Experimentally constrained $(p,\gamma)^{89}\text{Y}$ and $(n,\gamma)^{89}\text{Y}$ reaction rates relevant to p -process nucleosynthesis

A. C. Larsen,^{1,*} M. Guttormsen,^{1,†} R. Schwengner,^{2,‡} D. L. Bleuel,³ S. Goriely,⁴ S. Harissopulos,⁵ F. L. Bello Garrote,¹ Y. Byun,⁶ T. K. Eriksen,¹ F. Giacoppo,¹ A. Görgen,¹ T. W. Hagen,¹ M. Klintefjord,¹ T. Renstrøm,¹ S. J. Rose,¹ E. Sahin,¹ S. Siem,¹ T. G. Tornyi,¹ G. M. Tveten,¹ A. V. Voinov,⁶ and M. Wiedeking⁷

¹Department of Physics, University of Oslo, N-0316 Oslo, Norway

²Helmholtz-Zentrum Dresden-Rossendorf, 01328 Dresden, Germany

³Lawrence Livermore National Laboratory, Livermore, California 94551, USA

⁴Institut d'Astronomie et d'Astrophysique, Université Libre de Bruxelles, Brussels, Belgium

⁵Institute of Nuclear and Particle Physics, NCSR "Demokritos", Athens, Greece

⁶Department of Physics and Astronomy, Ohio University, Athens, Ohio 45701, USA

⁷iThemba LABS, P.O. Box 722, 7129 Somerset West, South Africa

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The nuclear level density and the γ -ray strength function have been extracted for ^{89}Y by using the Oslo method on $^{89}\text{Y}(p,p'\gamma)^{89}\text{Y}$ coincidence data. The γ -ray strength function displays a low-energy enhancement consistent with previous observations in this mass region ($^{93-98}\text{Mo}$). Shell-model calculations support the conclusion that the observed enhancement is due to strong, low-energy $M1$ transitions at high excitation energies. The data were further used as input for calculations of the $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ and $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$ cross sections with the TALYS reaction code. Comparison with cross-section data, where available, as well as with values from the BRUSLIB library, shows a satisfying agreement.

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I. INTRODUCTION

The quest for a detailed understanding of heavy-element nucleosynthesis is very intriguing, calling for extensive knowledge of nuclear properties at the relevant astrophysical energies. The astrophysical aspects are equally challenging; the most critical issue is to identify the correct astrophysical sites and conditions in which the heavy-element nucleosynthesis operates [1–5].

There are three main processes responsible for creating elements heavier than iron [6]: the slow-neutron-capture (s) process, the rapid-neutron-capture (r) process, and the p process, which in fact includes a variety of nuclear reactions such as photodisintegration of already-created s and r nuclides through, e.g., (γ,p) , (γ,n) , and (γ,α) reactions as well as their inverse capture reactions. To this end, only the s -process astrophysical sites are more or less clearly identified (low-mass asymptotic giant branch stars, $M \lesssim 8M_{\odot}$, and massive stars, $M > 8M_{\odot}$; see, e.g., Ref. [3]). The r -process astrophysical site is still a mystery; suggestions include the neutrino-driven wind following a core-collapse supernova, and neutron-star mergers [1,7]. Also, the p -process site remains rather elusive [2,4,8]; however, it is clear that conditions with temperatures reaching between 2 to 3 billion degrees must be reached for the p process to take place. The deep O-Ne layers of a massive star prior to or during its supernova phase remains the most popular suggestion, although some species, in particular the p isotopes of Mo and Ru, remain underproduced in the

present simulations. Type Ia supernovae¹ could also contribute to the production of p nuclei, but its contribution remains affected by uncertainties in galactic chemical evolution models as well as in the determination of the s -process enrichment prior to the p -process nucleosynthesis [2,8].

In addition to the uncertainties in the astrophysical modeling, significant uncertainties in the nuclear reaction rates bring further complications to our understanding of the nucleosynthesis. In order to estimate theoretically the abundance distribution, very large reaction-network calculations are required, typically including ~ 2000 nuclei and $\sim 20\,000$ cross sections for the p process, and ~ 5000 nuclei and $\sim 50\,000$ cross sections for the r process. Obviously, most of these cross sections have not been determined experimentally and will not be measured in the near future. Hence, the calculations rely on theoretical estimates of cross sections and reaction rates. These are usually derived within the Hauser-Feshbach theory [9], where the main input parameters include optical-model potentials, nuclear level densities, and γ -ray strength functions (see, e.g., Ref. [1,2] and references therein).

In this work, we investigated the yttrium isotope ^{89}Y by means of the reaction $^{89}\text{Y}(p,p'\gamma)^{89}\text{Y}$, utilizing the Oslo method [10] to extract the nuclear level density (NLD) and γ -ray strength function (γSF). The NLD data were previously reported in Ref. [11], and this work focuses on the γSF , and in particular the enhancement at low-energy γ rays.

*a.c.larsen@fys.uio.no

†magne.guttormsen@fys.uio.no

‡r.schwengner@hzdr.de

¹A type Ia supernova is spectroscopically defined by the absence of hydrogen emission lines and is believed to occur in binary systems, of which one of the stars is a white dwarf gradually accreting mass from its companion until a thermonuclear explosion takes place.

Extensive shell-model calculations are performed for the $M1$ strength, clearly showing a strong increase at low γ -ray energies, in accordance with previous experimental [12,13] and theoretical [14] findings in this mass region. We also investigate the impact of our results on capture cross sections and astrophysical reaction rates. More specifically, we consider the cases $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ and $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$, because these are relevant for the p process in this mass region and are also of interest for reaction networks in the context of stockpile stewardship [15].

The paper is organized as follows: In Sec. II, experimental details and an overview of the data analysis are given. Shell-model calculations and models for the γ -ray strength function are presented in Sec. III. In Sec. IV, calculated cross sections and reaction rates are shown and compared with existing data. Finally, a summary and outlook can be found in Sec. V.

II. EXPERIMENTAL DETAILS AND EXTRACTION OF LEVEL DENSITY AND γ -STRENGTH FUNCTION

A. Experimental details and unfolding of γ spectra

The experiment was performed at the Oslo Cyclotron Laboratory (OCL), utilizing a proton beam of 17 MeV. The beam impinged on a natural ^{89}Y target with thickness 2.25 mg/cm^2 . The beam current was typically $\approx 0.5\text{ nA}$, with about five days of beam time, including calibration runs on a natural Si target.

Charged ejectiles were measured with the silicon ring (SiRi) array [16], which is a ΔE - E telescope system of eight individual telescopes, each consisting of a $130\text{ }\mu\text{m}$ eightfold segmented front detector (ΔE) with a $1500\text{ }\mu\text{m}$ back detector (E) detector; thus, there are 64 individual silicon telescopes in the system. To reduce the amount of δ electrons from the target, a $10.5\text{-}\mu\text{m}$ -thick aluminium foil was placed in front of SiRi. The SiRi system was mounted in backward angles, covering the range $\theta = 126^\circ\text{--}140^\circ$ with an angular resolution of $\Delta\theta = 2^\circ$. Typical particle spectra from the experiment are shown in Fig. 1. It is seen that the charged-particle species are

clearly separated from each other. In the following, we gate on the protons, i.e., we consider the $^{89}\text{Y}(p,p'\gamma)^{89}\text{Y}$ reaction channel. The proton-energy resolution was $\approx 130\text{--}160\text{ keV}$ (full width at half maximum, or FWHM), determined from the ground-state peak and discrete peaks in the proton $\Delta E + E$ spectrum.

The γ rays were detected with the high-efficiency NaI(Tl) array CACTUS [17]. For this experiment, CACTUS comprised 26 cylindrical NaI detectors of diameter 12.7 cm and length 12.7 cm mounted on the spherical CACTUS frame at angles $\Theta = 37.4^\circ, 63.4^\circ, 79.3^\circ, 100.7^\circ, 116.6^\circ$, and 142.6° with respect to the beam direction. All NaI crystals are collimated with lead cones to reduce the Compton contribution and enhance the peak-to-total ratio. The total efficiency for $E_\gamma = 1.33\text{ MeV}$ was $\approx 14.1\%$. The trigger for the analog-to-digital converters of the CACTUS array was the signal from the thick E detector in SiRi; this was also the start signal for the time-to-digital converters of the NaI detectors.

By selecting the proton channel and requiring the coincident γ rays with the protons within a time window of $\approx 20\text{ ns}$, we obtain γ spectra for each excitation energy, which is found from the measured proton energy in SiRi corrected for the reaction kinematics. The excitation-energy versus NaI signals are shown in Fig. 2(a). Some pileup events are observed for $E_\gamma > E_x$; however, they are rare and do not contribute significantly. We also see contaminant γ lines from ^{12}C and ^{16}O , e.g., at $E_\gamma = 4.4, 6.1$, and 7.1 MeV . These lines are kept in the data set for unfolding, and later removed as explained in the following.

The γ -ray spectra for each excitation energy bin were corrected for the CACTUS response functions, i.e., removing the Compton, single-escape, double-escape, and backscattered annihilation events and correcting for the full-energy efficiency. The contamination lines were also unfolded to obtain only the full-energy peaks in the final, unfolded γ -ray spectra. Then, a three-dimensional scan of the area around the full-energy peaks was performed to get a correct estimate of the underlying spectrum, and finally removing the carbon

FIG. 1. ΔE - E plot for $^{89}\text{Y}+p$ (a), with the deposited energy in the thick E detector versus the thin ΔE detector for $\theta = 132^\circ \pm 1^\circ$, and (b) the sum of the deposited energy in the thin and thick detector for the same angle. The labeled peaks are the ground state (0), the first excited 909 keV level (1), the second excited 1507 keV level (2), and the third excited 1744 keV level (3).

FIG. 2. (a) Raw NaI spectra and (b) unfolded, primary γ spectra for each initial excitation-energy bin for $^{89}\text{Y}(p,p'\gamma)^{89}\text{Y}$; the area within the dashed lines are used in the further analysis, i.e., the data for $E_\gamma > 2.00$ MeV, $5.43 < E_x < 10.98$ MeV are selected. (c) Projection of the primary γ spectra for the range of excitation energies between the lines.

and oxygen lines. Experimental response functions have been recorded in-beam for γ transitions of ^{13}C , ^{16}O , ^{28}Si , and ^{56}Fe . The unfolding procedure is described in detail in Ref. [18]; the main advantage of this method is that the original, statistical fluctuations are preserved without introducing spurious fluctuations. This is obtained by applying a strong smoothing on the Compton background spectrum before subtracting it from the raw spectrum. This approach is justified by the fact that the Compton background varies slowly with γ -ray energy. Thus, we avoid spurious structures in the final, unfolded spectrum.

The experimental spin range for the present experiment can be inferred from both the singles-proton spectrum and the proton- γ coincidences. From the unfolded proton- γ coincidences, we clearly identify transitions for states with spin-parity up to $7/2^+$. From the singles-proton spectrum, we see that the 909 keV isomer with spin-parity $9/2^+$ is populated. This provides a lower limit for the populated spins.

B. Primary γ -ray spectra and functional form of level density and γ -ray strength

Once the γ spectra were properly unfolded, an iterative method [19] was applied to obtain the distribution of primary γ transitions from secondary and higher-order transitions. The principle of this method is that for a given excitation-energy bin j , this will contain all the γ -rays of the decay cascades from lower-lying bins $i < j$, and in addition the primary transitions for bin j . Thus, by subtracting a weighted sum of the spectra below j , the distribution of primary γ rays for bin j is obtained. Systematic uncertainties of this procedure is discussed in detail in Ref. [20]. The obtained matrix of primary γ -ray spectra is displayed in Fig. 2(b), and the projection of the primary γ rays for $E_x \approx 5.5\text{--}11.0$ MeV is shown in Fig. 2(c). We note that the spectrum in Fig. 2(b) is rather smooth, although some structures appear, particularly at lower E_γ values.

The data within the dashed lines are used for the extraction of NLD and γ SF in the next step. The boundaries are chosen for the following reasons: we need to make sure that the decay originates from a region of fairly high level density, and also that the primary γ spectra are indeed correct. An indicator for how well the primary- γ extraction procedure works is

calculated for each iteration and for each excitation-energy bin. In short, this indicator shows whether the primary spectrum corresponds to a γ multiplicity of 1, which it obviously should, and in such cases the indicator is unity. We allow for a variation in this indicator of $\pm 15\%$; if the deviation from unity is larger, we do not use the primary spectra from that excitation-energy region (see Ref. [19] for more details). Therefore, cuts are made in the matrices as shown in Fig. 2, ensuring relatively high initial excitation energies and thus high initial level density: $E_x = 5.44\text{--}10.97$ MeV, and $E_\gamma^{\text{low}} = 2.01$ MeV for ^{89}Y . The E_γ^{low} limit is necessary because of strong, discrete transitions being either subtracted too little or too much below this energy, resulting in vertical “ridges” or “valleys” in the primary γ matrix (see also the discussion in Ref. [20]).

To extract the NLD and the γ SF from the set of primary γ -ray spectra, we make use of the following relation [10]:

$$P(E_\gamma, E_x) \propto \rho(E_f) T(E_\gamma). \quad (1)$$

Here, $P(E_\gamma, E_x)$ is the experimental primary γ -ray matrix, as shown in Fig. 2, but where the primary γ spectra of each excitation-energy bin are normalized to unity to represent the probability of decay from that bin [10]. The matrix of primary γ spectra, $P(E_\gamma, E_x)$, is proportional to the level density at the final excitation energy $E_f = E_x - E_\gamma$, and to the γ -ray transmission coefficient $T(E_\gamma)$. The latter is dependent on the γ -ray energy only, in accordance with the generalized form of the Brink-Axel hypothesis [21,22]. The generalized Brink-Axel hypothesis has very recently been experimentally verified for γ transitions in the quasicontinuum [23]. The expression in Eq. (1) is valid for statistical decay, i.e., where the decay is independent of the formation of the compound state [24].

The functional form of the NLD and γ SF is determined through a least- χ^2 fit to the $P(E_\gamma, E_x)$ matrix, as described in Ref. [10]. The absolute normalization of the functions remains to be found, i.e., determining the parameters $\mathcal{A}$, $\mathcal{B}$, and α in the transformations

$$\rho(E_x - E_\gamma) = \mathcal{A} \exp[\alpha(E_x - E_\gamma)] \bar{\rho}(E_x - E_\gamma), \quad (2)$$

$$T(E_\gamma) = \mathcal{B} \exp(\alpha E_\gamma) \tilde{T}(E_\gamma), \quad (3)$$

which all give equally good fits to the experimental data. This normalization will be described in the following.

C. Normalization of level densities

To determine the absolute scaling $\mathcal{A}$ and the slope α of the level density, our data points are normalized to discrete levels [28] at low excitation energy. At the neutron-separation energy S_n , data on average s -wave neutron-resonance spacings D_0 [25] are usually used to calculate the total level density $\rho(S_n)$ for all spins and both parities. For the case of ^{89}Y , there are no neutron-resonance data available because ^{88}Y is unstable. To estimate reasonable normalization parameters we considered systematics of s -wave ($\ell = 0$) resonance spacings for this mass region by using the most recent evaluation of the Reference Input Parameter Library (RIPL-3, Ref. [25]). In addition, systematic errors due to the spin distribution at S_n must be taken into account [20]. This will be discussed in the following.

In our recent work [11], the $^{89,90}\text{Y}$ level densities were normalized by using the global systematics of Ref. [27] within the constant-temperature (CT) approximation of the level-density function. This approach gives a very good description of the functional form of experimental level densities above $\approx 2\Delta$ [29], where $\Delta \approx 1$ MeV is the pair-gap parameter. However, it yields rather low (and constant) values for the spin-cutoff parameter. Moreover, this is only one out of several possible values for the spin-cutoff parameter and hence for the upper normalization point. Also, because the spin distribution for excitation energies up to the neutron-separation energy is needed for the normalization of the γ SF later on, we here rely on an excitation-energy-dependent spin-cutoff parameter. We have chosen three different approaches, the first one using a phenomenological Fermi-gas (FG) spin-cutoff parameter following Ref. [27] (FG09):

$$\sigma_{09}^2(E_x) = 0.391A^{0.675}(E_x - 0.5Pa')^{0.312}. \quad (4)$$

Here, E_x is the excitation energy, A is the mass number, and Pa' is the deuteron pairing energy as defined in Ref. [27]. Second, we consider the FG spin-cutoff parameter of Ref. [26], where a rigid-body moment of inertia is assumed (FG05):

$$\sigma_{05}^2(E_x) = 0.0146A^{5/3} \frac{1 + \sqrt{1 + 4a(E_x - E_1)}}{2a}. \quad (5)$$

Here, a is the level-density parameter and E_1 is the excitation-energy backshift determined from global systematics of Ref. [26]. For the phenomenological spin-cutoff parameters, the spin distribution is given by the standard expression [30,31]:

$$g(E_x, J) \simeq \frac{2J+1}{2\sigma^2(E_x)} \exp\left[-\frac{(J+1/2)^2}{2\sigma^2(E_x)}\right]. \quad (6)$$

By using the phenomenological spin-cutoff parameters and assuming equiparity (as shown in Ref. [11]), the total level density at S_n can be estimated from D_0 through the

FIG. 3. Spin distributions for ^{89}Y at the neutron-separation energy $S_n = 11.482$ MeV for the three different normalization approaches. For FG09 and FG05, Eq. (6) is used with their respective spin-cutoff parameters in Eqs. (4) and (5). The HFB + c calculations assume no specific shape of the spin distribution but happens to be very similar to FG05 in this case.

expression [10]

$$\rho(S_n) = \frac{2\sigma^2}{D_0} \frac{1}{(J_t + 1) \exp[-(J_t + 1)^2/2\sigma^2] + J_t \exp[-J_t^2/2\sigma^2]}, \quad (7)$$

where J_t is the ground-state spin of the target nucleus in the neutron-resonance experiment. Finally, we use microscopic calculations within the Hartree-Fock-Bogoliubov plus combinatorial (HFB + c) approach [32]. The three spin distributions are shown in Fig. 3 for the case of ^{89}Y .

We observe that the spin distribution of Eq. (6) using σ_{05} is very broad and centered at significantly higher spins than when using σ_{09} . We consider therefore σ_{05} to give the upper limit, and σ_{09} as the lower limit in estimating $\rho(S_n)$ for $^{89,90}\text{Y}$. The calculated $\rho(S_n)$ values for Rb, Sr, Y, and Zr isotopes are given in Table I together with the applied input parameters. The resulting systematics for the level densities at S_n are shown in Fig. 4. Note that the predictions from the global systematics are fit to the semi-experimental data points through a common scaling factor of $0.31^{+0.09}_{-0.10}$ and $0.34^{+0.08}_{-0.07}$ for $\rho_{05}(S_n)$ and $\rho_{09}(S_n)$, respectively. By using the upper (lower) χ^2 uncertainty for the σ_{05} (σ_{09}) results, we get the following estimates for ^{89}Y ($S_n = 11.482$ MeV): $\rho_{09}^{\text{low}}(S_n) = 4.87 \times 10^4 \text{ MeV}^{-1}$, corresponding to $D_0 = 143$ eV for $\sigma_{09}(S_n) = 4.12$, and $\rho_{05}^{\text{up}}(S_n) = 9.33 \times 10^4 \text{ MeV}^{-1}$, corresponding to $D_0 = 100$ eV for $\sigma_{05}(S_n) = 5.45$. Finally, the HFB + c calculations (with no excitation-energy shift, $\delta = 0.0$ MeV where δ is defined in Ref. [32]) yield $\rho_{\text{HFB+c}}(S_n) = 7.09 \times 10^4 \text{ MeV}^{-1}$ with $D_0 = 121$ eV. A fit of Eq. (6) on the HFB + c calculations at S_n gives $\sigma_{\text{HFB+c}} = 4.89$, i.e., in between the two phenomenological approaches.

TABLE I. Neutron resonance parameters D_0 and $\langle\Gamma_{\gamma 0}\rangle$ from Ref. [25], and spin-cutoff parameters from global systematics of Refs. [26,27]; A_f is the final nucleus following neutron capture, J_t is the ground-state spin of the target nucleus, S_n is the neutron-separation energy, σ_{05} and σ_{09} are the spin-cutoff parameters from Eqs. (4) and (5), D_0 is the s -wave level spacing [25], and $\rho_{05}(S_n)$ and $\rho_{09}(S_n)$ are the total level densities calculated from σ_{05} and σ_{09} . Finally, ρ_{05}^{syst} and ρ_{09}^{syst} are the total level densities at S_n as predicted from the global systematics of Refs. [26,27], respectively.

A_f	J_t	S_n (MeV)	D_0 (keV)	$\sigma_{05}(S_n)$	$\rho_{05}(S_n)$ (10^4 MeV^{-1})	$\rho_{05}^{\text{syst}}(S_n)$ (10^4 MeV^{-1})	$\sigma_{09}(S_n)$	$\rho_{09}(S_n)$ (10^4 MeV^{-1})	$\rho_{09}^{\text{syst}}(S_n)$ (10^4 MeV^{-1})	$\langle\Gamma_{\gamma 0}\rangle$ (meV)
^{86}Rb	5/2	8.651	0.172(8)	5.12	6.11(28)	21.6	4.02	4.22(20)	12.4	250(10)
^{88}Rb	3/2	6.083	1.630(150)	4.80	0.78(7)	1.47	3.85	0.53(5)	1.14	170(30)
^{85}Sr	0	8.530	0.320(120)	4.89	15.3(57)	20.4	3.91	9.88(370)	7.88	240(80)
^{87}Sr	0	8.424	2.600(800)	5.07	2.02(62)	13.2	3.94	1.23(38)	4.01	260(80)
^{88}Sr	9/2	11.11	0.290(80)	5.22	3.00(83)	13.1	4.02	2.44(67)	6.40	150(40)
^{89}Sr	0	6.359	23.70(290)	4.72	0.19(2)	0.53	3.76	0.12(2)	0.39	190(50)
^{90}Y	1/2	6.857	3.700(400)	4.99	0.70(8)	2.25	3.93	0.44(5)	1.71	130(40)
^{91}Zr	0	7.194	6.000(1400)	4.95	0.83(19)	1.20	3.88	0.52(12)	0.91	130(20)
^{92}Zr	5/2	8.635	0.550(100)	5.03	1.85(34)	4.11	3.95	1.29(23)	2.60	140(40)
^{93}Zr	0	6.734	3.500(800)	4.84	1.37(31)	1.69	3.86	0.88(20)	1.31	135(25)
^{94}Zr	5/2	8.221	0.302(75)	4.95	3.29(82)	5.62	3.94	2.34(58)	3.63	157(20)
^{95}Zr	0	6.462	4.000(800)	4.79	1.17(24)	1.99	3.86	0.77(15)	1.54	85(20)
^{97}Zr	0	5.575	13.00(300)	4.66	0.34(8)	0.58	3.76	0.23(5)	0.72	65(15)

Due to the cutoff on the γ energy, E_γ^{low} (see Fig. 2), our level-density data points will reach a maximum excitation energy of $S_n - E_\gamma^{\text{low}}$. Following Ref. [11], we use an interpolation between our data points up to $\rho(S_n)$ of the constant-temperature form [30]:

$$\rho_{\text{CT}}(E_x) = \frac{1}{T} \exp \frac{E_x - E_0}{T}, \quad (8)$$

where E_0 is the excitation-energy shift and T is the constant nuclear temperature. For ^{89}Y , we used $(E_0, T) = (0.658, 0.95)$, $(0.069, 1.02)$, and $(0.098, 1.05)$ MeV for FG05, HFB + c, and FG09, respectively.

The normalized level density is shown in Fig. 5. Our data points follow the discrete levels well up to $E_x \approx 4.5$ MeV, which means that these levels are populated by primary γ

FIG. 4. Level densities at S_n in the Y mass region with global systematics of (a), (b) Ref. [26] (FG05) and (c), (d) Ref. [27] (FG09) (see Table I). The unknown level density for ^{89}Y is shown as a purple diamond. The global-systematics predictions are scaled with a factor of 0.31 and 0.34 for the FG05 and FG09 approaches, respectively. The error bands show the upper-limit scaling of (a), (b) 0.40 for FG05 and (c), (d) the lower limit of 0.27 for FG09.

FIG. 5. The extracted level density of ^{89}Y . The black points give the HFB + c normalization, the lower and upper limits are shown as azure lines. The inset shows a zoom of the high- E_x part.

decay from the higher-lying levels in the region; above this energy the level densities increase rapidly, while the known levels show a saturation when reaching $\approx 65 \text{ MeV}^{-1}$. As discussed in Ref. [11], the level density displays a remarkable constant-temperature behavior. The lower and upper limits for the normalization are also shown, representing the systematic errors. Note that the lower limit is very similar to the normalization used in Ref. [11] (within 18%). At low excitation energies there is obviously not much difference between the normalization options, because our data points are fixed to the discrete levels. At the neutron-separation energy the systematic uncertainty is at its maximum, within a factor of ≈ 2 (see inset of Fig. 5).

D. γ -ray strength function

The slope of the γ SFs is determined through the normalization of the level densities [see Eqs. (2) and (3)]. The absolute scale $\mathcal{B}$ can be found by use of the total, average radiative width $\langle \Gamma_{\gamma 0} \rangle$. The average radiative width of neutron s -wave capture resonances with spins $J_t \pm 1/2$ expressed in terms of the experimental $\mathcal{T}$ is given by [33]

$$\begin{aligned} \langle \Gamma_{\gamma 0}(S_n, J_t \pm 1/2, \pi_t) \rangle &= \frac{\mathcal{B}}{4\pi\rho(S_n, J_t \pm 1/2, \pi_t)} \\ &\times \int_{E_\gamma=0}^{S_n} dE_\gamma \mathcal{T}(E_\gamma) \rho(S_n - E_\gamma) \\ &\times \sum_{J=-1}^1 g(S_n - E_\gamma, J_t \pm 1/2 + J), \end{aligned} \quad (9)$$

where J_t and π_t are the spin and parity of the target nucleus in the (n, γ) reaction, $\rho(S_n - E_\gamma)$ is the normalized, experimental level density obtained in Sec. II C, and $\mathcal{T}$ is the experimental

FIG. 6. Normalized γ SF of ^{89}Y compared to $^{89}\text{Y}(\gamma, n) + ^{89}\text{Y}(\gamma, np)$ data from Refs. [36,37], and evaluated data from Ref. [38]. The black points are obtained with the HFB + c normalization, and the azure lines show the lower and upper limits.

transmission coefficient, which in principle includes all types of electromagnetic transitions: $\mathcal{T}_{E1} + \mathcal{T}_{M1} + \mathcal{T}_{E2} + \dots$. The sum runs over all final states with spins $J_t \pm 1/2 + J$, where $J = -1, 0, 1$ from considering the spins reached after one primary dipole transition with energy E_γ [see also Eq. (3.1) in Ref. [34]]. Note that the factor $1/\rho(S_n, J_t \pm 1/2, \pi_t)$ equals the neutron resonance spacing D_0 . From the normalized transmission coefficient, the γ SF is determined by

$$f(E_\gamma) = \frac{\mathcal{T}(E_\gamma)}{2\pi E_\gamma^3} \quad (10)$$

by using the fact that dipole transitions dominate the strength for the E_x region considered [34,35].

For ^{89}Y , we estimated the unknown $\langle \Gamma_{\gamma 0} \rangle$ from data in this mass region [25]. Specifically, we took the average value of the nuclei close in mass to ^{89}Y ; namely, $^{88,89}\text{Sr}$, ^{90}Y , and ^{91}Zr ; see Table I. With an uncertainty of $\approx 25\%$, we obtain $\langle \Gamma_{\gamma 0} \rangle(^{89}\text{Y}) = 150(38) \text{ meV}$. The assumed uncertainty of 25% is strongly guided by the comparison with photonuclear data from the reactions $^{89}\text{Y}(\gamma, n) + ^{89}\text{Y}(\gamma, np)$ [36–38]. These cross-section data are converted into γ SF by the relation [25]

$$f(E_\gamma) = \frac{\sigma_{(\gamma, n)}(E_\gamma)}{3\pi^2 \hbar^2 c^2 E_\gamma}, \quad (11)$$

again assuming that dipole radiation is dominant, which is reasonable in this E_γ region (see, e.g., Ref. [25] and references therein).

The normalized γ SF is shown in Fig. 6. The error bands include the uncertainty in $\langle \Gamma_{\gamma 0} \rangle$ as well as in the level density and the choice of spin distribution. The uncertainty in absolute value ranges from a factor of 1.8 at $E_\gamma = 2.0 \text{ MeV}$ to ≈ 4.8 at $E_\gamma = 11.3 \text{ MeV}$.

For γ -ray energies above 3–4 MeV, we find that the strength is increasing as a function of γ -ray energy, as expected for

the tail of the giant dipole resonance (GDR) [39]. We also observe a drop in strength for γ rays between ≈ 9.6 –11.0 MeV. This is understood by looking at the primary γ -ray matrix in Fig. 2(a), where it is clear that the upper-right corner in the triangle has significantly fewer statistics than for lower γ -ray energies. This could be due to less-direct feeding to the ground state, with spin-parity $1/2^-$, and more to the first excited $9/2^+$ state at 909 keV. We note that this behavior is in agreement with the (γ,γ') data from Ref. [40], which also display a reduction in intensity in the energy range $E_\gamma \approx 9.8$ –11.3 MeV. This indicates that, in this particular region, the extracted data are not representative of a general γ SF, because there is, very likely, a strong dependence on the final state(s) and the individual overlap with the initial and final state(s) of the transition(s).

Our data show an increase at decreasing γ energies for $E_\gamma < 4$ MeV. This phenomenon, hereafter called the *upbend*, has been subject of great interest recently, and was first discovered in iron isotopes about a decade ago [41]. In Ref. [20], simulations on ^{57}Fe using the DICEBOX code [42] suggest that an enhancement in the $E1$ strength could be invoked, although it was not present in the input $E1$ strength for the simulation. It was stressed that the low-energy enhancement could not be due to artifacts in the unfolding or the method for extracting primary transitions, because the same feature was seen by using the primary transitions directly from the DICEBOX simulations. Three main reasons for this behavior were pointed out:

- (i) There was no significant contribution from quadrupole ($E2$) transitions in the simulations.
- (ii) Specific restrictions were applied on the initial spin population.
- (iii) The level density was very low for high spins in this particular simulation.

Hence, the simulations indicated an increase in the low-energy part of the γ SF in cases where the level density is low and the reaction populates selectively high spins and a rather narrow spin range.

However, this hypothesis was disproved in Ref. [13], where a different and virtually model-independent technique was applied on data from the $^{94}\text{Mo}(d,p\gamma\gamma)^{95}\text{Mo}$ reaction. In contrast to the $^{96}\text{Mo}(^3\text{He},\alpha)^{95}\text{Mo}$ data from Ref. [12], where the reaction favors high- ℓ transfer, the (d,p) reaction mainly populates low- ℓ states. Nevertheless, the same shape of the γ SF was found [13]. Furthermore, by using the Oslo method, it was shown in Ref. [35] that the $^{56}\text{Fe}(p,p')^{56}\text{Fe}$ and $^{57}\text{Fe}(^3\text{He},\alpha)^{56}\text{Fe}$ data sets yielded very similar γ SFs. Therefore, considering the available data as of today, other explanations must be sought to explain the upbend feature.

The upbend has recently been shown to be dominantly of dipole nature in ^{56}Fe [35] and in $^{151,153}\text{Sm}$ [43], but the electromagnetic character is not known at present. Theoretical attempts suggest that both $E1$ as well as $M1$ strength may contribute. An enhancement of $E1$ type is predicted from a thermal coupling of quasiparticles to the continuum of unbound states at relatively high temperature [44], whereas an enhancement of $M1$ strength is found in shell-model

calculations [14,45]. Moreover, the presence of the upbend may enhance astrophysical (n,γ) rates of exotic neutron-rich nuclei by up to two orders of magnitude [46]. In the following, we present shell-model calculations as well as models for the $E1$ strength and compare with our data.

III. SHELL-MODEL CALCULATIONS AND MODELS FOR DIPOLE STRENGTH

The shell-model calculations were performed by means of the code RITSSCHIL [47] by using a model space composed of the $(0f_{5/2}, 1p_{3/2}, 1p_{1/2}, 0g_{9/2})$ proton orbits and the $(0g_{9/2}, 1d_{5/2}, 0g_{7/2})$ neutron orbits relative to a ^{68}Ni core. This configuration space was also applied in our earlier study of $M1$ strength functions in $^{94,95,96}\text{Mo}$ and ^{90}Zr [14]. In the present calculations, two protons were allowed to be lifted from the fp shell to the $0g_{9/2}$ orbit and two neutrons from the $0g_{9/2}$ to the $1d_{5/2}$ orbit. This resulted in dimensions up to 29 000. The additional inclusion of the $\nu(0g_{7/2})$ orbit has negligible influence on the low-energy part of the strength function but produces a few strong $M1$ transitions around 7 MeV dominated by the $\nu(0g_{9/2})^{-1}\nu(0g_{7/2})$ spin-flip configuration (cf. Ref. [14]). As these transitions do not describe the spin-flip peak completely up to high energy, we use a phenomenological description for the spin-flip resonance in the present work.

The calculations included states with spins from $J = 1/2$ to $21/2$ for ^{89}Y . For each spin the lowest 40 states were calculated. Reduced transition probabilities $B(M1)$ were calculated for all transitions from initial to final states with energies $E_f < E_i$ and spins $J_f = J_i, J_i \pm 1$. For the minimum and maximum J_i , the cases $J_f = J_i - 1$ and $J_f = J_i + 1$, respectively, were excluded. This resulted in more than 32 000 $M1$ transitions for each parity $\pi = +$ and $\pi = -$, which were sorted into 100 keV bins according to their transition energy $E_\gamma = E_i - E_f$. The average $B(M1)$ value for one energy bin was obtained as the sum of all $B(M1)$ values divided by the number of transitions within this bin.

The $M1$ strength functions were deduced by using the relation

$$f_{M1}(E_\gamma) = \frac{16\pi}{9(\hbar c)^3} \overline{B}(M1, E_\gamma) \rho(E_i). \quad (12)$$

They were calculated by multiplying the $B(M1)$ value in μ_N^2 of each transition by 11.5473×10^{-9} times the level density at the energy of the initial state $\rho(E_i)$ in MeV^{-1} and deducing averages in energy bins, as done for the $\overline{B}(M1)$ values (see above). The level densities $\rho(E_i, \pi)$ were determined by counting the calculated levels within energy intervals of 1 MeV for the two parities separately. The strength functions obtained for the two parities were subsequently added. When calculating the strength functions, gates were set on the excitation energy E_x that correspond to the ones applied in the analysis of the experimental data (see Sec. II). The resulting $M1$ strength function for ^{89}Y is shown in Fig. 7. The low-energy behavior of this strength function is very similar to that of the strength functions calculated for the neighboring nuclei $^{94,95,96}\text{Mo}$, ^{90}Zr [14] and for $^{56,57}\text{Fe}$ [45].

The low-energy enhancement of $M1$ strength is caused by transitions between many close-lying states of all considered

FIG. 7. The γ SF of ^{89}Y together with photonuclear data [36,37] and evaluated (γ, n) data from Ref. [38] compared with models for the dipole strength.

spins located above the yrast line in the transitional region to the quasicontinuum of nuclear states. Inspecting the wave functions, one finds large $B(M1)$ values for transitions between states that contain a large component (up to about 50%) of the same configuration with broken pairs of both protons and neutrons in high- j orbits, whereas states containing the unpaired $1p_{1/2}$ proton and proton excitations only are not depopulated by strong $M1$ transitions. The largest $M1$ matrix elements connect configurations with the spins of high- j protons recoupled with respect to those of high- j neutrons to the total spin $J_f = J_i, J_i \pm 1$. The corresponding main configurations for negative-parity states in ^{89}Y are generated by exciting neutrons over the shell gap at $N = 50$, such as $\pi(1p_{1/2}^{-1}\nu(0g_{9/2}^{-1}1d_{5/2}^1))$ or $\pi(1p_{1/2}^{-1}\nu(0g_{9/2}^{-2}1d_{5/2}^2))$ and by additional proton excitations within the (fp) shell, i.e., $\pi[(0f_{5/2}, 1p_{3/2})^{-1}1p_{1/2}^2\nu(0g_{9/2}^{-1}1d_{5/2}^1)]$ and also proton excitations over the subshell gap at $Z = 40$, $\pi[(0f_{5/2}, 1p_{3/2})^{-1}1p_{1/2}^0 0g_{9/2}^2\nu(0g_{9/2}^{-1}1d_{5/2}^1)]$. The positive-parity states require the excitation of an (fp) proton to the $0g_{9/2}$ orbit, for example $\pi(1p_{3/2}^{-1}1p_{1/2}^1 0g_{9/2}^1\nu(0g_{9/2}^{-1}1d_{5/2}^1))$. The orbits in these configurations have large g factors with opposite signs for protons and neutrons. Combined with specific relative phases of the proton and neutron partitions they cause large total magnetic moments.

For a comparison with the experimental data, an $E1$ contribution to the strength function has to be added. We have chosen two ways to estimate the $E1$ strength: (i) the phenomenological generalized Lorentzian (GLo) model [34], and (ii) a microscopic approach based on the quasiparticle-random-phase approximation (QRPA) [48,49]. For option (i), we apply the GLo model with a constant temperature for the final states T_f , in contrast to a variable temperature which depends on the final excitation energy. The choice of a constant temperature is in accordance with the Brink hypothesis [21] and our ansatz that the γ -transmission coefficient is, on average, independent of excitation energy in the statistical

FIG. 8. Same as Fig. 7 but with the QRPA $E1$ strength (see text).

E_x region. This is also in accordance with the constant-temperature level density found for ^{89}Y . The GLo model is then given by

$$f_{\text{GLo}}(E_\gamma, T_f) = \frac{1}{3\pi^2 \hbar^2 c^2} \sigma_{E1} \Gamma_{E1} \times \left[\frac{E_\gamma \Gamma(E_\gamma, T_f)}{(E_\gamma^2 - E_{E1}^2)^2 + E_\gamma^2 \Gamma(E_\gamma, T_f)^2} + 0.7 \frac{\Gamma(E_\gamma = 0, T_f)}{E_{E1}^3} \right], \quad (13)$$

with

$$\Gamma(E_\gamma, T_f) = \frac{\Gamma_{E1}}{E_{E1}^2} (E_\gamma^2 + 4\pi^2 T_f^2). \quad (14)$$

The parameters Γ_{E1} , E_{E1} , and σ_{E1} correspond to the width, centroid energy, and peak cross section of the GDR. For option (ii), the $E1$ strength is obtained from large-scale QRPA calculations on top of a Skyrme–Hartree–Fock–Bogoliubov description of the ground state. The QRPA calculations are performed in the spherical approximation, and a folding procedure is applied to obtain the correct spreading width of the GDR. For more details on the QRPA calculations, we refer the reader to Ref. [49]. The $E1$ calculations were taken from the BRUSLIB library [50].

We included an $M1$ spin-flip resonance with a standard Lorentzian form [25] by using parameters in accordance with a recent $(p, p')^{90}\text{Zr}$ experiment [51]. Strong $M1$ transitions were also observed in a photon-scattering experiment in the excitation-energy region ≈ 8 –10 MeV [52]. For the QRPA calculation, we had to shift the $E1$ strength with an energy shift of $\delta_{\text{QRPA}} = +1.5$ MeV, i.e., $E_\gamma^{\text{new}} = E_\gamma + \delta_{\text{QRPA}}$, so as to match the GDR data reasonably well around $E_\gamma = 12$ MeV. The resulting theoretical dipole strengths are displayed in Figs. 7 and 8 for the best reproduction of the HFB + c normalization. All the parameters used for the shown models are listed in Table II.

TABLE II. Parameters used for the model strength functions of ^{89}Y in Figs. 7 and 8.

Nucleus	Γ_{E1} (MeV)	E_{E1} (MeV)	σ_{E1} (mb)	T_f (MeV)	Γ_{M1} (MeV)	E_{M1} (MeV)	σ_{M1} (mb)	δ_{QRPA} (MeV)
^{89}Y	4.3	16.8	233.0	0.30	2.7	9.5	1.1	1.5

In the following, we use our experimentally inferred lower and upper limits on the level density and γSF as input for cross-section calculations of the $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$ and $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ reactions.

IV. CROSS SECTION AND RATE CALCULATIONS

As discussed in the introduction, level density and γSF are two key ingredients in the Hauser–Feshbach approach to calculate cross sections. In this work, we use the open-source nuclear reaction code TALYS-1.6 [54,55] for the cross-section and reaction-rate calculations for the $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ and $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$ reactions. Our approach is the following:

- (1) Making use of all the various models already implemented in TALYS for the level density, the γSF , and the proton and neutron optical potentials to investigate the spread in the resulting cross sections and reaction rates; i.e., inferring the lower and upper limits on these quantities inherent from the available models.
- (2) Calculating the cross sections and reaction rates with default input parameters in TALYS.
- (3) Implementing level densities and γSFs in accordance with our present data and the experimentally inferred lower and upper limits.

A. TALYS predictions for model uncertainties

We calculated all possible combinations of input level densities, γSF models, and optical potentials available in TALYS to estimate the minimum and maximum $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$ and $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ cross sections predicted by these models. For the (n,γ) reaction, the combinations are

- (i) a minimum cross section and rate with the temperature-dependent Hartree–Fock–Bogoliubov plus combinatorial level density of Ref. [56], the $E1$ strength function from the Hartree–Fock–BCS plus QRPA approach of Ref. [57] making use of the renormalization to the estimated $\langle\Gamma_\gamma\rangle = 170$ meV from a spline-fit interpolation table in TALYS, and the JLM neutron potential [58] (TALYS keywords *ldmodel 6, strength 3, gnorm -1., jlmomp y*).
- (ii) a maximum cross section and rate with the combined constant-temperature plus back-shifted Fermi gas model [59] for the level density with parameters according to the TALYS manual, the standard Lorentzian model (Brink–Axel) for the $E1$ strength [21,22], and a global parametrization of the neutron optical potential [60] (TALYS keywords *ldmodel1, strength2, localomp n*).

The corresponding combinations for the (p,γ) reaction are

- (i) a minimum cross section and rate with the temperature-dependent Hartree–Fock–Bogoliubov plus combinatorial level density of Ref. [56], the GLo model [34] with variable temperature as implemented in TALYS renormalized to the spline-fit $\langle\Gamma_\gamma\rangle$, and the global parametrization of the proton optical potential [60] (TALYS keywords *ldmodel 6, strength 1, gnorm -1., localomp n*).
- (ii) a maximum cross section and rate with the combined constant-temperature plus back-shifted Fermi gas model [59] for the level density with parameters according to the TALYS manual, the standard Lorentzian model (Brink–Axel) for the $E1$ strength [21,22], and the JLM proton potential [58] (TALYS keywords *ldmodel 1, strength 2, jlmomp y*).

We find that the level-density models in TALYS give a factor of ≈ 5 and ≈ 4 uncertainty for the (n,γ) and (p,γ) cross sections, respectively, while the corresponding numbers for the γSF models are ≈ 30 and ≈ 28 , respectively. The impact of the choice of optical-model potential has also been tested for the reactions of interest. More specifically, we have used the proton and neutron potentials of Koning and Delaroche [60] with global parameters as described in the TALYS manual, and also the semimicroscopic optical potential of the Jeukenne–Lejeune–Mahaux (JLM) type [58]; see the TALYS documentation for more details [54,55]. We did not adjust any parameters in the neutron or proton potentials but used the default parameters as implemented in TALYS-1.6. We have found that the ratio between the cross sections using the global potential and the JLM potential reaches a maximum deviation of $\approx 34\%$ and $\approx 48\%$ for the proton and neutron potentials, respectively, for the energy ranges $2.5 \times 10^{-6} \leq E_n \leq 5$ MeV and $0.75 \leq E_p \leq 5$ MeV. On average, the JLM potential gives a lower capture cross section than the global potential for the neutron capture, and opposite for proton capture.

The combined model uncertainties in the cross sections, including the optical-model uncertainties, reach a factor of ≈ 33 and ≈ 36 at maximum for the $(n,\gamma)^{89}\text{Y}$ and $(p,\gamma)^{89}\text{Y}$ cross sections. These can be considered as the intrinsic uncertainties for the models implemented in TALYS (note again that no information is available regarding the level spacing D_0 and the average radiative width $\langle\Gamma_\gamma\rangle$). The results are shown as thick, black lines in Figs. 9 and 10.

B. TALYS default predictions

If no keywords related to level density, γSF , or proton or neutron potential are specified in the TALYS input file, default values will automatically be invoked. For the $(n,\gamma)^{89}\text{Y}$ reaction, these are the neutron optical potential [60] with local parameters, the GLo model with variable temperature [34], and the combined constant-temperature plus backshifted-Fermi-gas model [59] for the level density. For the $(p,\gamma)^{89}\text{Y}$ reaction these default values are the same, except for the proton potential, for which there are no local parameters and the global parametrization of Ref. [60] is used. The results are displayed as black crosses in Figs. 9 and 10.

FIG. 9. (a) Input γ SFs of ^{89}Y , note that the TALYS default is for $E_n = 1 \text{ MeV}$, corresponding to an initial excitation energy of 12.48 MeV in the GLo model, and that the GLo model is by default normalized to a radiative width $\langle \Gamma_\gamma \rangle = 170 \text{ meV}$ taken from an interpolation routine in TALYS; (b) the resulting $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$ cross sections, (c) and the corresponding astrophysical reaction rates compared to the BRUSLIB (dashed magenta line, Ref. [50]) and the JINA REACLIB (green solid line, Ref. [53]). The minimum and maximum predictions from the models implemented in TALYS are also shown (thick, black lines).

It is interesting to note that, by default, there is an automatic rescaling of the γ SF in TALYS to match the estimated $\langle \Gamma_\gamma \rangle = 170 \text{ meV}$ from the spline-fit interpolation table. However, this procedure does not guarantee that the resulting (n,γ) cross section will be the same for different model combinations of the γ SF and level density. The default-estimate cross section (crosses in Fig. 9) with a scaling factor $gnorm = 1.44$ is very different from the lower-limit cross section with $gnorm = 0.28$ (lower, thick black lines in Fig. 9). This demonstrates that there is a delicate interplay with the adopted level density (with its spin distribution), γ SF and particle optical potential in the calculation of radiative cross sections.

C. Results from this work implemented in TALYS

Finally, we used our results to constrain the input level density and γ SF for the $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$ and $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$

cross sections and reaction rates. We have used the constant-temperature model for the level density with parameters given in Sec. II C above $E_x = 2.88 \text{ MeV}$, with a spin distribution and parameters according to the lower and upper normalizations as described in Sec. II C. This model reproduces our level-density data very well. Below that excitation energy, we use the known, discrete levels. For the γ SF data, we tuned the γ SF models to reproduce our lower and upper data points. Specifically, we used a low-lying $M1$ strength that corresponds to the shell-model results parametrized as an exponential function $f_{\text{upbend}} = C \exp(-\eta E_\gamma)$, an $M1$ spin-flip resonance, and a GLo $E1$ component so as to match our upper and lower limits. In addition, our γ SF data overshoot the $M1 + E1$ models in the energy range of $E_\gamma \approx 2.0\text{--}3.5 \text{ MeV}$, as seen in Figs. 7 and 8. Therefore, we added a small resonance (SR) of $M1$ type to get a reasonable agreement with the measured strength.

FIG. 10. (a) Calculated $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ cross sections shown as a blue-shaded band compared to data from Ref. [61], and (b) the corresponding astrophysical reaction rates compared to the BRUSLIB (dashed magenta line, Ref. [50]) and the JINA REACLIB (green solid line, Ref. [53]). The minimum and maximum predictions from the models implemented in TALYS are also shown (thick, black lines), as well as the result using default input parameters (black crosses).

TABLE III. Parameters used for the γ SF input models in the TALYS calculations.

	Γ_{E1} (MeV)	E_{E1} (MeV)	σ_{E1} (mb)	T_f (MeV)	Γ_{M1} (MeV)	E_{M1} (MeV)	σ_{M1} (mb)	Γ_{SR} (MeV)	E_{SR} (MeV)	σ_{SR} (mb)	C (10^{-8} MeV^{-3})	η (MeV^{-1})
Lower limit	3.5	16.8	233.0	0.30	2.7	9.5	0.24	2.2	2.8	0.10	4.0	2.5
Upper limit	6.6	17.8	233.0	0.30	2.7	9.5	2.00	2.2	2.8	0.18	5.0	2.5

The applied parameters are given in Table III. The input γ SF for ^{89}Y is shown in Fig. 9(a).

The resulting $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$ and $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ cross sections are displayed in Figs. 9(b) and 10(a), respectively, and shown as blue-shaded bands. We note that our estimated (n,γ) cross section is in the lower region of the uncertainty band provided by TALYS. Also, our error band does not match well with the default TALYS prediction, as our results indicate a significantly lower cross section.

For the $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ cross section, direct measurements exist [61]. Overall, we find a very good agreement between our calculations and the $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ data from Ref. [61], although the data indicate a slightly higher cross section for incoming proton energies $E_p > 4.3$ MeV. The overall good agreement, however, gives confidence in our approach to estimate radiative capture cross sections, as also demonstrated in Refs. [62,63].

We also estimated the astrophysical reaction rates, which are shown in Figs. 9(c) and 10(b). We compared to the BRUSLIB [50] and the JINA REACLIB [53] databases. We find that the BRUSLIB rates are rather similar to our estimated upper limits, especially for the (p,γ) reaction. In contrast, the JINA REACLIB rates are significantly higher both for the proton-capture and neutron-capture cases, especially for higher stellar temperatures.

The low-energy upbend has been shown to increase neutron-capture rates of very neutron-rich nuclei, if it is found to be present in such exotic isotopes [46]. For ^{89}Y , however, the impact of the upbend on the cross sections is very limited; for example, for the (p,γ) reaction, the relative change in cross section is at maximum 1%, and for the $^{88}\text{Y}(n,\gamma)$ reaction the maximum is 2%. However, provided that the upbend persists for very-neutron-rich Y nuclei, it may influence the (n,γ) rates of relevance for the r process.

V. SUMMARY AND OUTLOOK

In this work, we determined the nuclear level density and γ SF of ^{89}Y . We find that this nucleus exhibits a low-energy enhancement in the γ SF, in accordance with previous find-

ings in this mass region. Moreover, shell-model calculations describe the main part of the low-energy enhancement by $M1$ transitions, similar to previous shell-model results for Mo, Fe, and Zr isotopes. However, as (a contribution of) $E1$ transitions cannot be ruled out based on the available data, an experimental determination of the electromagnetic character of this enhancement is highly desired to firmly establish the mechanism behind this phenomenon.

We applied our new data as input for nuclear-reaction calculations, estimating radiative capture cross sections and reaction rates for the reactions $^{88}\text{Sr}(p,\gamma)^{89}\text{Y}$ and $^{88}\text{Y}(n,\gamma)^{89}\text{Y}$. For the radiative proton capture case, we find good agreement with our experiment-constrained calculations and direct measurements. Moreover, in comparison with widely used reaction-rate libraries, we note that the BRUSLIB predictions in general agree better with our results.

In the future, it would be very valuable to measure other nuclei in this mass region with the Oslo method, to be able to further constrain reactions such as $^{87}\text{Y}(n,\gamma)^{88}\text{Y}$ of importance to the p -process nucleosynthesis. To be able to reduce the errors associated with our approach, further developments of independent absolute-normalization techniques would be highly desirable, as well as complementary experiments, such as measurements of particle-evaporation spectra and γ -ray two-step cascade spectra to infer level densities and γ SFs, respectively.

The data are available in Ref. [64].

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