHM). The resulting $^{28}\text{Si}(\alpha, \alpha')$ excitation energy spectrum at $\theta_{\text{Lab}} = 6^\circ$ is shown in the middle panel of Fig. 2. At this angle, many of the states excited in the energy range 14 – 22 MeV were identified to have spin-parity $J^\pi = 2^+$ from angular distribution measurements and excitation of the ISGQR was found to dominate over other isoscalar modes ($E0, E3$). There is good correspondence of the main structures except for an overall shift of the excitation energies of a few hundred keV.

The $E2$ strength distribution in ^{28}Si has also been measured with α scattering by the Texas A&M group and overall fair agreement is found [13]. These data are not included in the further analysis because spectra for kinematics where $E2$ cross sections dominate were not available and the overall energy resolution was 250 – 300 keV (FWHM) only.

Additional comparison is made with $^{28}\text{Si}(e, e')$ data at low momentum transfers [14]. The lower panel in Fig. 2 displays the $B(E2)$ strength distribution extracted from a form factor decomposition. The structures at 18 – 19 MeV and around 21 MeV agree with the peaks in the α and proton scattering data. An additional pronounced structure is visible around 20 MeV absent in the hadron scattering data. In contrast to the dominance of isoscalar

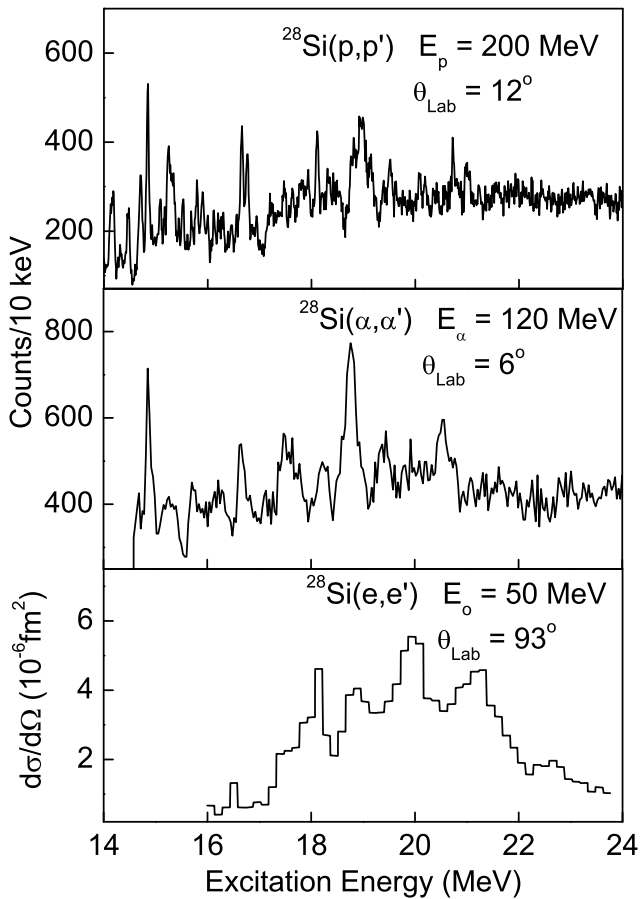


FIG. 2: Excitation energy spectra for the ISGQR in ^{28}Si from different probes: $^{28}\text{Si}(p,p')$ (present work), $^{28}\text{Si}(\alpha,\alpha')$ [12], and $^{28}\text{Si}(e,e')$ [14].

excitations in the proton and α scattering, electron scattering also excites isovector transitions indicating isospin $T = 1$ for the structure at 20 MeV.

III. MODEL CALCULATIONS

Isoscalar $E2$ strength distributions were calculated using the Random Phase Approximation (RPA) and the SRPA formalism [15, 16] with an effective interaction derived from the realistic Argonne V18 nucleon-nucleon interaction in the framework of the UCOM, as described in [17]. The same model was used in [8] to analyse the fine structure of the ISGQR in ^{40}Ca . A single-particle space of 13 harmonic-oscillator shells was used here to obtain the Hartree-Fock (HF) reference state and all HF single-particle states are included in the $1p1h$ and $2p2h$ spaces. The RPA and SRPA matrix elements are renormalized as described in [16] with occupation numbers taken from a $4\hbar\omega$ shell-model calculation. This hybrid model allows for the activation of sd states as particle and hole states to take into account the open-shell character of ^{28}Si . We

further exclude low-energy $1p1h$ and $2p2h$ configurations within the sd shell in order to avoid unphysical solutions owing to the approximate description of the sd shell.

The above models do not explicitly take into account deformation. Therefore, theoretical $E2$ distributions were calculated also in the RPA using an axially-deformed HF ground state and applying angular-momentum projection techniques [10]. Both the HF and the RPA calculations use the same realistic nucleon-nucleon interaction derived from the Argonne V18 potential by a unitary transformation in the framework of the (UCOM) and the similarity renormalisation group as described in [19, 20] and are supplemented by a phenomenological three-nucleon contact interaction. This Hamiltonian was introduced and tested in Ref. [21] for ground-state observables and applied to RPA calculations in closed-shell nuclei in Ref. [22] (we use the version labeled ‘S-UCOM(SRG)’). All calculations are performed in a harmonic-oscillator single-particle basis covering 15 oscillator shells. Further details on the deformed RPA approach employed in this work can be found in Ref. [10].

The experimental spectra have a finite energy resolution while the RPA and SRPA models provide a discretized strength distribution. Therefore, for the purposes of a direct comparison between experiment and theory independent of the resolution, the calculations were smoothed with a Gaussian function of width $\Delta E = 38$ keV (FWHM) and put into a bin size of 10 keV. Figure 3 summarizes the results of the RPA and SRPA predictions for ^{28}Si . It can be seen that the ISGQR is spread out widely. The differences between RPA and SRPA (middle and lower panels) results are small. The main effect is an overall shift of about 2 MeV lower in excitation in the SRPA calculation while the coupling to $2p2h$ states produces little additional fragmentation. The RPA calculation based on a deformed HF ground state (top panel) finds a similar distribution in energy but shows significantly more structure.

IV. EXTRACTION OF CHARACTERISTIC ENERGY SCALES

A. Wavelet analysis

The main aim of the present study is to extract characteristic energy scales within the excitation energy region of the ISGQR in the experimental data and compare those scales to the ones obtained from the corresponding theoretical strength functions. In order to quantitatively investigate the observed structures, wavelet analysis techniques [3, 5, 8] were utilized. A wavelet analysis of the data was accomplished using the Matlab programme [23].

A wavelet transform is defined as a convolution of the original energy spectrum $\sigma(E)$ multiplied by a scaled, shifted version of a wavelet function Ψ . The coefficients

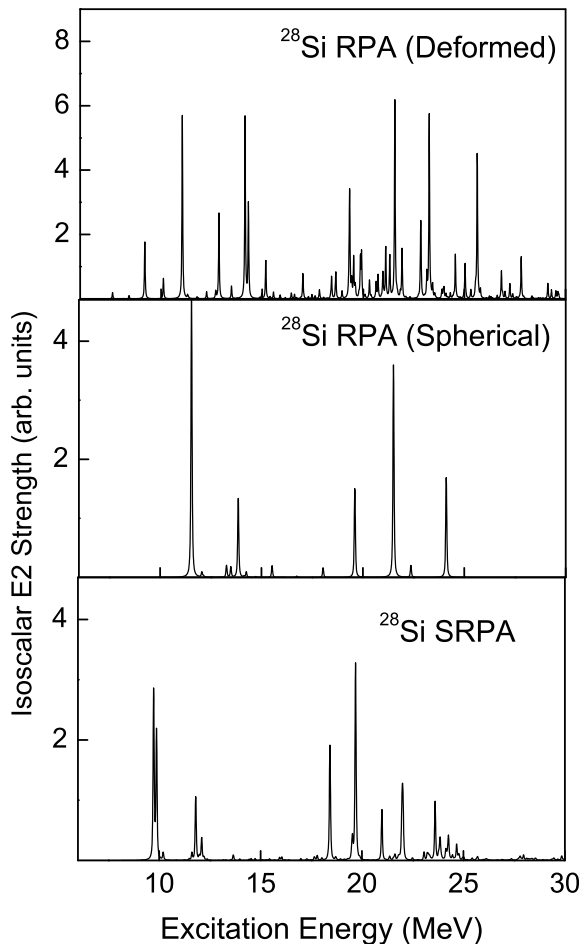


FIG. 3: RPA and SRPA isoscalar $E2$ strength distributions for ^{28}Si from the models described in the text. The discrete distributions are smoothed with the experimental energy resolution for direct comparison to the experimental data.

$(E_x, \delta E)$ of the wavelet transform are obtained as

$$C(E_x, \delta E) = \frac{1}{\sqrt{\delta E}} \int \sigma(E) \Psi\left(\frac{E_x - E}{\delta E}\right) dE, \quad (1)$$

where δE is the scale and E_x is the position along the energy scale, respectively.

The parameters, excitation energy, E_x , and scale, δE , can be varied continuously (Continuous Wavelet Transform, CWT) or in discrete steps j , where $\delta E = 2^j$, $j = 1, 2, 3, \dots$, and $E_x = \delta E$. A variety of wavelets provide a set of tools for performing many different tasks. The choice of a particular wavelet is based on its mathematical properties [23, 24] amongst which are moments, compact support and the regularity of a signal (the number of times that it is continuously differentiable). Complex wavelets produce a complex wavelet transform, allowing the phase of the result to be examined. The choice of a mother wavelet is motivated by the type of spectrum being analyzed. The Morlet mother wavelet used in this analysis is derived by taking a periodic wave and localiz-

ing it with a Gaussian envelope

$$\Psi(x) = \frac{1}{\pi^{1/4}} \exp(ikx) \exp\left(-\frac{x^2}{2}\right). \quad (2)$$

Here, the parameter k weighs the resolution in scales *versus* the resolution in localization. A value of $k = 5$ was chosen in the present work.

B. Application to ^{28}Si and ^{27}Al

CWT was applied to the excitation energy region of the ISGQR in the spectra of Fig. 1 with a Complex Morlet mother wavelet. It should be emphasised, however, that before the application input data preparation is essential in order to minimize end effects. In our case, the mean of the input data set was subtracted from the data set before applying the wavelet technique. By plotting the real part of the complex coefficients on a two-dimensional plot of scale *versus* excitation energy, the positions of the structures within the original energy spectrum can be identified as illustrated in the lower-right panel of Fig. 4 for the $^{28}\text{Si}(p, p')$ experimental data. The maximum scale was restricted to 5 MeV in order to show the detail in substructures existing at smaller energy scales. To obtain the corresponding power spectrum, the square root of sum of the squares of the complex coefficients are summed across onto the scale axis. The resulting power spectrum is shown in the lower-left panel of Fig. 4. It exhibits pronounced peaks representing characteristic energy scales of the structures in the spectrum.

It is instructive to compare the wavelet analysis results for the ISGQR in ^{28}Si from the (p, p') and (α, α') data sets (Fig. 5). The power spectra at the r.h.s. of Fig. 5 show very similar scales, i.e. maxima at a certain energy, as one would expect from the similarity of the spectra. The power maximum around 1 MeV exhibits comparable peaks but as a result of the better energy resolution the (p, p') data lead to more structure in the power spectrum at smaller scale values. The largest scales are affected by the shape of the physical background and the peak-to-background ratio leading to the observed differences in power values.

A summary of the power spectra from the CWT analysis for the present high-resolution (p, p') data in the energy region of the ISGQR in ^{28}Si and ^{27}Al and the corresponding theoretical isoscalar $E2$ strength distributions (see Fig. 3) is given in Fig. 6. The choice and suitability of wavelet functions for the analysis of nuclear giant resonances was discussed extensively in Ref. [3]. However, in the present analysis a Complex Morlet function was chosen since the resulting energy scales extracted are directly comparable to those obtained from a Fourier analysis. In previous analyses of the ISGQR in heavier nuclei [3, 4] a real Morlet wavelet was used and the extracted scales need to be divided by a factor of 0.813 in order to yield the corresponding Fourier scale. Characteristic scales deduced from the power spectra are summarized in Table. I.

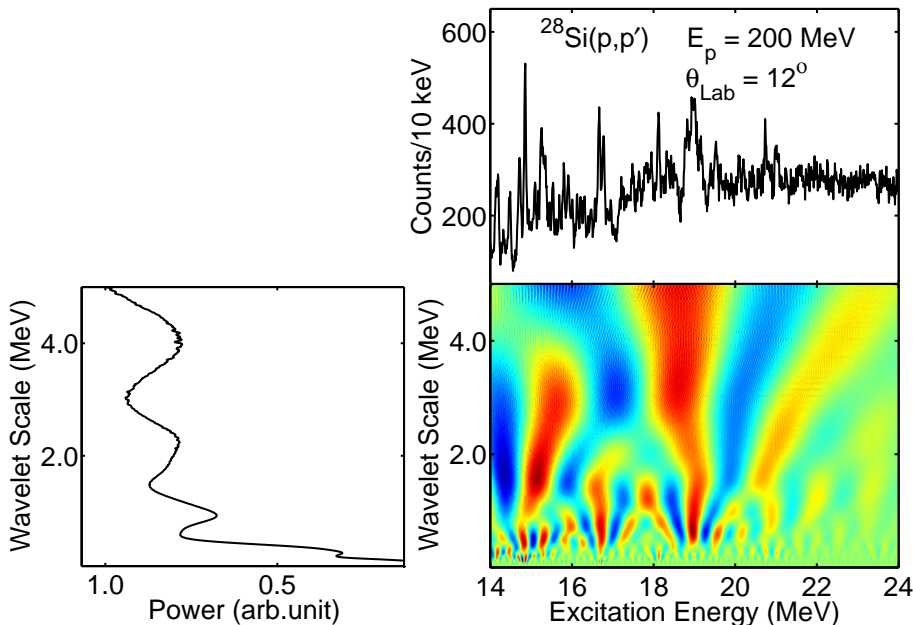


FIG. 4: Upper-right: Excitation energy spectrum of the $^{28}\text{Si}(p,p')$ reaction at a scattering angle $\theta_{\text{Lab}} = 12^\circ$ corresponding to the maximum cross-section for excitation of the ISGQR. Lower-right: CWT after applying a complex Morlet mother wavelet, red indicates large positive and blue large negative coefficients. Lower-left: Power spectrum in which the maxima of the peaks correspond to the values of characteristic energy scales.

In passing we note that an earlier performed CWT analysis of the $^{28}\text{Si}(p,p')$ and (α,α') spectra (but without the above-mentioned input data preparation) gave very similar results as in the present work [25]. Analogous to previous studies of scales of the ISGQR [4, 5, 8] the scales are grouped in three classes: < 300 keV (Class I), $300 - 1000$ keV (Class II) keV, and > 1000 keV (Class III).

One finds two broad (Class III) scales in the ^{28}Si data and in all calculations. However, a quantitative comparison is limited. In particular, the largest scale in the data is underestimated by all models. This may be related to neglect of coupling to the continuum which is known to

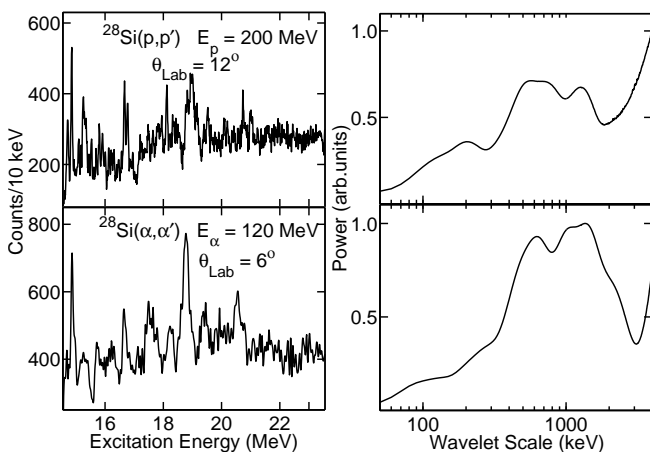


FIG. 5: Comparison of the CWT power spectra for the energy region of the ISGQR from the $^{28}\text{Si}(p,p')$ and (α,α') spectra.

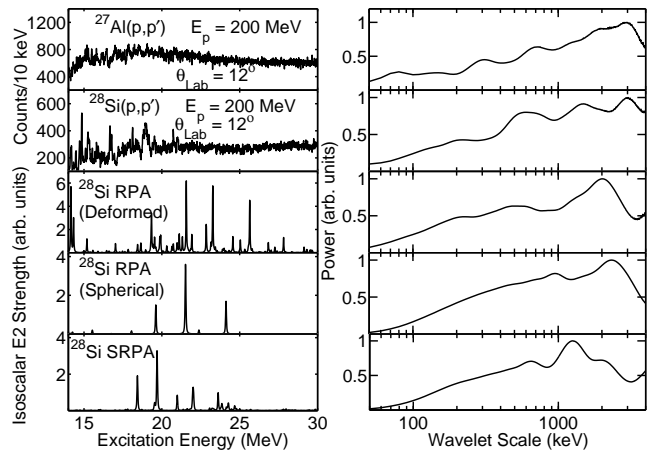


FIG. 6: L.h.s.: Excitation energy spectra of ^{27}Al and ^{28}Si at the scattering angle $\theta_{\text{Lab}} = 12^\circ$ corresponding to the maximum cross sections for excitation of the ISGQR and theoretical strength distributions from RPA and SRPA. R.h.s.: Corresponding power spectra from the CWT.

make significant contributions to the resonance width in light nuclei. One intermediate (Class II) scale is observed in the data as well as in the RPA calculations on a deformed g.s.. In contrast, the spherical RPA results show no Class II scales, but the SRPA predicts one scale in fair agreement with the experimental value. None of the three nuclear models is capable to reproduce both Class I scales in ^{28}Si , but the deformed RPA calculations account for the larger one indicating that scales between 200 and

TABLE I: Summary of energy scales of the ISGQR in ^{27}Al and ^{28}Si and comparison with RPA and SRPA calculations for ^{28}Si .

		Class I		Class II		Class III	
		< 300 keV		300 – 1000 keV		> 1000 keV	
^{27}Al	Exp.	80	140	310	700	2000	2800
^{28}Si	Exp.	130	220		600	1500	3000
^{28}Si	RPA def.		250		500	1250	2000
^{28}Si	RPA sph.					1000	2400
^{28}Si	SRPA				650	1250	2000

1000 keV result dominantly from Landau fragmentation.

Similar to the case of ^{40}Ca [8], no interpretation based on the model calculations can be given for the smallest scale in the data at 130 keV. Corresponding scales around 100 keV were found in all analyses of the ISGQR in heavier nuclei, but these could be clearly related to the coupling to low-energy phonons [4, 5]. Here, the SRPA is performed in the diagonal approximation (i.e. the coupling among $2p2h$ states is ignored), leading to many near-degenerate eigenstates. Inclusion of the $2p2h$ - $2p2h$ interaction pushes the spacing distribution towards the Wigner limit by level repulsion with a maximum average spacing at larger distances [26], which may produce a characteristic scale in the wavelet analysis. We have tested the idea with calculations in a limited energy range and model space for the case of ^{40}Ca and ^{28}Si , and found it to be qualitatively consistent with our speculation.

Another possible interpretation of the lowest scale are Ericson fluctuations [27], i.e., the scale represents the coherence width of overlapping 2^+ resonances. The coherence width in ^{28}Si has been discussed e.g. in Ref. [28] and amounts to about 60 keV in the excitation region of the ISGQR. While the experiments quoted in Ref. [28] usually cover a larger spin window, the selectivity of the present data on $J^\pi = 2^+$ states leads one to expect a larger value for the coherence width which could well be of the order of the experimentally observed scale.

Finally, we briefly discuss a comparison of the scales in ^{28}Si with those in the odd-mass isotone ^{27}Al . All scales found in ^{28}Si appear in ^{27}Al as well. Typically, the values in the latter are slightly larger except for the broadest scale. Thus, the coupling of the unpaired proton seems to have little influence on the fine structure. However, an additional scale at 80 keV appears in ^{27}Al which may

be interpreted to arise from the coupling of the unpaired proton (hole) to an even-even core. Therefore, we have observed that there is little influence of coupling a single particle (or hole) to an even-even core on the fine structure of the ISGQR which is yet to be proven by theoretical calculations.

V. SUMMARY

The ISGQR excited in the nuclei ^{27}Al and ^{28}Si has been investigated using inelastic proton scattering at 200 MeV. Characteristic energy scales have been extracted by applying wavelet analysis techniques. In order to reveal the physical nature of observed scales, a comparison to results from RPA and SRPA calculations has been made for the case of ^{28}Si where the open-shell ground state has been approximated by shell-model occupation numbers for levels near the Fermi surface. Alternatively, RPA calculations were performed based on a deformed HF ground state. Clearly, the RPA calculations in a deformed basis provide a more realistic description of the experimentally observed fragmentation of the ISGQR. They qualitatively account for all but the smallest experimentally observed scale and indicate Landau fragmentation driven by deformation as the most important mechanism of the fine structure. As for ^{40}Ca , the smallest experimental scale at 130 keV in ^{28}Si cannot be explained by any of the models. It may be related to the neglect of coupling among the $2p2h$ states in the SRPA calculations or it may result from Ericson fluctuations. A comparison of the ISGQR in the neighbouring odd-mass isotone ^{27}Al reveals little difference except for an additional scale at 80 keV which may be interpreted to arise from the coupling of the unpaired proton.

Acknowledgements

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