

R-MATRIX ANALYSIS AND STATISTICAL PROPERTIES OF DYSPROSIUM ISOTOPES IN THE NEUTRON ENERGY RANGE UP TO A FEW KEV



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June 2023



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Nuclear Criticality Safety Program

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May 2023

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managed by
UT-Battelle LLC
for the
US DEPARTMENT OF ENERGY
under contract DE-AC05-00OR22725

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ABBREVIATIONS

BW	Breit–Wigner
ENDF	Evaluated Nuclear Data File
LINAC	Linear Accelerator
NIST	National Institute of Standards and Technology
NCSP	Nuclear Criticality Safety Program
ORNL	Oak Ridge National Laboratory
RM	Reich–Moore
RPC	resonance parameter covariance
RPI	Rensselaer Polytechnic Institute
RRR	Resolved Resonance Region
ToF	Time-of-Flight

ACKNOWLEDGMENTS

This work was supported by the Nuclear Criticality Safety Program, which is funded and managed by the National Nuclear Security Administration for the US Department of Energy. The author also thanks Devin Barry and Robert Block for helpful discussions on the RPI measurements and related experimental corrections.

ABSTRACT

In support of the Nuclear Criticality Safety Program, a set of evaluated resonance parameters was generated for seven dysprosium isotopes in the neutron energy range from thermal up to a few keV. The evaluation methodology used the Reich-Moore approximation to fit, with the *R*-matrix code SAMMY, the high-resolution capture and transmission measurements on natural and enriched samples recently performed at the Rensselaer Polytechnic Institute Gaertner LINear ACcelerator facility. Additional transmission data measured on enriched samples by Liou at the Columbia University Nevis synchrocyclotron in the mid seventies were used to gauge the neutron widths above 15 eV. Thermal constants such as absorption and (in)coherent scattering cross sections and corresponding scattering lengths were calibrated to the National Institute of Standards and Technology's compilation except for $^{161,164}\text{Dy}$ isotopes.

1. INTRODUCTION

Named from the Greek *dysprositos* meaning “hard to get,” dysprosium is a rare earth element that can have significant effect in nuclear reactor applications due to its property of absorbing neutrons continuously and effectively for a long time. In fact, because of its large capture cross section, dysprosium is usually exploited in the design of nuclear fuel or reactor control rods, for which dysprosium’s slow-burnout property is necessary. Being a rare earth element, dysprosium’s capture cross section is also important in the study of nucleosynthesis in astrophysics analyses. Among other common uses, dysprosium (in compound forms or dysprosium alloys) is in great demand because of its favorable thermophysical properties, as dysprosium resists demagnetization at high temperatures.

In view of the interest in dysprosium metal for several distinct types of nuclear applications, it is important to have a thorough understanding of its interaction with neutrons. To describe the reaction cross sections in the thermal and low-energy range, resonance parameters are usually stored in evaluated nuclear data file (ENDF) libraries. In the recently released ENDF/B-VIII.0 nuclear data library [1], the set of dysprosium evaluations in the resolved resonance region (RRR) was performed in 2001 (after the release of the ENDF/B-VI.8 library) and successively adopted by two releases of ENDF library—namely, ENDF/B-VII.0 [2] and -VII.1 [3]. Evaluated by the unfavorable multilevel Breit–Wigner (BW) approximation, the set of resonance parameters was adopted from the work of Mughabghab [4], except minor corrections to the bound levels were performed so as to reproduce more accurately the values of the thermal capture cross sections and the bound coherent scattering lengths. Among other experimental data sets, Mughabghab’s set of resonance parameters were mainly evaluated on the basis of high-resolution neutron time of flight spectroscopy measurements performed at the Columbia University Nevis synchrocyclotron in the mid seventies [5].

Following the evaluation work performed by Lee in the mid 2000s [6] and the series of recent transmission and neutron capture yield measurements on enriched and natural Dy samples performed at the Gaertner LINear ACcelerator (LINAC) facility at Rensselaer Polytechnic Institute [7, 8], this work presents the newly evaluated set of neutron resonance parameters for total and capture cross sections on $^{156,158,160-164}\text{Dy}$ isotopes in the energy range of thermal energy up to a few keVs. The analysis of the experimental data sets uses the SAMMY [9] code, which performs a multi-level, multi-channel R -matrix fit to neutron data in the Reich–Moore (RM) formalism. Experimental conditions such as resolution function, finite sample size, nonuniform thickness, detector efficiencies and nuclide abundances of sample, multiple scattering, self-shielding, normalization, background, and Doppler broadening are taken into account. From both the capture and transmission data, we determined pertinent resonance parameters—and from these, the values of their systematics, such as strength functions, level spacings, and cumulative plots. In the thermal energy range, we utilized the *Atlas of Neutron Resonances* [10] as well as the coherent scattering lengths published by the National Institute of Standards and Technology (NIST) as a source of information on scattering and capture cross sections, as well as the resonance integral.

The report is organized as follows. The general conditions and evaluation methodology are outlined in Section 2, including two extensive subsections about the quantification of the external functions and thermal characteristics of the Dy isotopes; the experimental database is described in Section 3; the calculations and results of the R -matrix analysis are presented and discussed in Section 4, followed by a section on the uncertainty quantification. Statistical properties of the resonance parameters and their average values are examined in Section 5 and, for reproducibility purposes, a brief discussion on the repository generated to store and reproduce the results of this work is given in Section 4.1, followed by our conclusions and

recommendations in Section 7.

2. GENERAL CONDITIONS AND EVALUATION METHODOLOGY

2.1 DYSPROSIUM ISOTOPES

Naturally occurring dysprosium $^{\text{nat}}\text{Dy}$ consists of seven stable isotopes, $^{156,158,160-164}\text{Dy}$, plus an extremely long half-life isotope, ^{154}Dy ($t_{1/2} = 3 \times 10^6$ y). The even- A stable isotopes have spin-parity $I^\pi = 0^+$ that couples to s -wave neutrons to make compound nuclear states all characterized by the same spin parity $J^\pi = 1/2^+$, whereas the two odd- A stable isotopes, $^{161,163}\text{Dy}$ having $I^\pi = 5/2^+$ and $5/2^-$, respectively, couple to make two compound nuclear states $J^\pi = 2^+, 3^+$ and $J^\pi = 2^-, 3^-$. Table 1 shows four major Dy isotopes with a relatively homogeneous isotopic distribution. For each isotope, the upper energy limit E_{max} is based on the systematics of Liou’s data sets and corresponds to the energy regions where a large fraction of s -levels began to be experimentally missing. As can be seen, the resolved resonance range for odd- A isotopes extends over a smaller E_{max} than for even- A isotopes since the latter ones are naturally characterized by a larger level spacings. As reported in Table 1, the threshold energy for the inelastic channel of each isotope should be considered when determining the upper energy limit of the unresolved resonance region (URR). For each isotope, the RRR energy limit is well below that threshold energy; therefore, the total cross section is defined by only two energetically possible channels, elastic scattering and neutron capture. Although not planned in the list of nuclei relevant to NCSP, two minor isotopes, $^{156,158}\text{Dy}$, were included in the R -matrix analysis for completeness. This set of isotopes occurs in the mass region of the split $4s$ giant resonance in

Table 1. Dysprosium abundances (in %), upper energy limits for the RRR, and inelastic threshold energies (in keV).

No.	Isotope (I^π)	Abundance (%)	E_{max} (keV)	E_{inl} (keV)
1 (*)	^{156}Dy (0^+)	0.056	0.1	137.8
2 (*)	^{158}Dy (0^+)	0.095	0.1	98.9
3	^{160}Dy (0^+)	2.329	2.0	80.6
4	^{161}Dy ($5/2^+$)	18.889	1.0	25.6
5	^{162}Dy (0^+)	25.475	5.0	80.6
6	^{163}Dy ($5/2^-$)	24.896	1.0	73.4
7	^{164}Dy (0^+)	28.260	7.0	73.4

(*) Not planned by NCSP.

the S_0 strength function where the S_1 strength functions are expected to be smaller, so the observed level populations are nearly complete s -level populations, with little or no p -level contamination. In fact, with the current experimental data sets, only a few p -wave resonances were observed, and the evaluated set of resonance parameters is practically a pure single population of s levels.

2.2 QUANTIFICATION OF THE EXTERNAL FUNCTIONS

The R -matrix theory shows that the reaction cross sections analyzed over a limited energy range do not depend only on the “internal” levels but also on contributions from “external” levels below and above that energy range [11]. Therefore, before performing the R -matrix fit of the resonance parameters, it is important to quantify the contribution of the resonance analysis below the energy E_{min} (usually chosen as the zero energy) and above E_{max} , which is generally the upper energy limit where measured data can be still analyzed without losing statistical information on resonance levels. These two energies represent the energy limits of the analyzed neutron range, and the external contribution can be achieved quite conveniently by a resonance statistical representation of the external levels approximated by the tails of two very broad resonances with

equal strength [11]. Namely, for s -wave and $\bar{\Gamma}_\gamma \ll I$, this yields external levels with neutron widths

$$\Gamma_{n\pm}^J = \frac{3}{2} I S_0^J \sqrt{\frac{|E_{\pm}|}{1 \text{ eV}}}, \quad (1)$$

located symmetrically with respect to the mid-energy $\bar{E} = (E_{\max} + E_{\min})/2$ of the length of the interval $I = (E_{\max} - E_{\min})$ containing the local levels and calculated at energies $E_{\pm} = \bar{E} \pm \sqrt{3}I/2$. The value of the s -wave strength function S_0^J for different spin populations (two for odd- A isotopes such as $^{161,163}\text{Dy}$) can be obtained by the statistical analysis of the known widths of the resonance parameters. In Table 2, the values of the neutron widths and related energy levels calculated with Eq. (1) are reported along with the neutron s -wave strength functions obtained from ENDF/B-VIII.0 resonance parameters.

Table 2. External function parameters calculated with Eq. (1) including the strength functions S_0^J (in 10^{-4} units), the corresponding pole strength s_0^J (in 10^{-2} units)*, the length interval I (in keV) and related resonance widths ($\Gamma_{n\pm}$ in meV) for each channel spin J^π .

Isotope (I^π)	I	E_-	Γ_{n-}^J	E_+	Γ_{n+}^J	S_0^J	s_0^J	J^π
$^{160}\text{Dy} (0^+)$	2.0	-732	15909	2732	30734	1.96	6.01	$1/2^+$
$^{161}\text{Dy} (5/2^+)$	1.0	-366	2739	1366	5291	0.95	2.92	2^+
			2304		4452	0.80	2.46	3^+
$^{162}\text{Dy} (0^+)$	5.0	-1793	66924	6693	129286	2.15	6.58	$1/2^+$
$^{163}\text{Dy} (5/2^-)$	1.0	-366	3415	1366	6597	1.19	3.65	2^-
			2447		4727	0.85	2.61	3^-
$^{164}\text{Dy} (0^+)$	7.0	-2562	70688	9562	136559	1.33	4.07	$1/2^+$

(*) The strength function and the pole strength are related by the channel radius a_c and wave number $k_c \equiv k_c(E)$ for the channel c as $S_c^J = 2a_c k_c s_c^J \sqrt{1 \text{ eV}/E}$.

The R -matrix external level representation obtained by two broad resonances with equal strength is quite convenient; however, it is an approximation to the R -matrix function

$$R_{cc}(E) - R_c^\infty = 2s_c \left\{ \tanh^{-1} \left[\frac{E - \bar{E}}{I/2} \right] + i \frac{\bar{\Gamma}_\gamma}{I} \left[1 - \left(\frac{E - \bar{E}}{I/2} \right)^2 \right]^{-1} \right\} \simeq \frac{\gamma_{n-}^2}{E_- - E - i\bar{\Gamma}_\gamma/2} + \frac{\gamma_{n+}^2}{E_+ - E - i\bar{\Gamma}_\gamma/2}, \quad (2)$$

where $\bar{\Gamma}_\gamma$ is the average capture width, R_c^∞ is the distant-level parameter, and the reduced amplitude is related to resonance neutron widths as $\Gamma_{n\pm}^c = 2P_\ell(E_{\pm})\gamma_{cn}^2$ as defined in Eq. (1). Although the differences between the two representations are usually small over most of the energy range, the differences are not negligible towards the edges. To correct for the edge effects of the approximation of Eq. (2), it is necessary to include additional bound and external levels to those of Table 2. Starting from the resonance parameters of Table 2 plus one additional bound and external level for each channel group, a fit of the elastic cross sections calculated from the R -function $R_{cc'}(E) \equiv R_{cc'}$ was performed over the energy interval I for each isotope. The fit of the calculated elastic cross sections was performed by varying the neutron widths and keeping unchanged the average capture widths. The use of the elastic cross sections was necessary because the SAMMY code cannot account for the complex part of the R -function $R_{cc'}$. In doing so, one assumes that the correlation between the elastic and capture reaction channel is negligible.

This procedure yields the set of resonance parameters of Table 3 and the results of the fit, specifically for the ^{160}Dy isotope, showing the inclusion of the edge effects in the external function; these additional bound and external levels are plotted in Fig. 1. The edge effects of the elastic cross sections calculated by the SAMMY code using the R -function $R_{cc'}$ (in purple) are well reproduced by the cross sections (in light blue), reconstructed from the set of resonance parameters reported in Table 3. The results from the use of the broad levels defined in Eq. (1) are shown in green. One should note that the parameters reported in Table 3

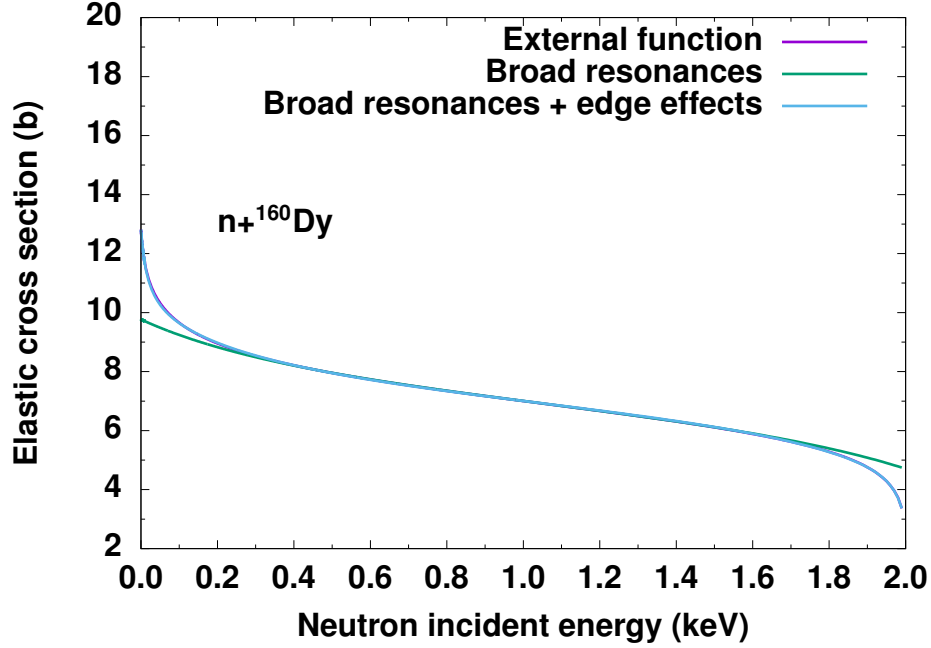


Figure 1. $n+^{160}\text{Dy}$ elastic cross sections showing the inclusion of the edge effects. The elastic cross sections calculated by the SAMMY code using the R -function $R_{cc'}$ are shown in purple. In light blue and green are the results of the elastic cross sections reconstructed from the resonance parameters with or without edge effects, respectively.

represent the initial set of external parameters. This is usually modified following the fit of the complete set of measured data and, most importantly, the calibration of the thermal values for scattering, capture, and fission (only for fissile nuclei) reaction channels.

2.3 DYSPROSIUM THERMAL CHARACTERISTICS

The calibration of the negative (or bound) energy levels and related resonance widths was achieved by fitting Block's transmission and capture data [7] measured on $^{\text{nat}}\text{Dy}$ samples together with the reference thermal characteristics of each isotope reported by Sears [12], which are also available on the NIST website. As shown in Fig. 3, Block's measured capture and transmission data nicely extend to the thermal energy below 1 eV and, together with Sears' thermal constants, can be excellent constraints in defining the resonance parameters—particularly for negative energies. In doing so, an important aspect of the fitting procedure is the inclusion of relevant experimental corrections related to the detector deficiencies for the capture yield measurement, as these corrections are particularly sensitive to the low-energy tail of the measured data. These are described in detail in Ref. [7] and used in SAMMY inputs as the scaling factor for each isotope.

Table 3. Resonance parameter for initial external function quantification with edge effects. Resonance parameters (in meV) and resonance energies (in eV) obtained by the fit of the elastic cross sections calculated from the R -function $R_{cc'}$ of Eq. (2).

Isotope (I^π)	E_{\min}	E_{\max}	E	Γ_n^J	$\bar{\Gamma}_\gamma$	J^π
^{160}Dy (0^+)	0.0	2.0	-565.21	10415.69	110	$1/2^+$
			-19.38	46.71		
			2023.03	662.08		
			2587.66	22293.35		
^{161}Dy ($5/2^+$)	0.0	1.0	-358.09	2455.13	120	2^+
			-12.71	23.96		
			1256.15	393.62		
			1347.29	4617.17		
			-361.69	1980.77	120	3^+
			-106.89	43.90		
			1009.96	134.35		
			1345.69	3736.60		
^{162}Dy (0^+)	0.0	5.0	-1294.41	38944.84	120	$1/2^+$
			-62.63	45.16		
			5032.44	1784.12		
			6253.58	86419.92		
^{163}Dy ($5/2^-$)	0.0	1.0	-410.60	2393.04	120	2^-
			1010.92	292.77		
			1349.55	6956.22		
			-382.47	3078.90	120	3^-
			-13.61	29.34		
			1276.95	73.73		
			1359.49	3549.83		
^{164}Dy (0^+)	0.0	7.0	-2036.79	46723.60	100	$1/2^+$
			-99.98	306.77		
			7102.11	1132.59		
			9015.64	98607.11		

This work started from the initial set of the external levels of Table 3 to find the current evaluated resonance parameters for the external energy levels and related channel widths of each isotope, as reported in Table 4. Highlighted in blue are those (negative) levels added to improve the fit of the measured data including thermal constants. Although thermal scattering cross sections and scattering lengths are intimately related,

Table 4. Resonance parameters for final external function quantification. Highlighted in blue are the energy levels added to improve the fit of the measured data including thermal constants. Units are the same as those in Table 2.

Isotope (I^π)	E_{\min}	E_{\max}	E	Γ_n^J	$\bar{\Gamma}_\gamma$	J^π
^{160}Dy (0^+)	0.0	2.0	-601.81	9545.25	103	$1/2^+$
			-51.98	326.19	99	
			-14.30	9.05	86	
			2023.00	662.62	110	
			2586.10	22795.48	110	
^{161}Dy ($5/2^+$)	0.0	1.0	-360.67	2454.73	120	2^+
			-19.83	29.88	120	
			1256.13	393.66	120	
			1347.11	4621.42	120	
			-360.67	2453.73	120	3^+
			-110.44	44.55	120	
			-2.50	15.37	150	
			1009.92	134.37	120	
^{162}Dy (0^+)	0.0	5.0	1345.10	3747.88	120	$1/2^+$
			-1255.35	13241.10	117	
			-47.40	38.58	111	
			5032.50	1779.22	120	
^{163}Dy ($5/2^-$)	0.0	1.0	6253.58	78659.40	120	2^-
			-552.74	4151.67	120	
			1007.24	294.93	120	
			1296.58	8631.93	120	3^-
			-584.20	1513.45	120	
			-4.65	3.20	120	
			1275.23	74.83	120	
^{164}Dy (0^+)	0.0	7.0	1275.50	4883.50	120	$1/2^+$
			-2037.05	46629.74	83	
			-102.50	310.25	83	
			-1.88	50.06	63	
			7102.11	1132.61	100	
^{164}Dy (0^+)	0.0	7.0	9015.64	98653.65	100	$1/2^+$

the latter should be always checked to ensure proper calibration with available measured data. In fact, even when the thermal cross sections are satisfactory, the scattering lengths may need additional calibration. Although the SAMMY code does not currently have the capability to fit the scattering length as measured data, this work attempted to calibrate both scattering lengths and scattering cross sections. The evaluated bound (in)coherent scattering lengths and thermal capture cross sections are reported in Table 5 together with the values of complex bound scattering lengths and related radii. In this evaluation work, the scattering or potential radius R' and the effective radius a_c that is usually used to calculate the penetration factors were kept the same for all channels of each isotope.

One can find reasonable agreement with the values reported by the NIST compilation except for the cases of the $^{161,164}\text{Dy}$ isotopes. For the most abundant isotope ^{164}Dy , there is a discrepancy of 200 b (about -7%) noticeable for the absorption cross section and 24 b (about $+8\%$) for the scattering cross section leading to a reduction of the thermal value for the total cross section of about 170 b (about -5%). The evaluated thermal capture value was found by fitting the recently measured activation data [13]. For the scattering channel, only a few measured data are available, and the NIST value differs by about 24 b with the ENDF/VIII.0 library. The value of about 331 b evaluated in this work was calibrated and constrained by the simultaneous fit of transmission and capture data for $^{\text{nat}}\text{Dy}$ together with thermal capture and scattering constants for six other isotopes. The choice of a larger ^{164}Dy thermal scattering value than that reported by the NIST compilation was also motivated by recent and accurate capture activation measurements. In fact, the ^{164}Dy scattering thermal cross section obtained by the difference between the total cross section, $\sigma_{\text{tot}} = 3000_{-45}^{+70}$ b, measured at thermal by Vertebny [14, 15], and the thermal capture cross section, $\sigma_{\gamma} = 2650 \pm 25$ b, measured by Farina [13], yields $\sigma_s = 350_{-37}^{+65}$ b. From this, and assuming Farina's recent measured data set is reliable, the NIST compilation value for the ^{164}Dy scattering cross sections is clearly found outside the considerably large derived uncertainty range. On a different note related to the 4th most abundant isotope, the scattering cross section of the ^{161}Dy isotope is still 2σ within the NIST compilation and was adopted from the ENDF/B-VIII.0 library based on the goodness of the fit of the natural data. This choice is also consistent with the evaluation work by Lee [6].

The following paragraphs report the details of the derivation for the equations used to calculate the thermal constant values reported in Table 5. For the scattering lengths, a convenient way to show their

Table 5. (In)coherent bound scattering^(*) and thermal absorption cross sections (in barn) with corresponding bound (in)coherent scattering lengths^() (in fm) compared to the values reported by NIST^(***). The scattering radius a_c (in fm) for each isotope is also reported.**

	a_c		σ_{γ}	σ_s	σ_{coh}	σ_{incoh}	b_{coh}	b_{incoh}	b_+	b_-
^{156}Dy	7.50	Present	33	4.1	4.1	0.0	5.7	0.0	n/a	n/a
		NIST	33	4.7	4.7	0.0	6.1	0.0	n/a	n/a
^{158}Dy	7.40	Present	43.1	6.5	6.5	0.0	7.2	0.0	n/a	n/a
		NIST	43.0	5.0	5.0	0.0	6.0	0.0	n/a	n/a
^{160}Dy	7.46	Present	55	5.6	5.6	0.0	6.7	0.0	n/a	n/a
		NIST	56	5.6	5.6	0.0	6.7	0.0	n/a	n/a
^{161}Dy	7.47	Present	605	17.5	14.5	3.0	10.7- i 0.17	4.9- i 0.12	8.71- i 0.16	2.08
		NIST	600	16.0	13.3	3.0	10.3	4.9	n/a	n/a
^{162}Dy	5.90	Present	195	0.28	0.28	0.0	-1.50	0.0	n/a	n/a
		NIST	194	0.25	0.25	0.0	-1.41	0.0	n/a	n/a
^{163}Dy	7.50	Present	125	3.2	3.2	0.27	4.9	1.3	3.5	1.4
		NIST	124	3.3	3.1	0.21	5.0	1.3	n/a	n/a
^{164}Dy	7.51	Present	2650	327	327	0.0	51.3- i 0.74	0.0	n/a	n/a
		NIST	2840	307	307	0.0	49.7- i 0.79	0.0	n/a	n/a

^(*)By definition, the scattering and coherent cross section are identical for nuclei with spin $I = 0$.

^(**)The imaginary component of the calculated bound scattering lengths b is reported for values >0.1 .

^(***)As discussed in subsection 2.2, the bound scattering lengths and related cross sections can be converted to free quantities by the factor $(A + 1)/A$ which, for the Dy isotopes, introduces a difference of about 0.6%. In this work, the contribution of the neutron-electron interaction was not included.

relationship to R -matrix parameters is to expand the collision matrix in the low energy limit. The theoretical

values of the reported scattering lengths were calculated on the basis of the collision matrix U_{cc}^J for the elastic channel ($c=c'$) computed from the set of the evaluated resonance parameters and scattering radius R' for each Dy isotope. In the reduced R -matrix approximation, the RM collision function for each s -wave incident neutron channel $J=(I \pm 1/2)^\pi$ (i.e., one for $I=0$, two for $I \neq 0$);

$$U_{cc}^J = e^{-2ik_c a_c} [1 - ik_c a_c R_{cc}^J]^{-1} [1 + ik_c a_c R_{cc}^J], \quad (3)$$

where the quantity a_c is the radius for the incident channel c (usually chosen to be equal to the potential scattering radius R'), and $k_c = \alpha_c \sqrt{E}$ is the wave number defined for the incident (neutron) channel, namely, for $\alpha_c = 2.1968 \times 10^{-4} A_c / (A_c + 1)$ (in $\text{fm}^{-1} \text{eV}^{-1/2}$): E is expressed in eV, and α_c is the nucleus–neutron mass ratio for particle pair of the incident channel. The reduced- R complex function for this one-channel case is

$$R_{cc}^J = \sum_{\lambda} \frac{\gamma_{\lambda(J,c=n)}^2}{E_{\lambda} - E - i\Gamma_{\lambda\gamma}/2}, \quad (4)$$

where the summation extends over the nuclear energy (positive and negative) levels E_{λ} related to the formal energy states of the nuclear compound system with reduced neutron width amplitude $\gamma_{\lambda n}$ and capture widths $\Gamma_{\lambda\gamma}$. Omitting the index c being the quantities corresponding to the (neutron) elastic scattering, the collision function in the low-energy limit, $x=ka \rightarrow 0^+$, can be written as

$$U^J = 1 + 2ix(R^J - 1) - 2x^2(R^J - 1)^2 + o(x^3). \quad (5)$$

Comparing the detailed and usual form of the elastic scattering wave function for s -wave neutron interaction, one can obtain the relation between the scattering wave amplitude and the RM collision function,

$$F^J = ia(1 - U^J)/(2x), \quad (6)$$

which is related to the free scattering length for the channel spin J as

$$a^J = -\lim_{x \rightarrow 0^+} F^J \approx a[1 - \Re(R^J) - i\Im(R^J)] = a(1 - R^J). \quad (7)$$

For nuclei with target spin $I = 0$, the free coherent scattering length is the quantity defined in Eq. (7); however, for nuclei with target spin $I \neq 0$, the free coherent scattering length

$$a_{\text{coh}} = \begin{cases} a^J & I = 0 \\ g_+ a^{J_+} + g_- a^{J_-} & I \neq 0 \end{cases}$$

is the sum of the partial scattering lengths weighted by their spin statistical factors $g_+ = (I + 1)/(2I + 1)$, $g_- = I/(2I + 1)$ related to $J_+ = I + 1/2$ and $J_- = |I - 1/2|$, respectively. The same quantities are also used to compute the incoherent scattering length, defined as

$$a_{\text{incoh}} = \begin{cases} 0 & I = 0 \\ \sqrt{g_+ g_-} (a^{J_+} - a^{J_-}) & I \neq 0. \end{cases}$$

Since the measured or recommended data are usually reported as bound (in)coherent scattering length $b_{(\text{in})\text{coh}}$, the relation used in this work to convert the measured quantities to free (in)coherent scattering lengths is $a_{(\text{in})\text{coh}} = b_{(\text{in})\text{coh}} A/(A + 1)$, from which one can calculate the free scattering cross section

$$\sigma_s = 4\pi \begin{cases} |a^J|^2 & I = 0 \\ g_+ |a^{J_+}|^2 + g_- |a^{J_-}|^2 & I \neq 0, \end{cases}$$

where, for $I \neq 0$, the result is given by the unitary condition of the spin statistical factors, $g_+ + g_- = 1$, and the sum of the coherent and incoherent cross sections,

$$\sigma_{\text{coh}} = 4\pi |g_+ a^{J_+} + g_- a^{J_-}|^2 \quad \text{and} \quad \sigma_{\text{incoh}} = 4\pi g_+ g_- |a^{J_+} - a^{J_-}|^2. \quad (8)$$

In the case of the Dy isotopes, the absorption cross section is given by the capture reaction channel only, and, as reported in Table 5, the value calculated from the free scattering length is defined by

$$\sigma_\gamma = \frac{4\pi}{k} \begin{cases} |\Im(a^J)| & I = 0 \\ g_+ |\Im(a^{J_+})| + g_- |\Im(a^{J_-})| & I \neq 0, \end{cases}$$

where the wave number k is calculated at the thermal neutron energy $E_{\text{th}} = 0.0253$ eV.

3. THE EXPERIMENTAL DATABASE

To improve the accuracy of the Dy evaluations in both thermal and resolved resonance neutron energy ranges, a comprehensive experimental database is needed to evaluate the set of resonance parameters. As schematically reported in Table 6, one can distinguish two major experimental campaigns for the Dy isotopes. Chronologically, the first one was performed by Liou [5] at the Columbia University Nevis synchro-cyclotron. This included transmission and capture measurements of $^{160-164}\text{Dy}$ isotopes on enriched Dy_2O_3 samples of different thicknesses focusing on epithermal incident neutron energy range, that is, above 15 eV and up to several keVs. The second experimental campaign was generated within the needs of the NCSP and performed at the Gaertner LINAC Center at RPI over thermal and epithermal ranges. Among these recent measurements, neutron capture yields of $^{161-164}\text{Dy}$ isotopes were measured in the energy range from 10 eV up to 1 keV. The capture yield measurements were performed on isotopically-enriched dysprosium metallic samples using the Time-of-Flight (ToF) method with a 16 segment sodium iodide multiplicity detector [8]. Within the same campaign, an additional series of capture and transmission measurements was performed on $^{\text{nat}}\text{Dy}$ samples in the thermal neutron energy region up to 20 eV [7]. More details on both series of measurements can be found in Refs. [7, 8]. Among the neutron transmission experimental data available for Dy isotopes, Liou's measured data represent a relevant portion because of their extended energy range and sample isotopic enrichment. In fact, Liou's transmission data are currently the only source of information to evaluate the neutron widths of the Dy isotopes in the energy range above 15 eV. Unfortunately, the capture data for this set of measurements are missing, and there is no uncertainty quantification for the transmission data reported in the EXFOR library. Moreover, the large number and magnitude of negative values found in the reported $^{160,163,164}\text{Dy}$ total measured cross sections implies an over-correction of the background contribution. Among possible background corrections, one might have been applied to eliminate the contribution of the oxygen cross sections due to the use of oxide samples Dy_2O_3 in the measurements. In this regard, it is unclear whether the number of atoms/barn reported is related to the specific enriched isotope or to the oxide sample: this affected our ability to correctly calculate the total number of atoms/barn of the samples. Moreover, especially in the low-energy region above 15 eV, the energy spectrum of the Dy isotopes contains several energy levels for which the neutron absorption is maximum. For these levels, Liou's data were reported as saturated resonances because of the choice of the sample thickness, and an accurate fit of the resonance widths was not possible. At least for the resonance levels large enough to be visible with measurements with $^{\text{nat}}\text{Dy}$ sample, this problem was partially resolved by Shin's transmission data.

Because Liou's transmission data were reported as total cross sections for the whole set of Dy isotopes, the conversion to transmission data was necessary to convolute in the R -matrix analysis the experimental

Table 6. Selected measurements and related sample configurations such as thickness n (a/b), path length L , and enrichment used in the evaluation of the resonance parameter of Dy isotopes.

Author (year)	Facility	Sample	Type	n (10^{-3} a/b)	Energy range	Path length L (m)
Liou (1975)	NEVIS	^{160}Dy	Trans [*]	0.881	15 eV–2 keV	202.05
		^{161}Dy		1.644, 7.451	15 eV–1 keV	
		^{162}Dy		1.533, 6.934	15 eV–5 keV	
		^{163}Dy		1.403	15 eV–1 keV	
		^{164}Dy		10.54, 2.337	15 eV–7 keV	
Shin (2017)	RPI	^{161}Dy	Yield	0.6359	12 eV–1 keV	25.57
		^{162}Dy		0.6445	10 eV–1 keV	
		^{163}Dy		0.6503	10 eV–1 keV	
		^{164}Dy		0.6196	10 eV–1 keV	
Block (2017)	RPI	$^{\text{nat}}\text{Dy}$	Trans	1.610	0.01–20 eV	14.97
		$^{\text{nat}}\text{Dy}$	Yield	0.8050	0.01–20 eV	25.44
		$^{\text{nat}}\text{Dy}$	Trans	0.7840	4 eV–2 keV	25.60
		$^{\text{nat}}\text{Dy}$	Yield	1.6304	10 eV–1 keV	25.57

(*) Converted to transmission data from the reported total cross sections.

configuration of Liou’s measurements. In this regard, the experimental resolution and Doppler broadening parameters were also needed as input parameters for an accurate description of the shape of the experimental data. For Liou’s data, only partial information on the set of experimental resolution parameters was found. This consisted of the minimum channel width of $\Delta_G = 40$ ns and the flight path length $L = 202.5$ m. With this information together with the relation between the energy and the speed of the neutron of mass m ,

$$t(E) = \sqrt{\frac{m}{2E}} L, \quad (9)$$

the values of the crunch data table were directly extracted from Liou’s data sets to obtain the energy dependence of the experimental resolution. As shown in Fig. 2 for ^{160}Dy measured data, the number of the energy boundaries $B_{\Delta_G}(E)$ and their related values are clearly distinguishable.

Although the crunch tables were correctly deduced from the data, the impact of the resolution broadening on the transmission data reconstructed including the experimental resolution parameters was not particularly evident. This might suggest that the reported measured data are not exactly the raw measured data but total cross sections reconstructed from the resonance parameters. This would explain the lack of experimental uncertainty analysis in the EXFOR library.

The recent measurements performed on $^{\text{nat}}\text{Dy}$ samples at the RPI facility for both transmission and capture yield data supplied additional and relevant information in evaluating the low-energy data in the thermal range and resonance levels up to 20 eV as shown in Fig. 3. However, the natural data measured in the epithermal range up to 2 keV could not match the neutron resonance level information of Liou’s isotopically-enriched measurements. In this regard, considerable spectroscopic information obtained from Liou’s measurements supported RPI capture yield measurements performed on isotopically-enriched samples.

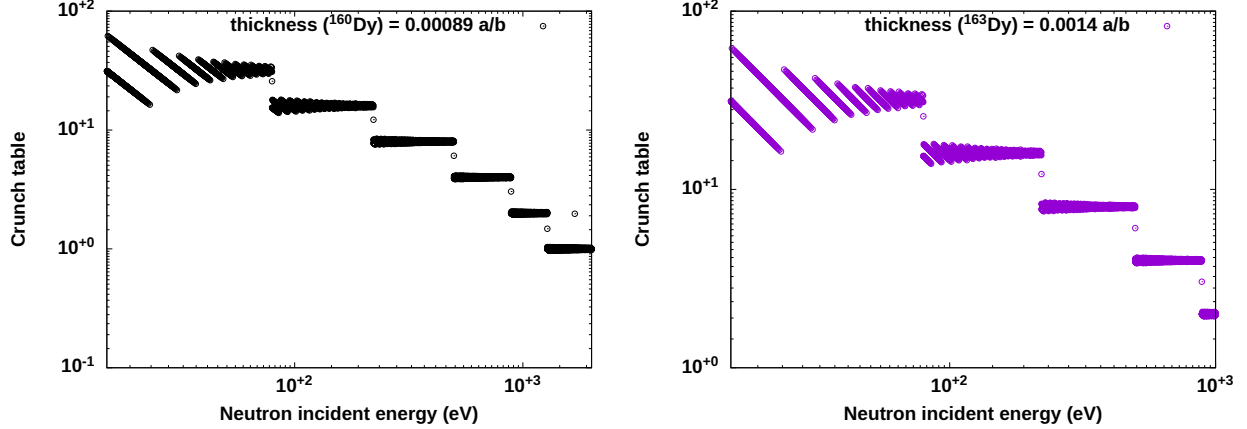


Figure 2. Energy dependent values of the crunch table for a channel width $\Delta_G=40$ ns as a function of the incident neutron energy for the even- A isotope ^{160}Dy and odd- A isotopes ^{163}Dy .

4. RESULTS OF THE R-MATRIX ANALYSIS

The set of RM resonance parameters obtained by the multi-isotopic and multi-channel SAMMY R -matrix analysis of Dy experimental data in the energy range from thermal to several keVs contains 664 energy levels, 634 in the analyzed energy range, and 30 external resonances including 17 negative levels. The set of resonance parameters for the natural sample is given in the appendix, and most of the energy levels were considered as induced by s -wave neutrons. Only a few p -wave resonances were included for two even- A isotopes $^{162,164}\text{Dy}$. In deriving the evaluated resonance parameters, particular care was devoted to the thermal energy region up to 20 eV. Here, the set of resonance parameters was constrained by the thermal values recommended by NIST (with the exception of ^{164}Dy) and by Block's transmission and capture measurements that nicely extended in the low-energy range up to 0.01 eV. In addition to the negative levels defined in Section 2.2, the fitting procedure started from the combination of the set of resonance parameters reported in the ENDF/B-VIII.0 library and Shin's paper [8]. In the latter work, 46 new energy levels (29 for ^{161}Dy and 17 for ^{163}Dy) were reported and adopted in this work, although their existence can be sometimes debatable due to the lower and lower statistics of both measured Shin's capture and Liou's transmission data for increasing incident neutron energy. Moreover, in Shin's work, 12 energy levels (six for ^{161}Dy , two for ^{163}Dy , and four for ^{164}Dy) listed in the ENDF/B-VIII.0 library were not observed. These updates were also adopted by this work with the exception of the energy level at 12.7 eV for the ^{161}Dy isotope. This level was reported and confirmed by Block's transmission measurements [7] and not observed in Shin's capture measurements most likely because of the poor statistics of the measurements in this energy range.

As shown in Fig. 3, the capture yield (left) and transmission (right) measurements on $^{\text{nat}}\text{Dy}$ samples were invaluable to evaluate the thermal neutron energy region and the first few resonances up to about 20 eV. The energy-dependent residuals (bottom) are overall well within a reasonable acceptance range except below 0.1 eV. The large deviation is due to the non-physical discontinuity in the experimental uncertainty (top) visible at about 0.1 eV. This effect occurs in both transmission and capture measurements approaching a minimal uncertainty $\lesssim 1\%$ and close to 0.1%, respectively. Although generally unremarkable, other discontinuities are visible in the energy regions at about 4 eV for the capture yield data and about 5 eV for the transmission data. Despite the fact that such discontinuities in the experimental uncertainty at about

0.1 eV and 4–5 eV are related to the crunch tables of the measured data, these generate residual variations up to about 10-sigmas and mainly below 0.1 eV. These large residuals are clearly not possible to address due to the extremely low uncertainty coupled with the constrained behavior ($\propto 1/\sqrt{E}$ and $\propto 1/\sqrt{E} + \text{const}$ for the capture and total reaction channel, respectively) of the evaluated data in the thermal region. In this regard, the experimental configuration of Block's and Shin's capture measured data depends on the detector efficiencies [7, 8] that are taken into account in the SAMMY *R*-matrix analysis as a multiplicative factor defined for each isotope, accordingly scaling the calculated neutron capture yield. As expected, these factors are very impactful in describing the low-energy tail of the measured data, and, in this work, they were considered constants. In the epithermal energy region, the energy-dependent residuals of the capture yield

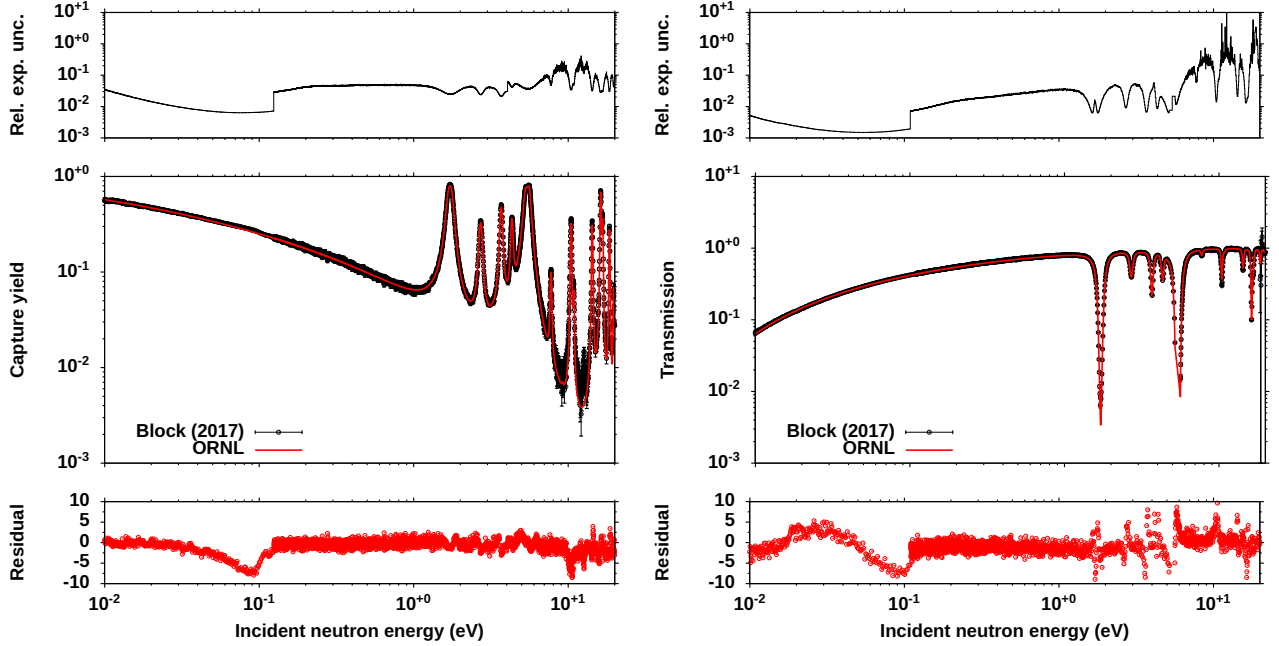


Figure 3. $n + {}^{\text{nat}}\text{Dy}$ capture yield (left) and transmission data (right) measured in thermal energy range at the RPI facility (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

and transmission data shown at the bottom of Fig. 4 are on average within 2-sigmas variation range. This is facilitated also by the experimental uncertainty (top) that is larger than previous data measured in the thermal energy range. To avoid clutter in the plotted data, the transmission data measured in the energy region between 300–400 eV where one of the filters' contribution was visible were excluded. Since these measurements performed on ${}^{\text{nat}}\text{Dy}$ samples, only the most relevant resonances were measurable. In this regard, Liou's transmission data [5] as well as Shin's capture yield data [8] measured on enriched samples allow a detailed analysis of the resonance levels for each isotope, as shown in the following figures. As explained in Section 3, Liou's measured data were converted from total cross section to transmission data and, because of no experimental uncertainty was reported, an estimated uncertainty of 1% was assigned to the total cross sections. The transmission data uncertainty was, therefore, based on total cross section estimated uncertainty through the relation $T(E) = e^{-n\sigma(E)}$ together with the corresponding energy-dependent partial derivatives. Clearly, this procedure generates inconsistent uncertainty for both negative total cross sections reported in EXFOR and large saturated resonance levels where the transmission data are nearly zero, and

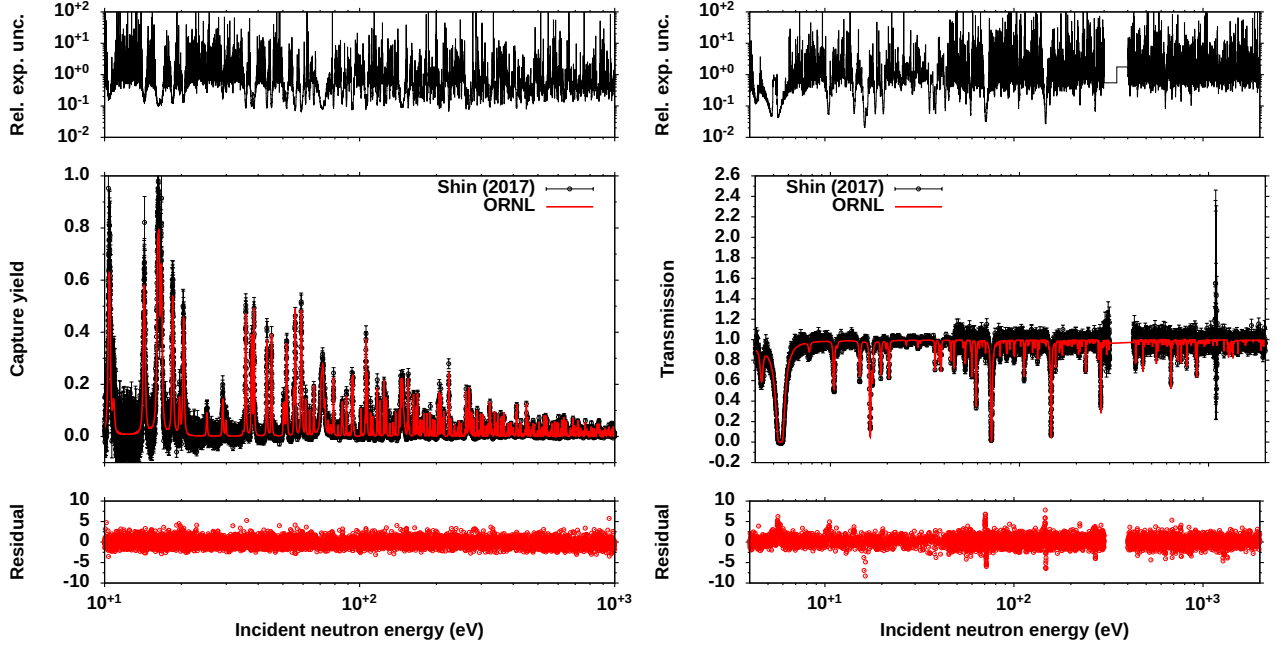


Figure 4. $n + {}^{\text{nat}}\text{Dy}$ capture yield (left) and transmission data (right) measured in epithermal energy range at the RPI facility (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

uncertainty could thus be in the range of negative values. Moreover, because the experimental standard deviation is inversely proportional to the square root of the number of events, the assignment of a constant uncertainty does not reflect the fact that in energy regions where large resonance levels are measured, the number of events is considerably higher than in other energy regions (e.g., the valleys). For instance, an expected energy-dependent behavior is clearly shown in the RPI measurements such as Fig. 4, in which a minimum in the capture yield uncertainty corresponds to the peak of the resonance level. In the same figure, the transmission data show a similar behavior—keeping in mind that, by their definition, the minima correspond to the resonance peaks in the total cross sections. For these reasons, Liou’s transmission data were not included in the optimization procedure and were used only for their detailed spectroscopic information. In Figs. 5,6,8,10,11, Liou’s transmission measured data (in black) with the estimated uncertainties (top) and corresponding residuals (bottom) are plotted together with the calculated data (in red)*.

4.1 UNCERTAINTY QUANTIFICATION

The uncertainty quantification of the evaluated resonance parameters was performed by the Bayesian procedure implemented in the SAMMY code. A sequential Bayesian fit including all (negative and positive) energy levels and related resonance widths for seven isotopes was performed over the set of measured data reported in Table 6 as well as thermal values. The retroactive scheme option was then applied to approximate the covariance matrix obtained in the previous step by assuming the resonance parameters are not very different from one step to another. The generation of a full resonance parameter covariance (RPC) matrix, including the isotope–isotope correlations, would have been the preferable solution. However, due to the

*In these plots, the data associated with negative reported experimental values are eliminated.

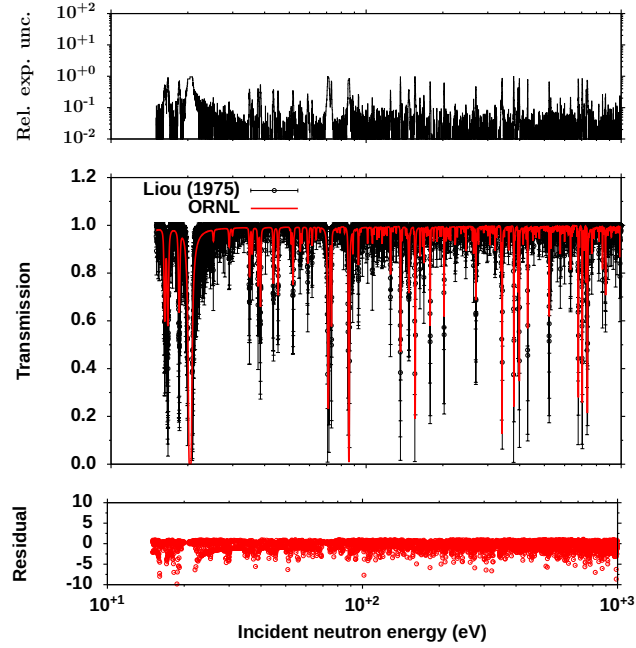


Figure 5. $n+^{160}\text{Dy}$ Liou's transmission data measured in the neutron energy range up to 1 keV (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

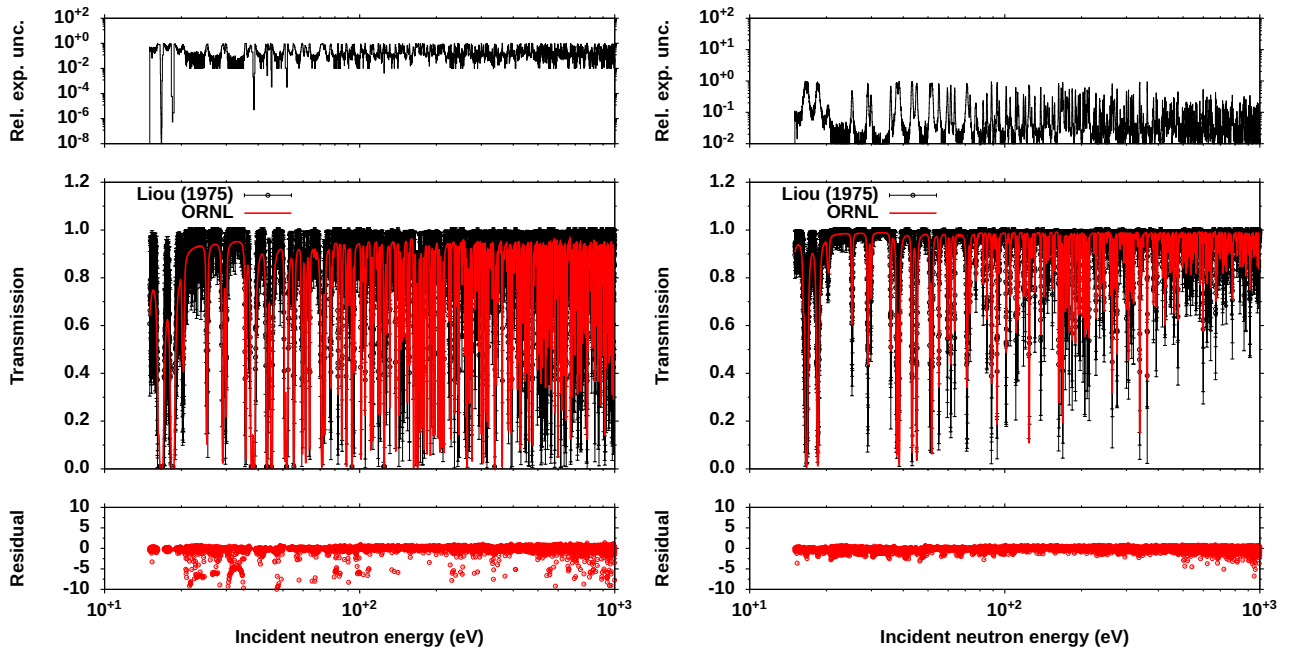


Figure 6. $n+^{161}\text{Dy}$ Liou's thick (left) and thin (right) transmission data measured in the neutron energy range up to 1 keV (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

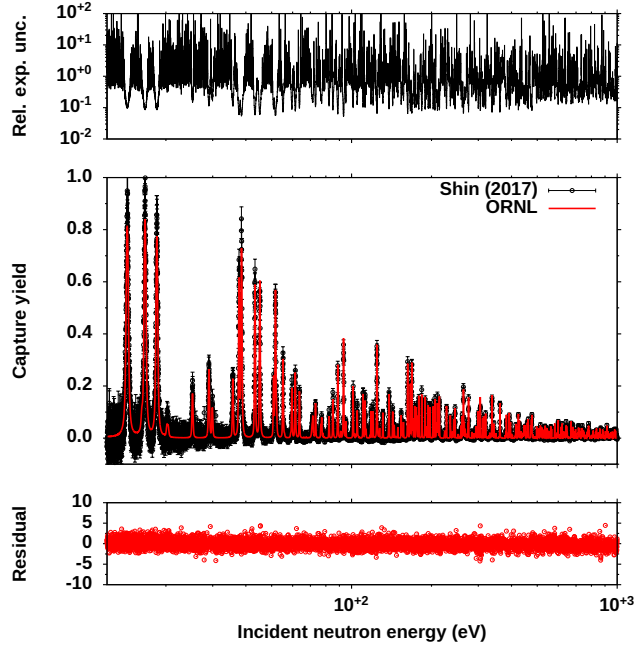


Figure 7. $n+^{161}\text{Dy}$ capture yield measured in epithermal energy range at the RPI facility (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

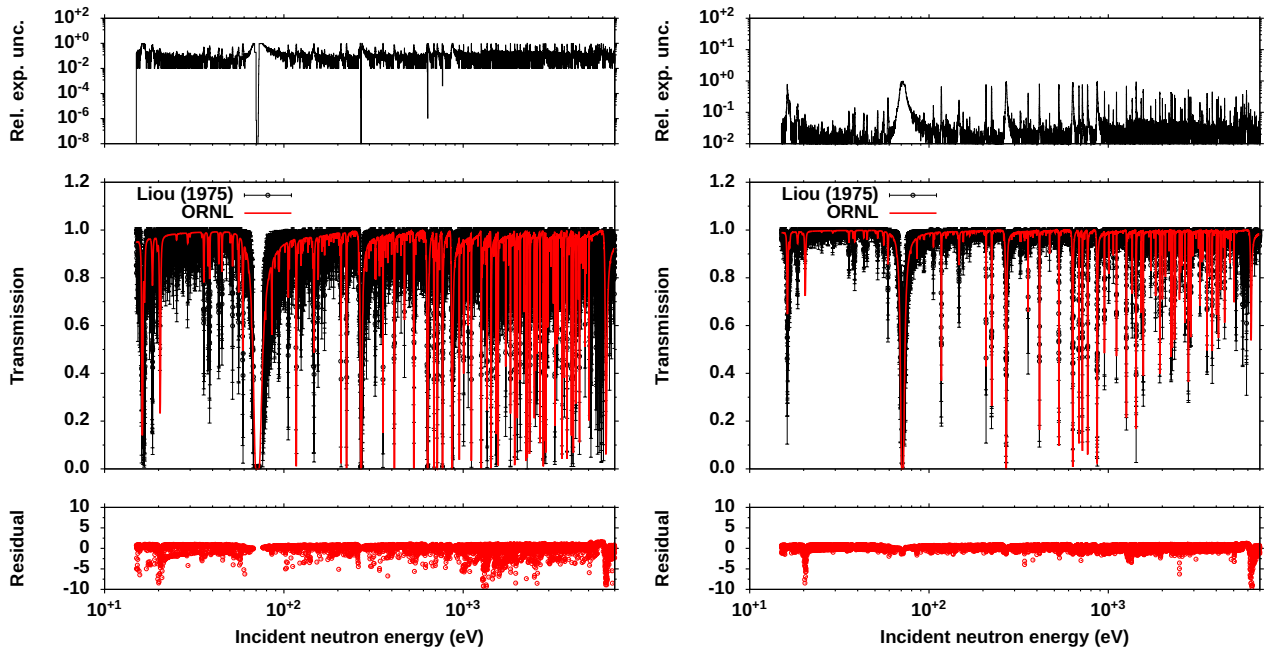


Figure 8. $n+^{162}\text{Dy}$ Liou's thick (left) and thin (right) transmission data measured in the neutron energy range up to 1 keV (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

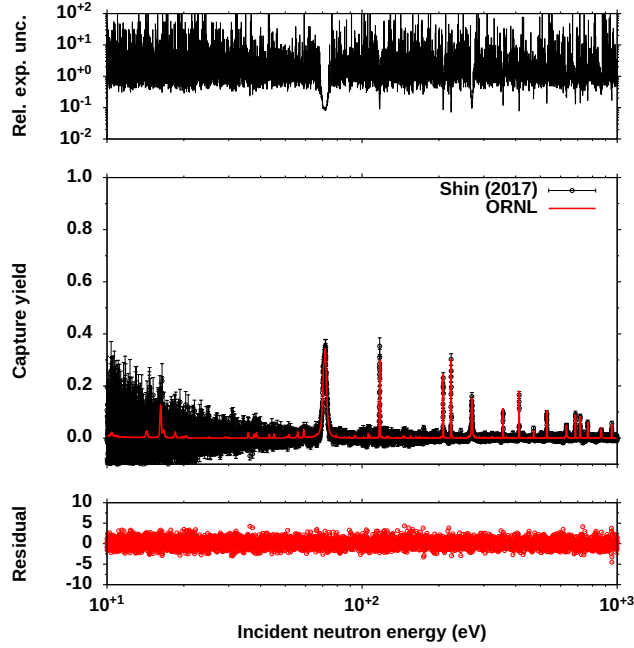


Figure 9. $n+^{162}\text{Dy}$ capture yield measured in epithermal energy range at the RPI facility (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

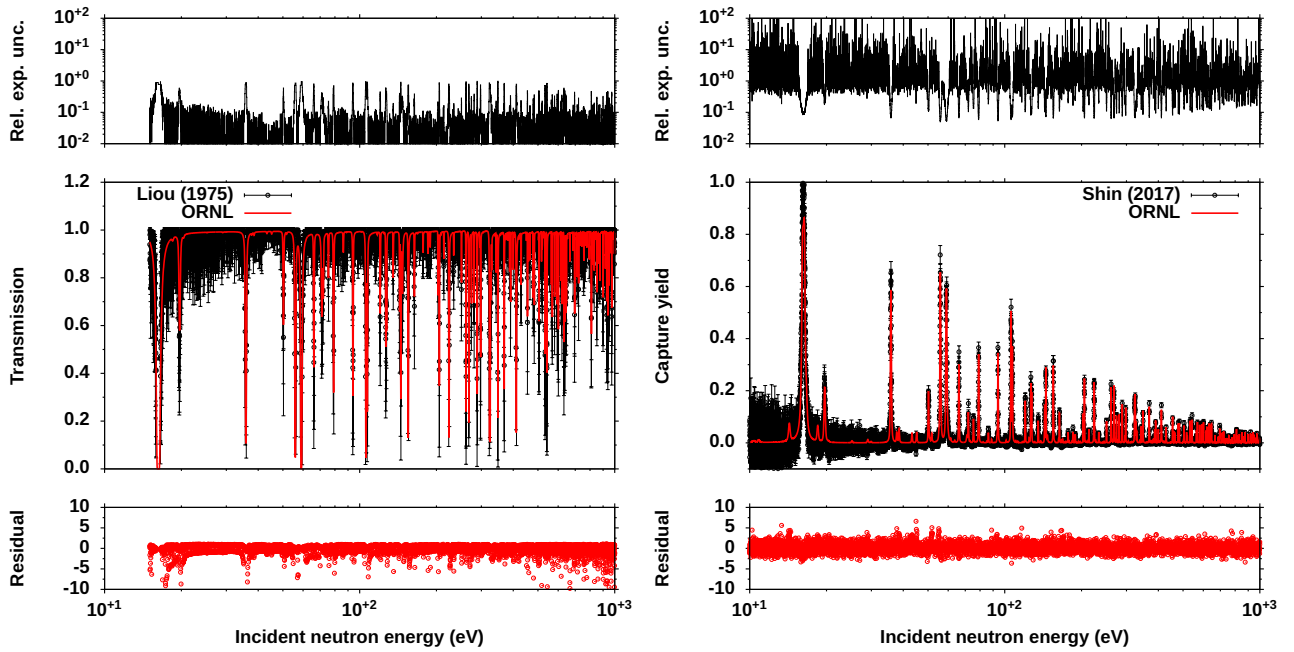


Figure 10. $n+^{163}\text{Dy}$ Liou's transmission and Shin's capture data measured in the neutron energy range up to 1 keV (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

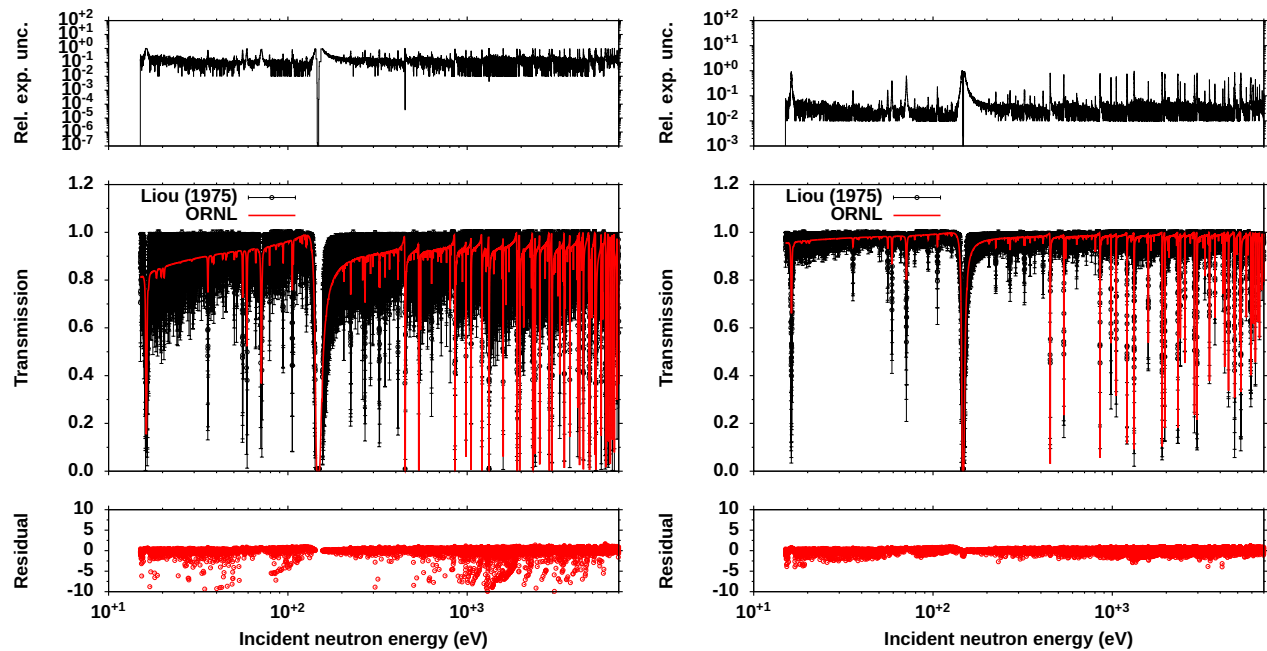


Figure 11. $n+^{164}\text{Dy}$ Liou's thick (left) and thin (right) transmission data measured in the neutron energy range up to 1 keV (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

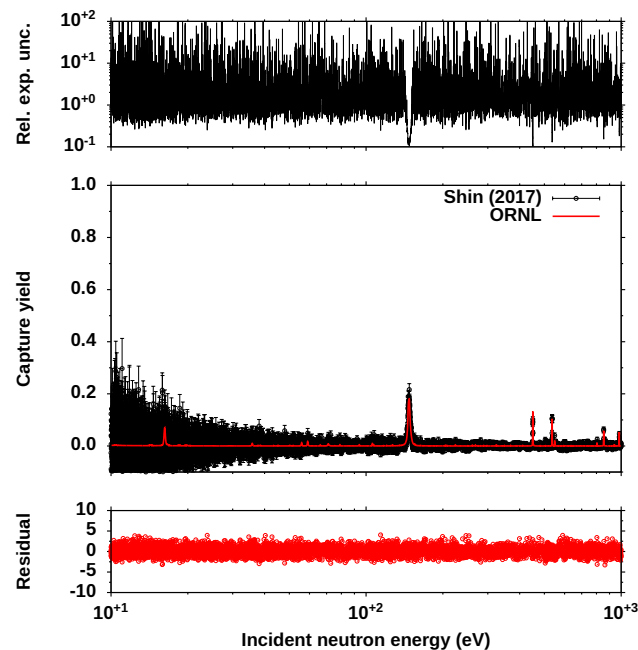


Figure 12. $n+^{164}\text{Dy}$ capture yield measured in epithermal energy range at the RPI facility (black dots) compared to theoretical data (solid red line). Relative experimental uncertainty (top) and residual (bottom) are shown in black solid lines and red dots, respectively.

current recommendations to avoid the submittal of ENDF files for natural nuclei, seven RPC matrices (one for each isotope) were generated. This was performed by extracting from the full RPC matrix the sub-matrix corresponding to the specific isotope and stored in a ENDF-formatted file with LRF=3 and LCOMP=1 options. Clearly, this procedure disregards the possible isotope–isotope correlations arising primarily by the fit of experimental data measured on $^{\text{nat}}\text{Dy}$ samples. For simplicity but also clarity, the cross section uncertainties and related correlation matrices calculated for the single-isotope RPC for all available reaction channels of each isotope are reported in a group representation averaged over 189 energy bins as shown in Figs. 13–19. Except for two minor isotopes $^{156,158}\text{Dy}$ and ^{162}Dy , the evaluated cross sections feature an uncertainty from about 0.5% to approximately 1–2% in the thermal energy range up to about a few eVs[†]. This relatively low uncertainty is driven by the very low uncertainty ($\ll 1\%$) reported by the measurements on natural samples as shown in Fig. 3. In the same energy region, the correlations are generally 100% because of the strong correlations of resonance parameters for negative levels usually associated with large neutron widths. Above a few eVs, the cross section uncertainties range between 1 and 10%, with mostly short-range positive and negative correlations. The plots also show the cross-reaction correlation matrix of the elastic and capture channels that is generally populated by large portions of negative correlations indicating anti-correlations between the two reaction channels.

[†]To ensure that the cross section uncertainties were not below 0.1%, a multiplication factor was applied to two isotopes, $^{161,164}\text{Dy}$, over a short energy range. The energy-dependent uncertainty file as implemented in the SAMMY code is shown in Fig. 44

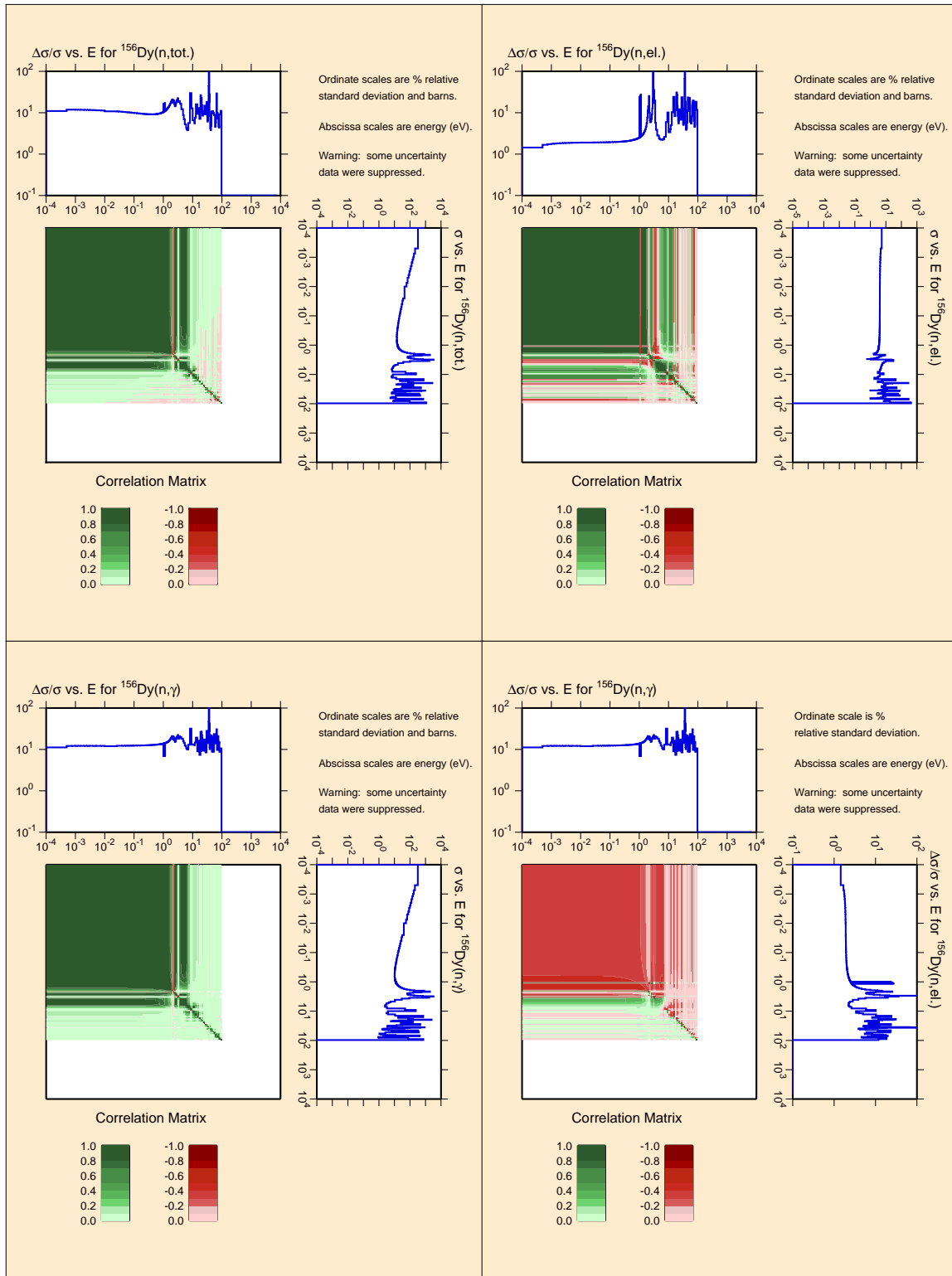


Figure 13. 189-group representation of ^{156}Dy cross sections and related correlation matrices for total, elastic, and capture reaction channels.

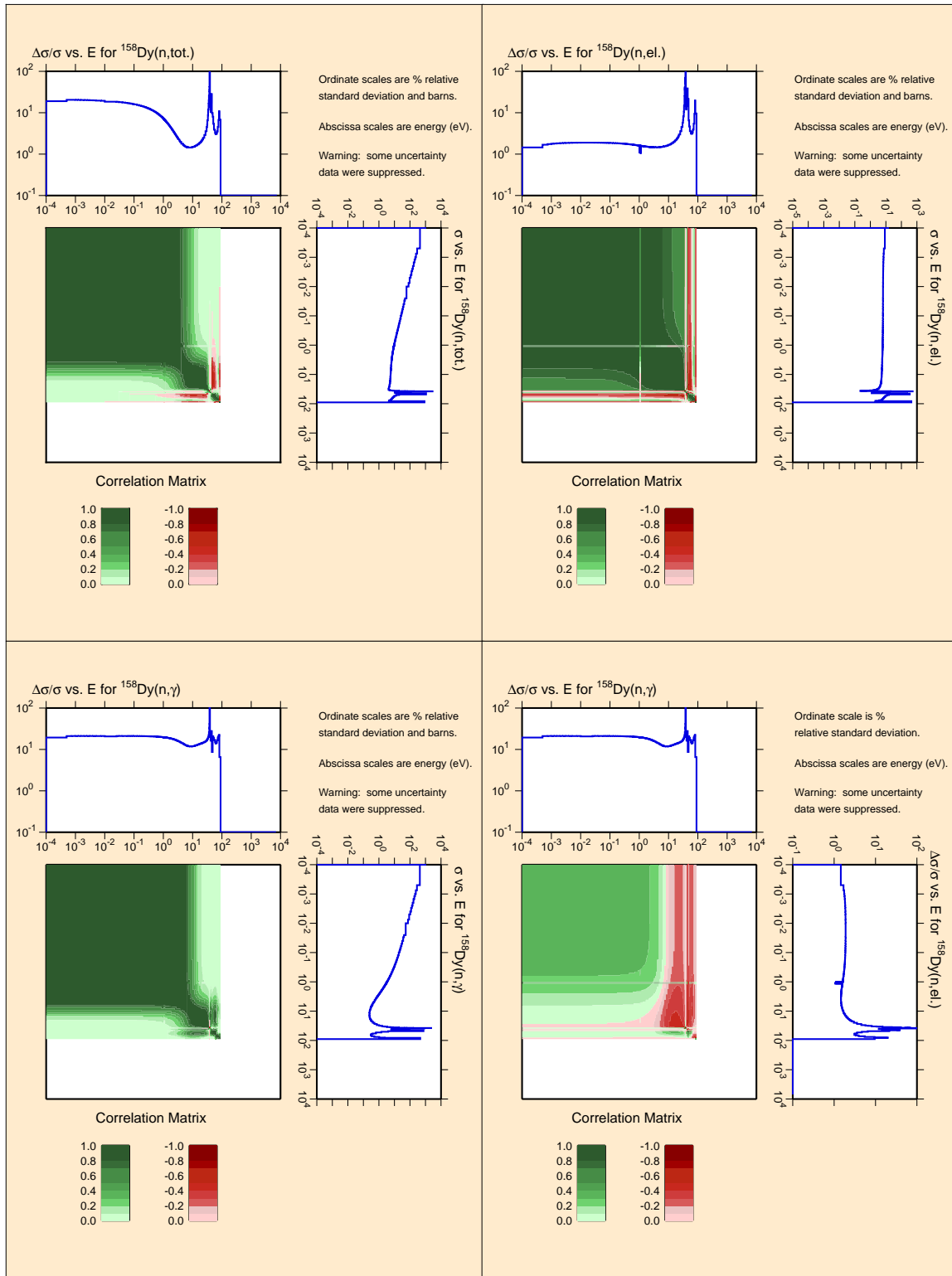


Figure 14. 189-group representation of ^{158}Dy cross sections and related correlation matrices for total, elastic, and capture reaction channels.

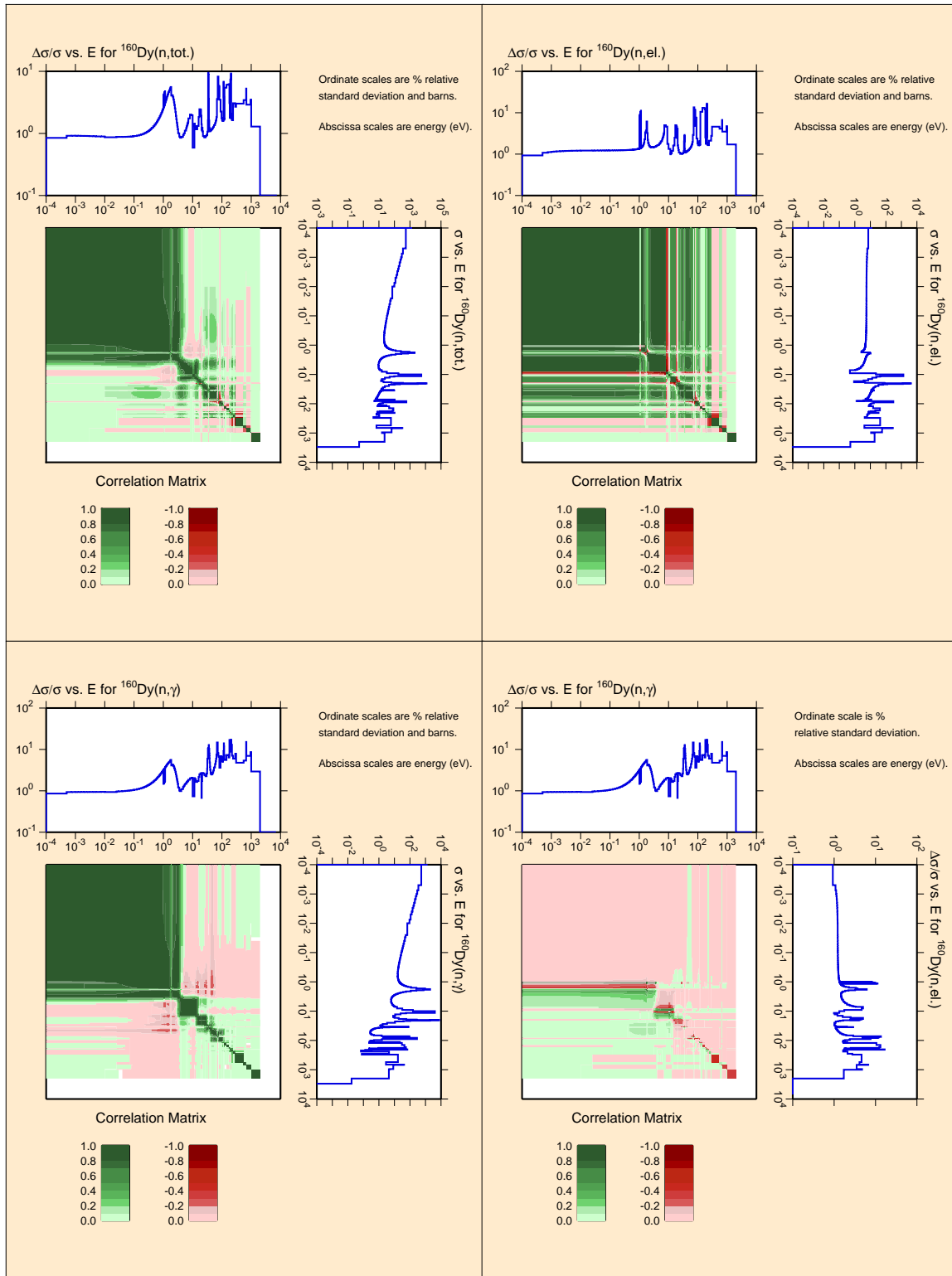


Figure 15. 189-group representation of ^{160}Dy cross sections and related correlation matrices for total, elastic, and capture reaction channels.

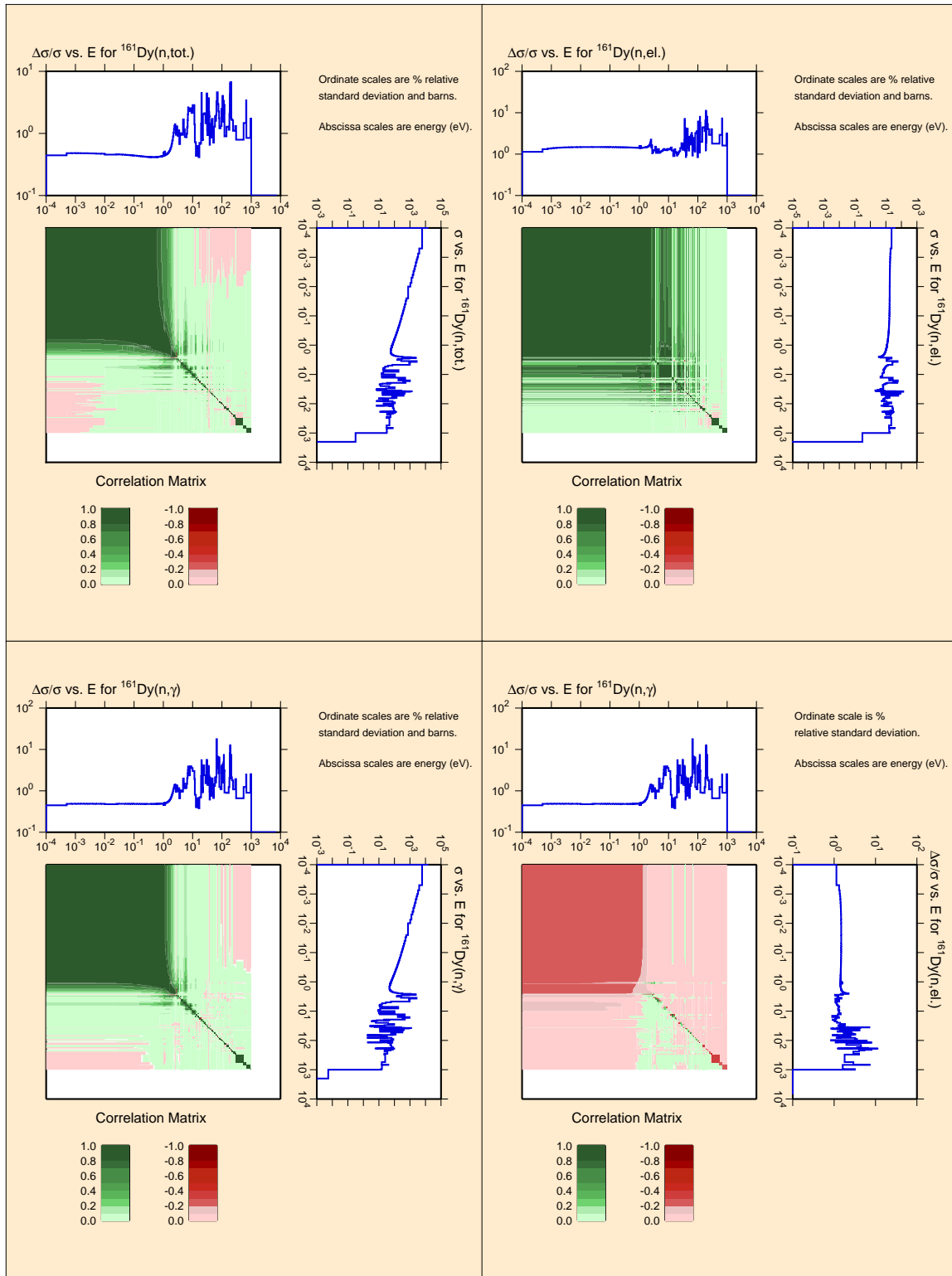


Figure 16. 189-group representation of ^{161}Dy cross sections and related correlation matrices for total, elastic, and capture reaction channels.

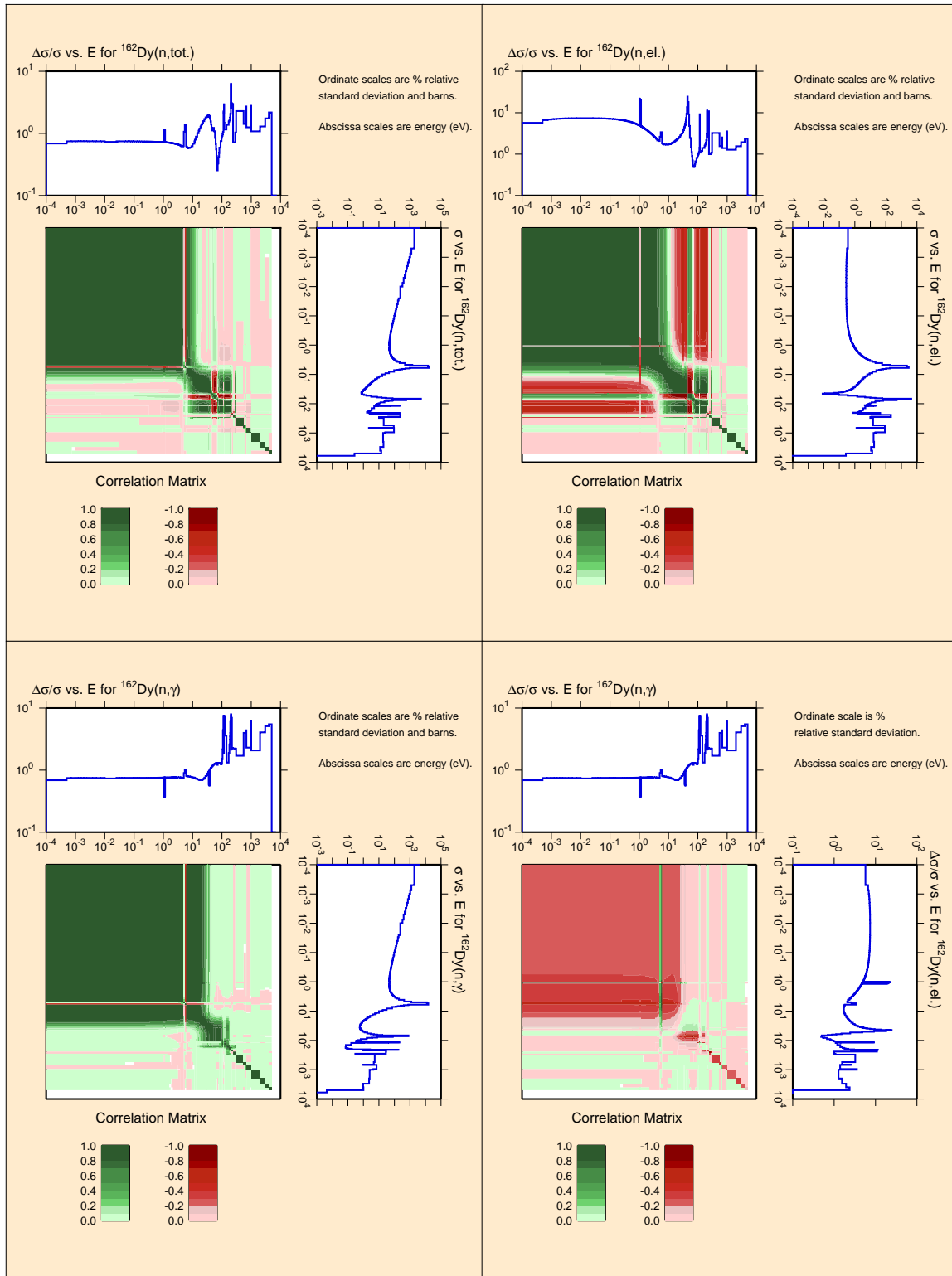


Figure 17. 189-group representation of ^{162}Dy cross sections and related correlation matrices for total, elastic, and capture reaction channels.

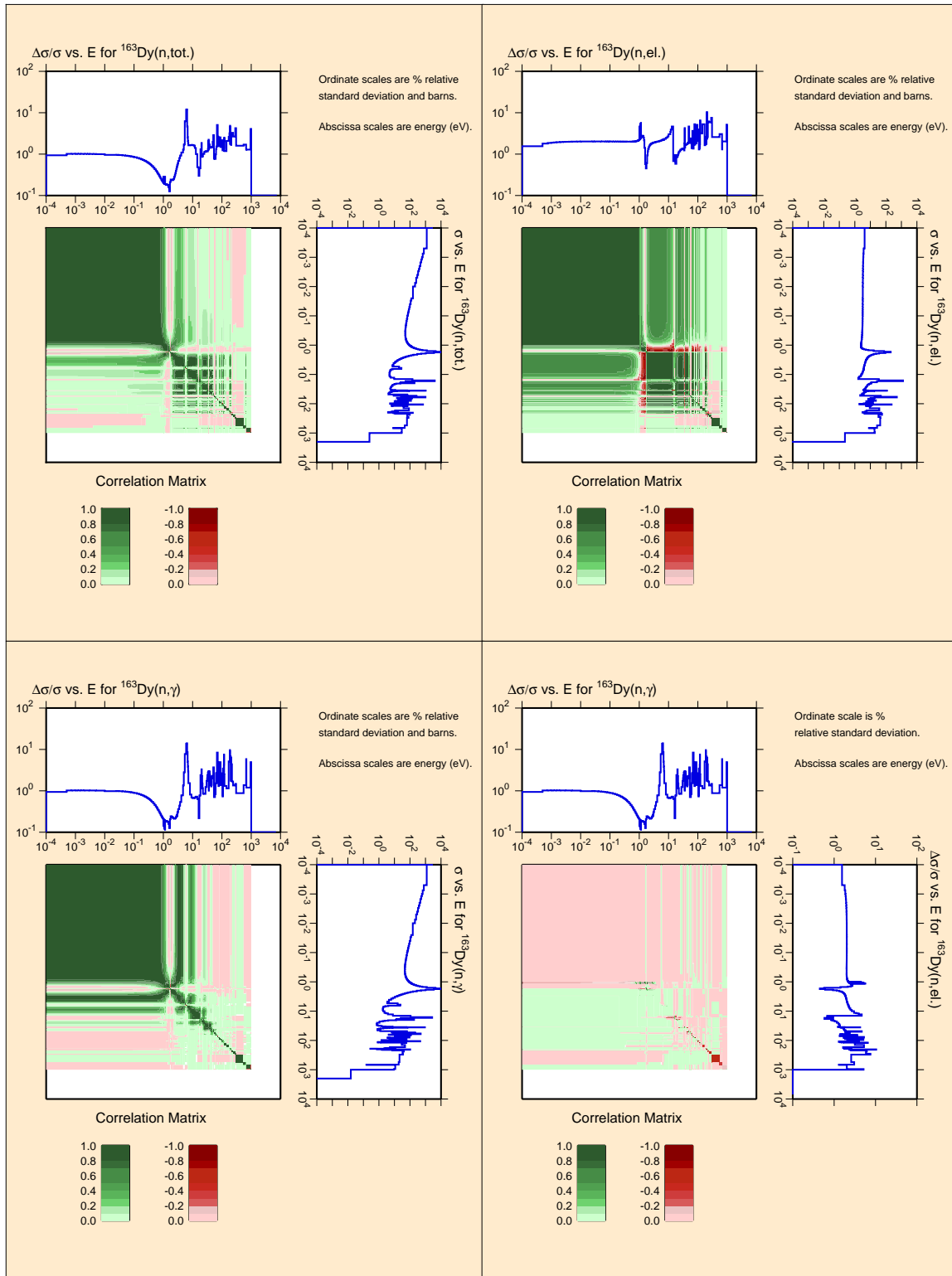


Figure 18. 189-group representation of ^{163}Dy cross sections and related correlation matrices for total, elastic, and capture reaction channels.

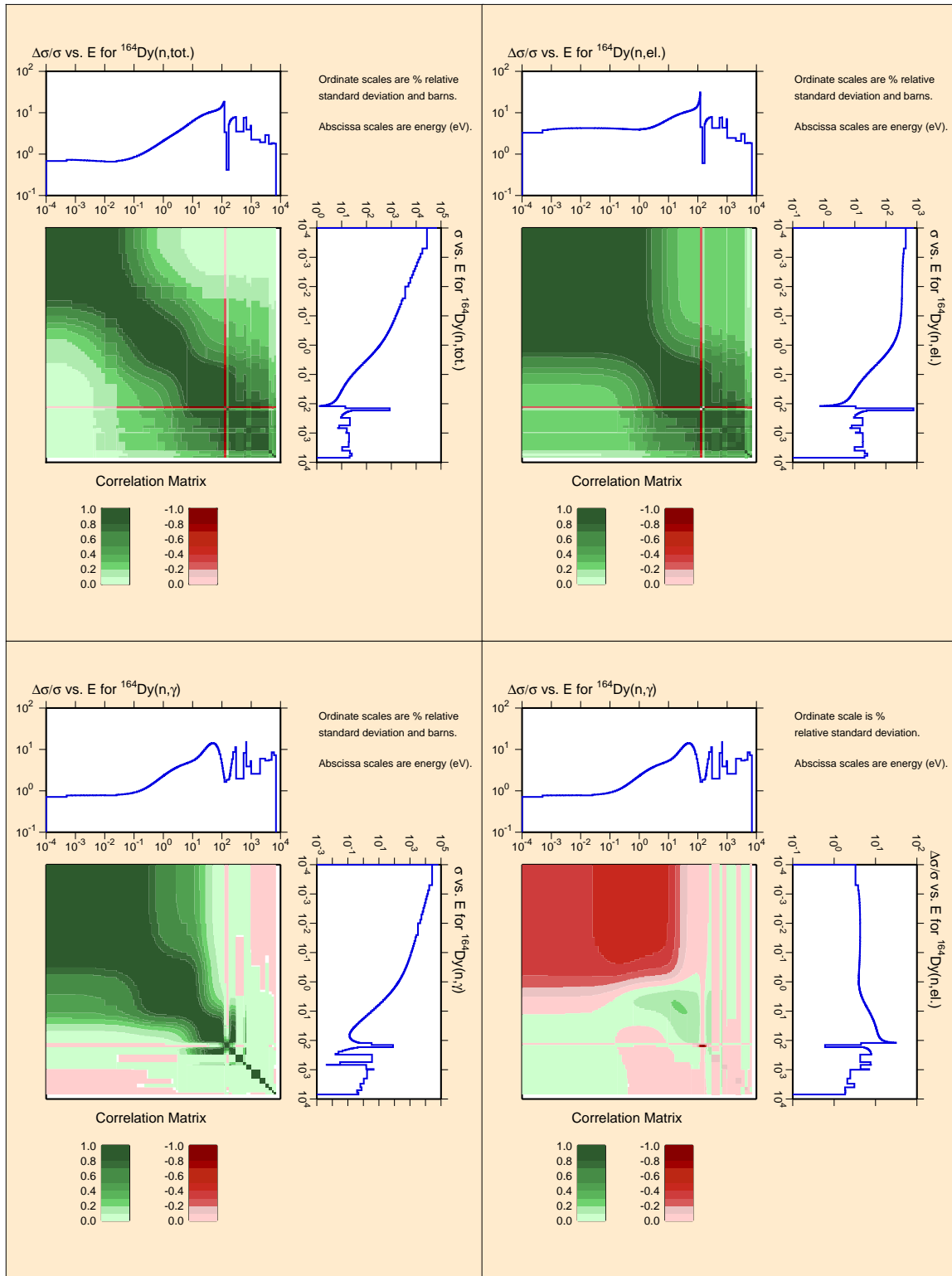


Figure 19. 189-group representation of ^{164}Dy cross sections and related correlation matrices for total, elastic, and capture reaction channels.

5. STATISTICAL PROPERTY OF THE RESONANCE PARAMETERS

The statistical analysis of the resonance parameters can be seen as a test to verify the nuclear reaction theory concerning the Wigner distribution for the level spacings, the Porter–Thomas distribution for the reaction widths, and related reaction channel multiplicity. Therefore, in addition to presenting our results for the observed neutron resonance energies, E , neutron and capture widths, Γ_n^J and Γ_γ^J , respectively, it is useful to report the study of their statistical properties, such as level spacing systematics and strength functions (or corresponding pole strength). In Figs. 20–26, cumulative plots of the number of observed s -wave vs energy for seven isotopes, $^{156,158,160-164}\text{Dy}$, are displayed with the estimated average level spacings and strength functions. These values were calculated from the fit of the observed energy levels and neutron widths over the entire positive energy range, although the choice of the upper energy in the fit can clearly impact the fitted values of the level spacing $\langle D_0 \rangle$ and strength function $\langle S_0 \rangle$ (or pole strength $\langle s_0 \rangle$). The estimate of the average s -level spacing $\langle D_0 \rangle$ reported in the figures was obtained as the inverse of the slope of a straight line fitted to the observed energy levels as

$$\frac{1}{\langle D_0 \rangle} = \frac{\lambda}{\Delta E_\lambda}, \quad (10)$$

where $\Delta E_\lambda = \lambda(E_\lambda - E_1)/(\lambda - 1)$ and $\lambda > 1$ is the energy level index for monotonically increasing positive energies whose upper limit is the total number of s -levels $\Lambda \equiv \sup\{\lambda\}$. For the strength function, the following relation was used:

$$S_l \Delta E_\Lambda = \sum_{\lambda>1}^{\Lambda} g_\lambda \Gamma_{n\lambda}^{0J(l)} v_l / (2l + 1), \quad (11)$$

where $\Gamma_{n\lambda}^{0J(l)} = \Gamma_{n\lambda}^{J(l)} / \sqrt{1\text{eV}/E_\lambda}$ is the reduced width and $v_l = \rho/P_l$ with $\rho = kR$ and $P_l \equiv P_l(\rho)$ is the penetrability factor. For s -wave, v_l is clearly unitary because $P_0 = \rho$.

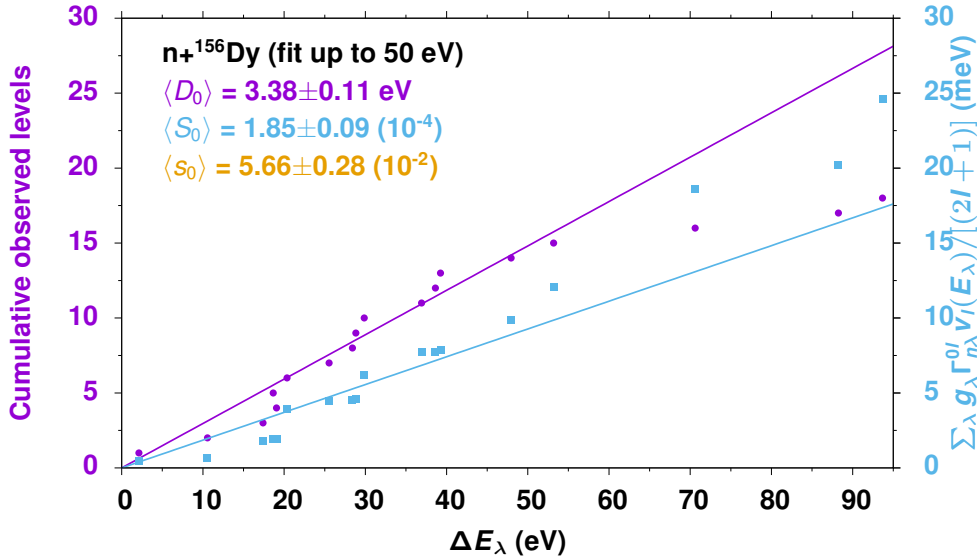


Figure 20. Plot of the cumulative number of observed s -wave (purple dots) vs energy for $n+^{156}\text{Dy}$. The values of average s -level spacings $\langle D_0 \rangle$ shown in the plot represent the inverse of the slope of a straight line fitted to the data (purple and blue lines) up to 50 eV. In blue, the fit of the neutron strength function S_0 is also shown together with the value of the related pole strength s_0 (in yellow).

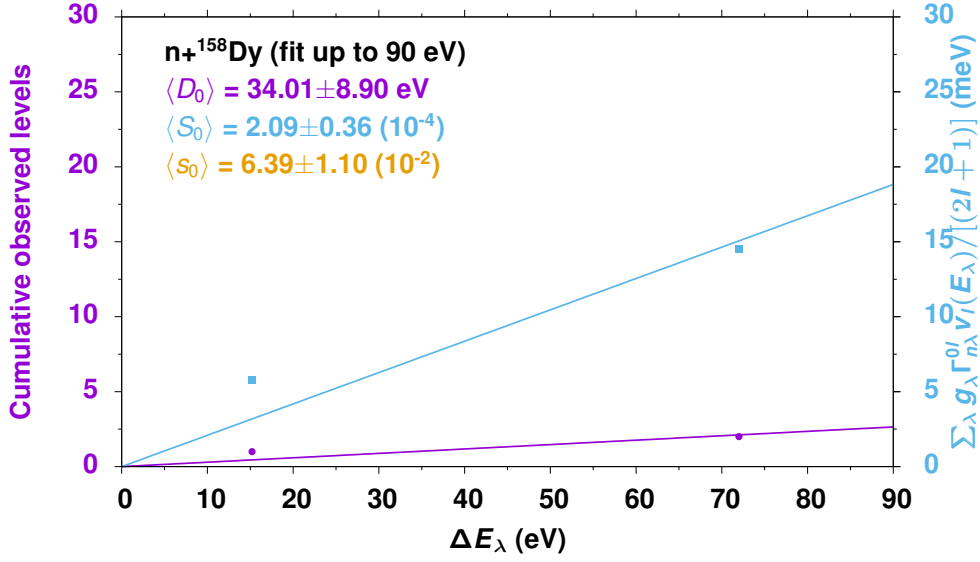


Figure 21. Plot of the cumulative number of observed s -wave (purple dots) vs energy for $n+^{158}\text{Dy}$. The values of average s -level spacings $\langle D_0 \rangle$ shown in the plot represent the inverse of the slope of a straight line fitted to the data (purple and blue lines) up to 90 eV. In blue, the fit of the neutron strength function S_0 is also shown together with the value of the related pole strength s_0 (in yellow).

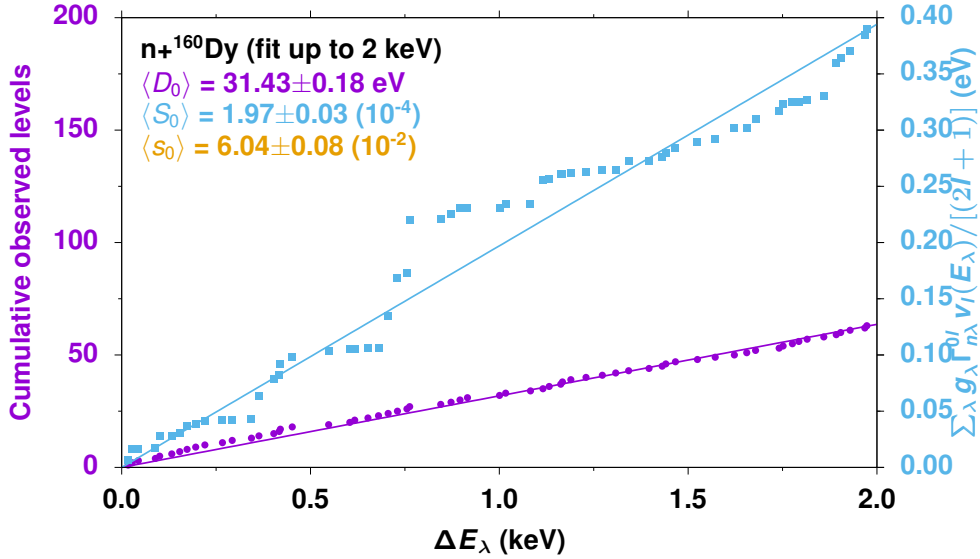


Figure 22. Plot of the cumulative number of observed s -wave (purple dots) vs energy for $n+^{160}\text{Dy}$. The values of average s -level spacings $\langle D_0 \rangle$ shown in the plot represent the inverse of the slope of a straight line fitted to the data (purple and blue lines) up to 90 eV. In blue, the fit of the neutron strength function S_0 is also shown together with the value of the related pole strength s_0 (in yellow).

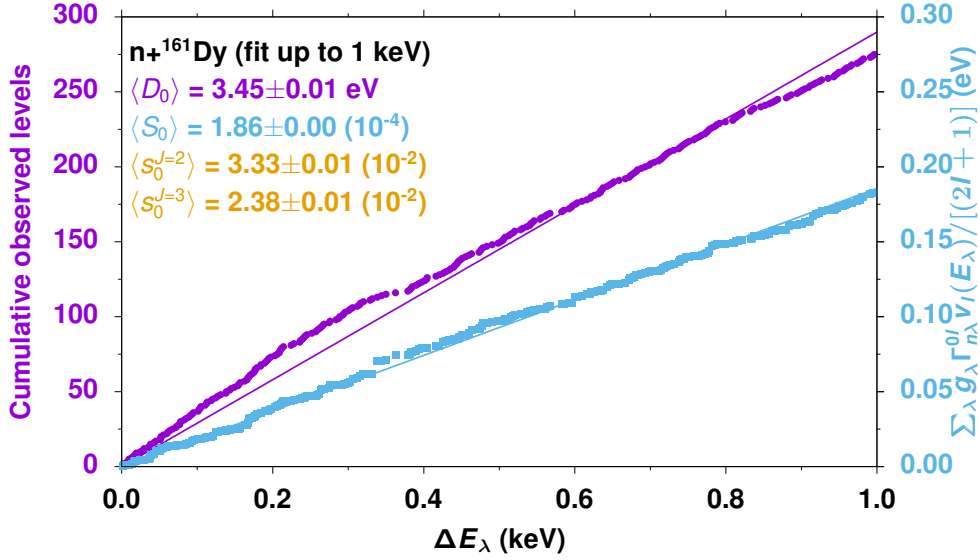


Figure 23. Plot of the cumulative number of observed s -wave (purple dots) vs energy for $n+^{161}\text{Dy}$. The values of average s -level spacings $\langle D_0 \rangle$ shown in the plot represent the inverse of the slope of a straight line fitted to the data (purple and blue lines) up to 1 keV. In blue, the fit of the neutron strength function S_0 is also shown together with the value of the related pole strength s_0 (in yellow).

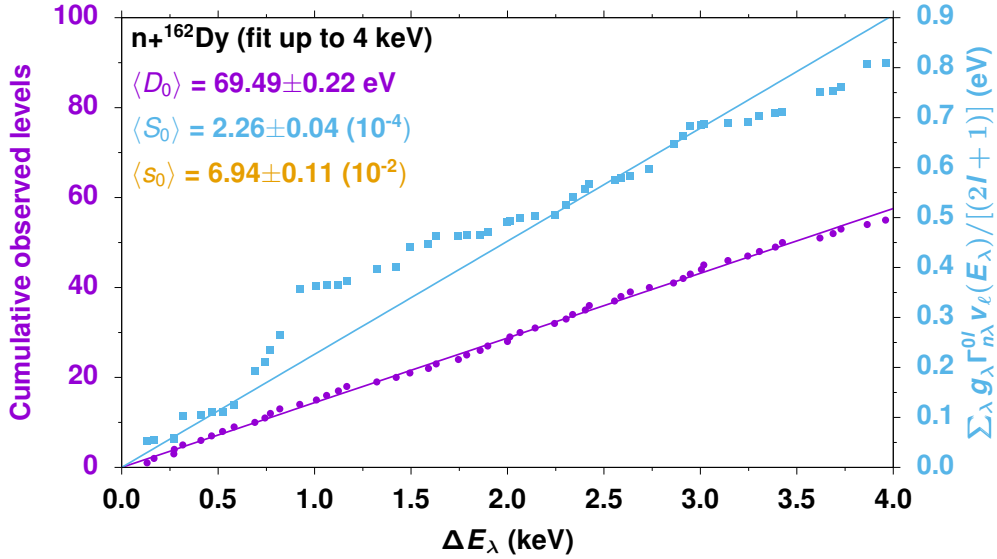


Figure 24. Plot of the cumulative number of observed s -wave (purple dots) vs energy for $n+^{162}\text{Dy}$. The values of average s -level spacings $\langle D_0 \rangle$ shown in the plot represent the inverse of the slope of a straight line fitted to the data (purple and blue lines) up to 5 keV. In blue, the fit of the neutron strength function S_0 is also shown together with the value of the related pole strength s_0 (in yellow).

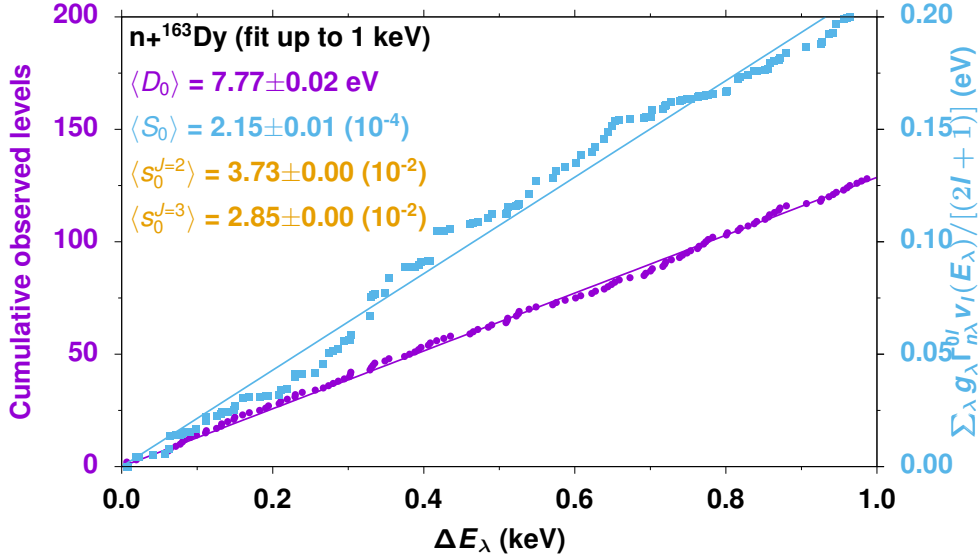


Figure 25. Plot of the cumulative number of observed s -wave (purple dots) vs energy for $n+^{163}\text{Dy}$. The values of average s -level spacings $\langle D_0 \rangle$ shown in the plot represent the inverse of the slope of a straight line fitted to the data (purple and blue lines) up to 1 keV. In blue, the fit of the neutron strength function S_0 is also shown together with the value of the related pole strength s_0 (in yellow).

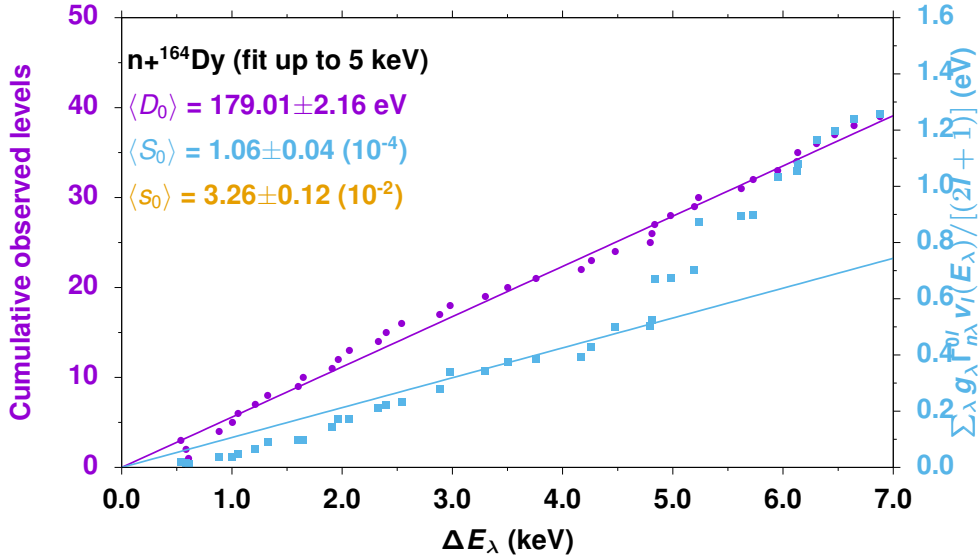


Figure 26. Plot of the cumulative number of observed s -wave (purple dots) vs energy for $n+^{164}\text{Dy}$. The values of average s -level spacings $\langle D_0 \rangle$ shown in the plot represent the inverse of the slope of a straight line fitted to the data (purple and blue lines) up to 7 keV. In blue, the fit of the neutron strength function S_0 is also shown together with the value of the related pole strength s_0 (in yellow).

6. REPOSITORY

The set of directories and related files used to generate the set of Dy evaluations has been stored in a GitHub repository found at the following link <https://code-int.ornl.gov/xtp/dysprosium[‡]>. When possible, the structure of the repository is designed to have a set of directories following the logical steps of the evaluation procedure. Each directory contains the files, scripts, and subdirectories necessary to reproduce the files submitted to ENDF repository. These folders are briefly described below and reported in such a way that roughly reproduces the execution order.

- **parameter** contains different versions of parameter files with a name convention **dy-ver.par**, where **ver** is the version label—for instance, **endf**, **beta1**, **final**. In addition to the resonance parameters (level energies, channel radii and widths), the files are set up to contain quantum number information and related particle pairs for seven Dy isotopes. As explained later, the isotopic abundance of each isotope is specified in the input files together with other experimental corrections.
 - **geraspin** is designed to generate the quantum number information as reported in the parameter files above. This is performed by running the **dy.csh** script, which uses existing input files for the SAMQUA code of the type **dy-wave.inp**, where **wave** is the input variable of the script and refers to the number of partial waves setup for each input file. The file **masses.gp** uses GNUPLOT to print the isotope masses as reported in the parameter files.
 - **coher** computes thermal properties such as thermal cross sections and scattering lengths as reported in Table 5 for each isotope from a specific **dy-ver.par** file.
 - **stat** generates plots of the statistical properties of the resonance parameters from a specific **dy-ver.par** file by using GNUPLOT after running the SAMDIST code. This is performed by running the **run_stat.csh** script with **ver** as input variable. Assumptions on spin group numbers listed in the parameter file are made. Based on these, the input files to run the SAMDIST code are automatically generated in **testdist.csh**. The script **stat.csh** uses files with extension **gp** to generate plots of resonance parameter statistical properties from a specific parameter file.
- **exfor** contains the measured data files with extension **exf** retrieved from the EXFOR library. These are used to generate input files for SAMMY with extension **twenty**. The files with extension **exf** and **twenty** contain the same information except for minor changes like negative values for the data type such as cross section and neutron capture yield, or inclusion of data uncertainty if not explicitly given in the EXFOR library.
 - The subdirectory **crunch** contains the plots for the crunch tables generated by the GNUPLOT script **crunch.gp**, which reads the measured data files. The derived crunch tables are the same used in the SAMMY input files.
- **inputs** contains the SAMMY inputs corresponding to each experiments. These files are set up with alpha-numeric cards to solely contain parameters of the experimental configurations. The quantum-number information, i.e., particle pair definitions with related channel spins and radii, are included in the parameter file in the **parameter** directory.

[‡]Available only internally for ORNL users.

- **thermal** is a directory containing data files in **twenty** stored in the subdirectory **data** and referring to different compilation of thermal values for each Dy isotope.
- **runs** is defined to calculate cross section or yield data for a set of measured data listed in the script **sammy.csh**. The data are reconstructed for a specific parameter **ver** found in the script.
- **cov** contains scripts to both perform the fit of selected measured data and generate ENDF-formatted file 2 and 32 for each isotope to be used in the final assembly of the ENDF file. The resonance parameter information and associated covariance matrix is extracted from the **SAMMY.COV** binary file which a multi-isotope covariance matrix obtained by the global fit of seven Dy isotopes.
- **dy1[**]** are directories containing **SAMMY.LPT** and **SAMMY.LST** files for a specific set of resonance parameters.
- **ndf** is a directory for the generation of a set of ENDF files for each Dy isotope with minimal information for cross section and covariance processing purposes. The information includes MF=2 and MF=32 sections starting from a parameter file **dy-ver.par** stored in the **parameter** directory. The ENDF file includes a section for MF=3 and MF=33 with zeros to allow processing with a code such as NJOY. The script to generate and process minimal ENDF files including only cross section information such as MF=2 and MF=3 sections for each Dy isotope is **ndf.csh**. The input parameter of this script is the specific parameter version **ver**. The script **ndf_cov.csh** is used similarly to **ndf.csh** but designed to include information for MF=32 and MF=33 sections in the ENDF file to quantify processed cross section covariances as well. The processing is performed by calling the script **xcs-res.constE-dy.csh**.
- **report** is a directory storing the files and figures used to generate the present ORNL/TM report.
- **nndc** contains the final ENDF-formatted files, **n-066_Dy_iso.endf**, submitted to the National Nuclear Data Center.

7. CONCLUSIONS

The set of the RM parameters for neutron-induced resonances of seven Dy isotopes, $^{156,158,160-164}\text{Dy}$, in the incident energy range from thermal to several keVs was generated by applying the SAMMY evaluation procedure. The multi-level multi-channel *R*-matrix shape fitting analysis was performed on available experimental data for natural capture yield and transmission data as well as isotopically enriched capture yield data. Experimental conditions such as resolution function, finite size sample, detector efficiencies, and nuclide abundances of sample, multiple scattering, self-shielding, normalization, background, and Doppler broadening were taken into account.

In addition to the guidance provided by Sears' (or NIST) compilation [12] for the thermal constants, the experimental database in the thermal and low-energy region up to 15 eV consisted of both high-resolution capture yields and transmission data [7]. Due to their extension to thermal energies, these data were extremely relevant in the resonance parameter evaluation. In this region, the evaluated thermal scattering and capture constant for the ^{164}Dy isotope shows differences with NIST compilation up to 8%. Also for ^{161}Dy isotopes, this work reports a value of the scattering cross section 10% higher than the NIST's reported value but in agreement with other evaluated works [6]. Particular emphasis should be devoted in future experimental campaigns to measuring scattering lengths for these two isotopes, particularly for the most abundant isotope, ^{164}Dy , with the largest total thermal cross section among the seven Dy isotopes. Moreover, future evaluation works should develop the capability to include scattering lengths in the fitting procedure, among other possible measured quantities. Uncertainty quantification of the total, elastic scattering, and capture cross section for all isotopes has been reported, for convenience, in 189-group representation as shown in Figs. 13–19. The uncertainty in the thermal energy region were about 1% constrained by the transmission, and capture measurements on natural samples [7] were reported to have a similar uncertainty.

In the energy region above 15 eV, although high-resolution capture yield data [8] are available, high-resolution transmission data are still needed for isotopically enriched samples. The currently available Liou's transmission data [5] are poorly documented in the uncertainty quantification analysis, and it seems very likely that the reported data communicated to the EXFOR database are not measured data but total cross section data reconstructed from resonance parameters fitted to the measured transmission data. Moreover, Liou's measurements campaign included capture measurements that are not available in the EXFOR library. The lack of modern transmission data measured on isotopically enriched samples in the neutron energy above 15 eV also precludes any attempt to extend the RRR to the currently reported upper energy ranges.

Integral benchmarks reporting reactivity coefficients such as k_{eff} are usually used to validate underlying nuclear data for a given configuration of materials—including fissile actinides and other chosen nuclei—for instance, to test their natural propensity to absorb neutrons. Reactivity coefficients of interest for criticality safety applications can be, by design, very sensitive to the thermal and epithermal region up to a few eVs. Therefore, the present evaluation based on Block's recent transmission data (measured from thermal up to 15 eV) represents a perfect example of nuclear data ready for validation purposes. However, in the specific case of Dy isotopes, there is a very limited number of modern benchmarks that can be used to perform a conclusive validation over a comprehensive suite of cases. Therefore, no particular validation test was performed on the evaluated data presented in this work.

Finally, this work represents one of the first attempts to generate a fully reproducible evaluation in the RRR as this is the primary goal of the Working Party on International Nuclear Data Evaluation Co-operation Subgroup 49.

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APPENDIX A. SAMMY inputs

The SAMMY inputs containing the experimental configuration for each measured data included in the evaluation work are reported in Figs. 27–42. For Liou’s transmission data, the experimental configuration consists of the crunch data, the sample thickness, and related isotopic abundances. In addition to these, detector efficiencies are usually reported for RPI’s capture measurements together with parameters for the resolution function.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy160      159.925204 0.00001 2.0000E+03      1      1 0 0 0 0      0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

      293.6      202.05      0.0012      0.0000      -0.025
      4.e-2      6
      8.0e+1      3.5e+1      2.4e+2      1.6e+1      5.5e+2      8.0e+0      9.0e+2      4.0e+0
      1.2e+3      2.0e+0      1.0e+3      1.0e+0
      7.45613 8.881e-4      1.e-2
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00020000 1.0000E-5 0 1
157.924415 .00020000 1.0000E-5 0 2
159.925204 .69460000 1.0000E-5 0 3
160.926939 .17830000 1.0000E-5 0 4 5
161.926805 .06450000 1.0000E-5 0 6 7 8
162.928737 .03550000 1.0000E-5 0 910
163.929181 .02670000 1.0000E-5 0111213

#MISCEllaneous parameters follow
#TZERO 1 1 .00084733 .00100000 1.0001034 .00100000 202.05000

```

Figure 27. SAMMY input for n+¹⁶⁰Dy Liou's transmission data.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy161      160.926939  0.00001 1.0000E+03    1    1 0  0 0  0      0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    202.05    0.0012    0.0000    -0.025
    4.e-2     5
    8.0e+1    3.5e+1    2.4e+2    1.6e+1    5.5e+2    8.0e+0    9.0e+2    4.0e+0
    1.2e+3    1.0e+0
    7.46999  7.451e-3
                                1.e-2
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00020000 1.0000E-5 0 1
157.924415 .00020000 1.0000E-5 0 2
159.925204 .00350000 1.0000E-5 0 3
160.926939 .95620000 1.0000E-5 0 4 5
161.926805 .02530000 1.0000E-5 0 6 7 8
162.928737 .00900000 1.0000E-5 0 910
163.929181 .00560000 1.0000E-5 0111213

```

Figure 28. SAMMY input for n+¹⁶¹Dy Liou's transmission thick sample data.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy161      160.926939  0.00001 2.0000E+02    1    1 0 0 0 0    0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    202.05    0.0012    0.0000    -0.025
    4.e-2    2
    8.0e+1    3.5e+1    2.4e+2    1.6e+1
    7.46999  1.644e-3                1.e-2
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00020000 1.0000E-5 0 1
157.924415 .00020000 1.0000E-5 0 2
159.925204 .00350000 1.0000E-5 0 3
160.926939 .95620000 1.0000E-5 0 4 5
161.926805 .02530000 1.0000E-5 0 6 7 8
162.928737 .00900000 1.0000E-5 0 910
163.929181 .00560000 1.0000E-5 0111213

```

Figure 29. SAMMY input for n+¹⁶¹Dy Liou's transmission thin sample data.

```

Dy / capture / rpi data / Original filename Dy161.dat ( or very similarly Dy161_
RPI_Cap.txt)
Dy161 160.926939 1.00000 2.0000E+03 1 1 0 0 0 0 0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
INFINITE SLAB
DOUBLE
YIELD
DO NOT SHIFT ENERGY
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

      293.6 25.5686 0.005800 0.0600000 -0.01800 5.0 0.001
      0.01280 4
      99.91198 8.000 499.67926 4.000 999.63580 2.000 9999.79785 1.000
      7.469997 6.359E-4 1.000E-3
CAPTURE
0.0287000 1.0200 0.0 1.0200 0.0

DETECTOR EFFICIENCIES
1.000000001 0.0000000 0 1
1.000000001 0.0000000 0 2
1.000000000 0.0000000 0 3
1.000000000 0.1000000 0 4 5
0.86405800 0.1000000 0 6 7 8
0.93911600 0.1000000 0 910
0.88117600 0.1000000 0111213

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00020000 1.0000E-5 0 1
157.924415 .00020000 1.0000E-5 0 2
159.925204 .00350000 1.0000E-5 0 3
160.926939 .95660000 1.0000E-5 0 4 5
161.926805 .02530000 1.0000E-5 0 6 7 8
162.928737 .00860000 1.0000E-5 0 910
163.929181 .00560000 1.0000E-5 0111213

```

Figure 30. SAMMY input for n+¹⁶¹Dy RPI's capture data.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy162      161.926805  0.00001 1.5000E+04    1    1 0 0 0 0    0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    202.05    0.0012    0.0000    -0.025
    4.e-2     6
    8.0e+1    3.5e+1    2.4e+2    1.6e+1    5.5e+2    8.0e+0    9.0e+2    4.0e+0
    1.2e+3    2.0e+0    1.5e+4    1.0e+0
    5.90000  6.934e-3    1.e-2
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00010000 1.0000E-5 0 2
159.925204 .00800000 1.0000E-5 0 3
160.926939 .01240000 1.0000E-5 0 4 5
161.926805 .96130000 1.0000E-5 0 6 7 8
162.928737 .01790000 1.0000E-5 0 910
163.929181 .00720000 1.0000E-5 0111213

```

Figure 31. SAMMY input for $n+^{162}\text{Dy}$ Liou's transmission thick sample data.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy162      161.926805  0.00001 1.5000E+04    1    1 0  0 0  0      0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    202.05    0.0012    0.0000    -0.025
    4.e-2     6
    8.0e+1    3.5e+1    2.4e+2    1.6e+1    5.5e+2    8.0e+0    9.0e+2    4.0e+0
    1.2e+3    2.0e+0    3.0e+3    1.0e+0
    5.90000  1.533e-3                                1.e-2
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00010000 1.0000E-5 0 2
159.925204 .00800000 1.0000E-5 0 3
160.926939 .01240000 1.0000E-5 0 4 5
161.926805 .96130000 1.0000E-5 0 6 7 8
162.928737 .01790000 1.0000E-5 0 910
163.929181 .00720000 1.0000E-5 0111213

```

Figure 32. SAMMY input for n+¹⁶²Dy Liou's transmission thin sample data.

```

Dy / capture / rpi data / Original filename Dy162.dat ( or very similarly Dy162_
RPI_Cap.txt)
Dy162      161.926805  10.0000 1.0000E+03      1      1 0 0 0 0      0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
INFINITE SLAB
DOUBLE
YIELD
DO NOT SHIFT ENERGY
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

      293.6  25.5686   0.005800 0.0600000  -0.01800      5.0      0.001
      0.01280      4
      0.999795  256.0   9.9884176   128.0   49.982803   16.00   9999.71582   2.000
      5.900000  6.445E-4      1.000E-3
CAPTURE
      0.0295000      0.7545      0.0      0.7545      0.0

DETECTOR EFFICIENCIES
1.000000001 0.00000000 0 1
1.000000001 0.00000000 0 2
1.000000000 0.00000000 0 3
1.15733000 0.10000000 0 4 5
1.000000000 0.10000000 0 6 7 8
1.08687000 0.10000000 0 910
1.01981000 0.10000000 0111213

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00010000 1.0000E-5 0 2
159.925204 .00080000 1.0000E-5 0 3
160.926939 .01240000 1.0000E-5 0 4 5
161.926805 .96170000 1.0000E-5 0 6 7 8
162.928737 .01780000 1.0000E-5 0 910
163.929181 .00710000 1.0000E-5 0111213

```

Figure 33. SAMMY input for n+¹⁶²Dy RPI's capture data.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy163      162.92873  0.00001 1.0000E+03   1   1 0 0 0 0   0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    202.05    0.0012    0.0000    -0.025
    4.e-2     5
    8.0e+1    3.5e+1    2.4e+2    1.6e+1    5.5e+2    8.0e+0    9.0e+2    4.0e+0
    1.2e+3    1.0e+0
    7.49754  1.403e-3
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00010000 1.0000E-5 0 2
159.925197 .00030000 1.0000E-5 0 3
160.926933 .00360000 1.0000E-5 0 4 5
161.926798 .01230000 1.0000E-5 0 6 7 8
162.928731 .96840000 1.0000E-5 0 910
163.929174 .01520000 1.0000E-5 0111213

```

Figure 34. SAMMY input for n+¹⁶³Dy Liou's transmission data.

```

Dy / capture / rpi data / Original filename Dy163_RPI_Cap.txt
Dy163      162.928737  10.0000 1.0000E+03    1    1 0 0 0 0      0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
INFINITE SLAB
DOUBLE
YIELD
DO NOT SHIFT ENERGY
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

      293.6  25.5686   0.005800 0.0600000 -0.01800    5.0      0.001
      0.01280    4
  99.91198   8.000   499.67926   4.000   999.63580   2.000   9999.79785   1.000
  7.497540 6.503E-4      1.000E-3
CAPTURE
0.0289000    1.0300    0.0    1.0300    0.0

DETECTOR EFFICIENCIES
1.00000001 0.0000000 0 1
1.00000001 0.0000000 0 2
1.00000000 0.0000000 0 3
1.06483000 0.1000000 0 4 5
0.92007700 0.1000000 0 6 7 8
1.00000000 0.1000000 0 910
0.93830400 0.1000000 0111213

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00010000 1.0000E-5 0 2
159.925204 .00030000 1.0000E-5 0 3
160.926939 .03600000 1.0000E-5 0 4 5
161.926805 .01230000 1.0000E-5 0 6 7 8
162.928737 .96860000 1.0000E-5 0 910
163.929181 .01520000 1.0000E-5 0111213

```

Figure 35. SAMMY input for n+¹⁶³Dy RPI's capture data.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy164      163.929181  0.00001 4.0000E+03    1    1 0 0 0 0    0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    202.05    0.0012    0.0000    -0.025
    4.e-2     6
    8.0e+1    3.5e+1    2.4e+2    1.6e+1    5.5e+2    8.0e+0    9.0e+2    4.0e+0
    1.2e+3    2.0e+0    1.5e+4    1.0e+0
    7.51121  1.054e-2    1.e-2
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00020000 1.0000E-5 0 2
159.925204 .00020000 1.0000E-5 0 3
160.926939 .00150000 1.0000E-5 0 4 5
161.926805 .00350000 1.0000E-5 0 6 7 8
162.928737 .01030000 1.0000E-5 0 910
163.929181 .98420000 1.0000E-5 0111213

```

Figure 36. SAMMY input for n+¹⁶⁴Dy Liou's transmission thick sample data.

```

Dy / Total / Liou(75) / ENTRY 10525 / Exp. corrections : background
Dy164      163.929181  0.00001 4.0000E+03   1   1 0  0 0  0       0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6      202.05      0.0012      0.0000      -0.025
    4.e-2      6
    8.0e+1      3.5e+1      2.4e+2      1.6e+1      5.5e+2      8.0e+0      9.0e+2      4.0e+0
    1.2e+3      2.0e+0      4.0e+3      1.0e+0
    7.51121    2.337e-3      1.e-2
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00020000 1.0000E-5 0 2
159.925204 .00020000 1.0000E-5 0 3
160.926939 .00150000 1.0000E-5 0 4 5
161.926805 .00350000 1.0000E-5 0 6 7 8
162.928737 .01030000 1.0000E-5 0 910
163.929181 .98420000 1.0000E-5 0111213

```

Figure 37. SAMMY input for n+¹⁶⁴Dy Liou's transmission thin sample data.

```

Dy / capture / rpi data / Original filename Dy164.dat (or very similarly Dy164_R
PI_Cap.txt)
Dy164      163.929182  1.00000 2.0000E+03      1      1 0  0 0  0          0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
INFINITE SLAB
DOUBLE
YIELD
DO NOT SHIFT ENERGY
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

      293.6  25.5686   0.005800 0.0600000  -0.01800      5.0      0.001
      0.01280      4
    99.91198   8.000  499.67926   4.000   999.63580   2.000  9999.79785   1.000
      7.511213 6.196E-4      1.000E-3
CAPTURE
0.0310000      0.7855      0.0      0.7855      0.0

DETECTOR EFFICIENCIES
1.000000001 0.0000000 0 1
1.000000001 0.0000000 0 2
1.000000000 0.0000000 0 3
1.13485000 0.1000000 0 4 5
0.98057400 0.1000000 0 6 7 8
1.06575000 0.1000000 0 910
1.000000000 0.1000000 0111213

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00010000 1.0000E-5 0 1
157.924415 .00010000 1.0000E-5 0 2
159.925204 .00010000 1.0000E-5 0 3
160.926940 .00140000 1.0000E-5 0 4 5
161.926805 .00350000 1.0000E-5 0 6 7 8
162.928738 .01030000 1.0000E-5 0 910
163.929181 .98450000 1.0000E-5 0111213

```

Figure 38. SAMMY input for n+¹⁶³Dy RPI's capture data.

```

Dy / capture / rpi data / Original filename DyNat.dat ( or very similar DyNat_20
mil_RPI_Cap)
Dynat      162.499472  1.00000 1.0000E+03    1    1 0 0 0 0      0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
DOUBLE
INFINITE SLAB
YIELD
DO NOT SHIFT ENERGY
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

      293.6  25.5686   0.005800 0.0600000  -0.01800    5.0      0.001
      0.01280      4
99.99477   8.000  499.98755   4.000   999.85242   2.000   9999.69434   1.000
      7.088000 1.6304E-3      0.00100
CAPTURE
0.0604000      1.9070      0.0 1.9070      0.0

DETECTOR EFFICIENCIES
1.000000001 0.0000000 0 1
1.000000001 0.0000000 0 2
1.000000000 0.0000000 0 3
1.00190000 0.1000000 0 4 5
0.86570000 0.1000000 0 6 7 8
0.94090000 0.1000000 0 910
0.88285000 0.1000000 0111213

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00056000 1.0000E-5 0 1
157.924415 .00095000 1.0000E-5 0 2
159.925204 .02329000 1.0000E-5 0 3
160.926939 .18889000 1.0000E-5 0 4 5
161.926805 .25475000 1.0000E-5 0 6 7 8
162.928737 .24896000 1.0000E-5 0 910
163.929181 .28260000 1.0000E-5 0111213

```

Figure 39. SAMMY input for $n + {}^{\text{nat}}\text{Dy}$ RPI's epithermal capture data.

```

Dy / capture / rpi data / Original filename Dy_tcnat_10-mil_CSISRS.twenty
Dynat      162.499472  0.00001 2.0000E+01  1  1 0 0 0 0 0
FGM
TWENTY
EV
DEBUG
GENERATE PLOT FILE AUTOMATICALLY
DO NOT SOLVE BAYES EQUATIONS
ENERGY UNCERTAINTIES ARE AT END OF LINE
EXPONENTIAL FOLDING
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
DOUBLE
YIELD
FINITE SLAB
DO NOT SHIFT ENERGY
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    25.444    0.0055    0.0000   -1.0
    0.125         3
    0.1237    64.0      4.0459    4.0      20.0051    1.000
    7.50000    8.050e-4
CAPTURE
0.0254000    2.5400      1.67259

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00060000 1.0000E-5 0 1
157.924415 .00100000 1.0000E-5 0 2
159.925204 .02340000 1.0000E-5 0 3
160.926939 .18910000 1.0000E-5 0 4 5
161.926805 .25510000 1.0000E-5 0 6 7 8
162.928737 .24900000 1.0000E-5 0 910
163.929181 .28180000 1.0000E-5 0111213

MISCELLANEOUS PARAMETERS FOLLOW
DELTE 0 00 6.28-5  1.22-5  0.234  0.016  -.038  0.003

DETECTOR EFFICIENCIES
0.83000000 0.0300000 0 1
0.83000000 0.0200000 0 2
0.81500000 0.0150000 0 3
0.90600000 0.0200000 0 4 5
0.80700000 0.0200000 0 6 7 8
0.89400000 0.0150000 0 910
0.78000000 0.0250000 0111213

#0.0254000    1.9000
#0.0254000    2.5400      1.67259 ! As in the EXFOR sheets
#  7.50000    8.050e-4      ! As discussed in May 2021 with D. Barry
#  7.50000    7.840e-4      ! As in the EXFOR sheets
#                      ! As in the Paper

```

Figure 40. SAMMY input for $n+^{nat}\text{Dy}$ RPI's thermal capture data.

```

Dysprosium / transmission / rpi data / Original filename Dy_ttnat_10mil_CSISRS.t
wenty
Dynat      162.499472  4.00001 2.0000E+03    1    1 0 0 0 0    0
FGM
TWENTY
EV
GENERATE PLOT FILE AUTOMATICALLY
PRINT THEORETICAL VALUES
PRINT DEBUG INFORMATION
DO NOT SOLVE BAYES EQUATIONS
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
DO NOT SHIFT RPI RESOLUTION FUNCTION TO CENTER
ENERGY UNCERTAINTIES ARE AT END OF LINE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    25.597    0.0055    0.0000
    7.50000  7.840e-4
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00060000 1.0000E-5 0 1
157.924415 .00100000 1.0000E-5 0 2
159.925204 .02340000 1.0000E-5 0 3
160.926939 .18910000 1.0000E-5 0 4 5
161.926805 .25510000 1.0000E-5 0 6 7 8
162.928737 .24900000 1.0000E-5 0 910
163.929181 .28180000 1.0000E-5 0111213

RPI RESOLUTION PARAMETERS FOLLOW
BURST 0      60.      0.1
CHANN 0      44.68    500.000    0.100
CHANN 0      262.17    62.500    0.100
CHANN 0      1998.1    31.250    0.100

MISCELLANEOUS PARAMETERS FOLLOW
DELTE 0 00 -1.27-6  4.49-6    0.187    0.063    -.021    0.003

```

Figure 41. SAMMY input for $n+^{nat}\text{Dy}$ RPI's epithermal transmission data.

```

Dysprosium / transmission / rpi data / Original filename Dy_ttnat_20mil_CSISRS.t
wenty
Dynat      162.499472  0.00001 2.0000E+01    1    1 0 0 0 0    0
FGM
TWENTY
EV
GENERATE PLOT FILE AUTOMATICALLY
PRINT THEORETICAL VALUES
PRINT DEBUG INFORMATION
DO NOT SOLVE BAYES EQUATIONS
REICH-MOORE FORMALISM IS WANTED
QUANTUM NUMBERS ARE IN PARAMETER FILE
DO NOT SHIFT RPI RESOLUTION FUNCTION TO CENTER
ENERGY UNCERTAINTIES ARE AT END OF LINE
USE NO CUTOFF FOR DERIVATIVES OR CROSS SECTIONS
#IMPLICIT DATA COVARIANCE MATRIX IS INCLUDED

    293.6    14.973    0.0055    0.0000    0.0
    7.50000  1.610e-3
TRANSMISSION

NUCLIDE MASSES AND ABUNDANCES FOLLOW
155.924284 .00060000 1.0000E-5 0 1
157.924415 .00100000 1.0000E-5 0 2
159.925204 .02340000 1.0000E-5 0 3
160.926939 .18910000 1.0000E-5 0 4 5
161.926805 .25510000 1.0000E-5 0 6 7 8
162.928737 .24900000 1.0000E-5 0 910
163.929181 .28180000 1.0000E-5 0111213

RPI RESOLUTION PARAMETERS FOLLOW
BURST 0 2100. 21.
TAU 00000320.936 0.02426 318.461 0.02922 235.721
TAU 00 82.641 0.00625 82.004 0.00752 60.698
LAMBD00000 679.984 -226.205 21.009
LAMBD 146.197 48.634 4.517
A1 00000 -.000780 0.02407 -.000630 3.531 0.00112
A1 .00005 0.0 .0 0.0 0.00006
EXPON00000 935.91 -69.068 .005 0.39485 0.00075
EXPON 0.801 0.396 .00001 0.00042 0.00002
CHANN 0 0.1096 32000. 0.100
CHANN 0 5.301 2000. 0.100
CHANN 0 19.94 500. 0.100

NORMALization and "constant" background follow
1.00671608-.00625250-.00045116 0. 0. 0 0 0 0 0 0
.010000000 .00100000 .00100000 0. 0. 0.

```

Figure 42. SAMMY input for $n+^{nat}\text{Dy}$ RPI's thermal transmission data.

APPENDIX B. Resonance parameters

The SAMMY parameters file containing the particle-pair definitions, spin group information, and resonance parameters each Dy isotope is shown in Fig. 43. The multiplication factors to calibrate the evaluated uncertainties are reported in Fig. 44. The details on the format of both files can be found in the SAMMY manual [9].

PARTICLE PAIR DEFINITIONS									
Name=156Dy+n	Particle a=neutron	Particle b=156Dy	Shift=0	155.92428359					
Za= 0	Sb= 0.0	Ma=	Particle b=158Dy						
Na=158Dy+n	Particle a=neutron	Particle b=158Dy	Shift=0	157.92441482					
Za= 0	Sb= 0.0	Ma=	Particle b=160Dy						
Na=160Dy+n	Particle a=neutron	Particle b=160Dy	Shift=0	159.92520358					
Za= 0	Sb= 0.0	Ma=	Particle b=162Dy						
Name=162Dy+n	Particle a=neutron	Particle b=162Dy	Shift=0	160.92693943					
Za= 0	Sb= 2.5	Ma=	Particle b=164Dy						
Na=164Dy+n	Particle a=neutron	Particle b=164Dy	Shift=0	161.92680451					
Za= 0	Sb= 0.0	Ma=	Particle b=163Dy						
Name=163Dy+n	Particle a=neutron	Particle b=163Dy	Shift=0	162.92873722					
Za= 0	Sb= -2.5	Ma=	Particle a=neutron	Shift=0	163.92918082				
Name=164Dy+n	Particle a=neutron	Particle b=164Dy	Shift=0						
Za= 0	Sb= 0.0	Ma=							
Sa= 0.5									
SPIN GROUPS									
1 1 156Dy+n	0 0.5 0.006000								
2 1 158Dy+n	0.5 0.001000								
3 1 160Dy+n	0 0.5								
4 1 162Dy+n	0 0.5								
5 1 164Dy+n	0 2.0 0.189100								
6 1 161Dy+n	0 3.0 0.189100								
7 1 162Dy+n	0 0.5 0.255100								
8 1 162Dy+n	0 0.5								
9 1 162Dy+n	-1.5 0.255100								
10 1 163Dy+n	-2.0 0.249000								
11 1 164Dy+n	-3.0 0.249000								
12 1 164Dy+n	0 0.5								
13 1 164Dy+n	-1.5 0.281800								
14 1 164Dy+n	1 0.5								
RESONANCES are listed next									
2.150000000 83.6000000 .200000000									
3.210000000 83.6000000 .800000000									
9.190000000 83.6000000 .600000000									
17.240000000 83.6000000 .430000000									
17.700000000 83.6000000 .100000000									
19.600000000 83.6000000 2.400000000									
24.500000000 83.6000000 2.400000000									
27.400000000 83.6000000 .400000000									
28.100000000 83.6000000 .400000000									
29.300000000 83.6000000 9.700000000									
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1862.719000	105.836500	1295.73800	3	3	3
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3513.800000	116.800000	13.00000000	1	1	1	8
3566.900000	116.800000	13.00000000	1	1	1	8
-521.739000	120.171300	4151.47100	1	1	1	10
-464829800	120.567100	311924400	1	1	1	10
1.718878617	101.580971	2.10009344	1	1	1	9
5.896487000	108.516800	.032793660	1	1	1	9
6.257694000	120.879400	.040264060	1	1	1	9
16.24160000	62.2511200	29.0239100	1	1	1	10
19.68895000	114.427100	1.02464100	1	1	1	10
35.80051000	107.274000	13.5988700	1	1	1	9
50.26940000	108.335200	3.35220300	1	1	1	10
55.93295000	136.524100	26.8833800	1	1	1	10
59.05320000	104.516800	106.813100	1	1	1	9
66.14400000	108.445700	8.54483500	1	1	1	10
72.04150000	108.547600	4.36537300	1	1	1	9
73.42878000	108.366800	3.65031100	1	1	1	9
82.30282000	108.423100	2.50594500	1	1	1	10
94.14237000	108.223200	19.3890700	1	1	1	10
105.8724000	118.880800	66.8361300	1	1	1	10
107.2066000	110.030900	39.4068200	1	1	1	9
120.3983000	108.523900	9.74491800	1	1	1	10
126.6425000	108.668600	16.8567400	1	1	1	10
127.5046000	108.623500	13.2406900	1	1	1	10
135.3777000	108.585400	5.72828900	1	1	1	10
143.4806000	108.545200	16.7510000	1	1	1	9
145.0107000	103.754600	54.2522200	1	1	1	9
155.1077000	99.9663100	106.001300	1	1	1	9

163.8931000	108.528800	11.2000800	1	1	1	10
177.2511000	108.583200	3.73201200	1	1	1	10
185.2036000	108.572900	4.35937200	1	1	1	10
189.0368000	108.596400	4.22452900	1	1	1	9
202.9841000	105.962500	22.1691000	1	1	1	9
205.3569000	100.329900	54.1867200	1	1	1	10
213.8396000	108.601300	6.35007600	1	1	1	10
223.6205000	101.166300	190.920400	1	1	1	10
224.6202000	108.186600	6.10448800	1	1	1	10
225.6692000	108.600200	6.17839400	1	1	1	10
230.4300000	108.600800	15.7035000	1	1	1	10
261.3150000	99.9760100	107.625800	1	1	1	10
268.1567000	98.8800700	133.537000	1	1	1	10
274.3579000	115.936100	37.4991900	1	1	1	10
281.1887000	110.552300	27.8244200	1	1	1	10
289.0532000	94.7451200	145.695200	1	1	1	10
295.9477000	108.595700	7.85981800	1	1	1	9
297.3158000	108.267900	12.9405000	1	1	1	9
298.0892000	92.5848900	92.5263000	1	1	1	10
323.0735000	103.616600	255.973200	1	1	1	10
324.7490000	100.999300	260.120200	1	1	1	10
327.1781000	108.711000	33.9232300	1	1	1	10
329.6652000	108.612000	19.2883500	1	1	1	9
343.0807000	108.634700	20.0278300	1	1	1	10
348.3942000	93.9921100	239.130300	1	1	1	10
358.4392000	108.613000	13.1804100	1	1	1	10
375.4251000	108.596400	3.15806400	1	1	1	10
382.4495000	108.593000	6.09037200	1	1	1	10
387.2756000	112.142900	32.9710600	1	1	1	10
390.6223000	89.9603700	45.0655000	1	1	1	10
400.3766000	109.384600	34.5865600	1	1	1	10
403.0539000	108.596500	4.04289500	1	1	1	9
411.3714000	175.671000	629.927600	1	1	1	10
420.8477000	108.599400	2.50380900	1	1	1	10
429.6393000	67.3738000	36.1346700	1	1	1	10
455.1084000	147.802200	75.8308700	1	1	1	10
459.5126000	107.763600	22.4578900	1	1	1	10
465.5376000	108.608300	20.6512800	1	1	1	10
479.4098000	148.501900	85.0241800	1	1	1	9
483.9689000	126.141200	94.8026900	1	1	1	9
490.3730000	108.596500	2.50380900	1	1	1	10
504.4627000	108.524800	22.4532500	1	1	1	9
505.2247000	62.0862800	127.122300	1	1	1	10
516.5177000	114.002800	25.8456500	1	1	1	10
519.3703000	101.205100	19.5291000	1	1	1	10
520.4253000	108.262300	14.1011200	1	1	1	9
533.6283000	122.047300	248.793200	1	1	1	10
542.6806000	125.148800	342.624300	1	1	1	10
564.6313000	87.3671500	76.6732200	1	1	1	9
569.4025000	157.131700	157.703300	1	1	1	10
581.1365000	126.213300	73.2086300	1	1	1	10
594.8218000	112.358600	86.8534700	1	1	1	10
601.5340000	143.858200	199.544400	1	1	1	10
616.2194000	73.0699600	57.2289700	1	1	1	10
621.0201000	122.831600	134.527100	1	1	1	10
637.5430000	135.689000	115.689700	1	1	1	10
643.4141000	77.2175900	111.842200	1	1	1	10
646.7747000	151.128800	102.098300	1	1	1	10
652.6199000	108.600700	7.4035900	1	1	1	10
666.8271000	108.265200	49.4645200	1	1	1	10
685.2264000	108.272300	6.61804100	1	1	1	10
687.5984000	108.597900	22.2573000	1	1	1	10
693.9699000	108.270100	22.5901300	1	1	1	10
696.3310000	118.939000	199.856200	1	1	1	9
710.3155000	108.601700	10.1011900	1	1	1	9
712.8034000	94.2501700	90.3537400	1	1	1	9

714.1198000	108.265900	20.7607000	1	1	1	9
721.7421000	75.1755000	73.0770300	1	1	1	10
732.4391000	108.602900	17.0294400	1	1	1	10
736.4246000	108.599500	17.3933000	1	1	1	10
742.4096000	108.596200	19.7738600	1	1	1	10
747.2669000	108.599300	8.78106200	1	1	1	9
756.4203000	82.3923000	41.7097600	1	1	1	10
758.4774000	108.270700	13.1859900	1	1	1	9
764.5082000	108.600100	22.0335300	1	1	1	9
767.4592000	108.608400	14.1231700	1	1	1	9
770.4300000	108.600800	15.7035000	1	1	1	10
776.7194000	107.978100	23.2317200	1	1	1	10
794.5185000	110.023300	47.4432900	1	1	1	10
796.0830000	108.276100	23.0774300	1	1	1	9
810.2264000	138.754500	362.761200	1	1	1	9
823.1801000	117.840500	69.7845200	1	1	1	9
823.4070000	109.262600	41.9144200	1	1	1	10
831.5636000	109.685200	105.423400	1	1	1	10
834.9037000	108.272200	11.4485000	1	1	1	10
845.9350000	108.597600	11.2424300	1	1	1	10
850.4771000	108.265500	24.1206400	1	1	1	9
851.9654000	108.472900	192.603700	1	1	1	9
854.4091000	112.619100	75.3143100	1	1	1	9
864.0582000	108.265300	26.6709100	1	1	1	10
869.6601000	103.151200	59.3741200	1	1	1	9
880.4392000	108.600800	15.7035000	1	1	1	10
900.7551000	97.278100	210.223300	1	1	1	9
919.1644000	105.014100	73.9816300	1	1	1	10
920.7148000	108.268400	41.1907200	1	1	1	10
930.4786000	98.3948900	240.323000	1	1	1	10
936.0401000	108.791000	32.5148300	1	1	1	9
940.1376000	86.3417000	208.223800	1	1	1	10
949.9583000	103.472200	169.591600	1	1	1	10
954.6061000	108.272400	15.3286100	1	1	1	9
958.1560000	109.495500	31.2926600	1	1	1	10
967.4671000	111.397700	60.7538500	1	1	1	9
973.6790000	120.808200	295.082300	1	1	1	10
980.8597000	108.599500	11.3241100	1	1	1	9
997.3068000	108.807300	66.3128800	1	1	1	10
1007.240000	120.007300	239.312800	1	1	1	9
1102.430000	108.600800	15.7035000	1	1	1	9
1175.269000	120.001100	78.8319500	1	1	1	10
1206.580000	120.087400	8631.93300	1	1	1	9
-2037.04900	83.6000100	46629.7700	1	1	1	11
-102.499200	83.6000500	310.255600	1	1	1	11
-1.87976800	63.5308900	50.065500	1	1	1	11
147.1067000	90.4412600	832.878300	1	1	1	11
450.5417000	72.6695300	268.867800	1	1	1	11
536.6550000	63.5467500	106.479500	1	1	1	11
549.1917000	114.237000	6.9076800	1	1	1	11
854.3820000	82.3533600	520.078800	1	1	1	11
983.7230000	110.415000	105.845300	1	1	1	11
1052.182000	105.091300	230.530700	1	1	1	11
1207.943000	114.332300	666.624300	1	1	1	11
1364.068000	114.464800	924.569300	1	1	1	11
1482.782000	113.902200	220.033300	1	1	1	11
1897.089000	114.328600	2037.20200	1	1	1	11
1960.861000	114.303300	1229.56600	1	1	1	11
2065.076000	114.208200	46.1358700	1	1	1	11
2320.220000	114.276300	1905.32200	1	1	1	11
2397.211000	114.236800	421.286600	1	1	1	11
2537.213000	114.242000	659.521700	1	1	1	11
2871.766000	114.278200	2496.20800	1	1	1	11
2970.108000	114.248000	3154.50500	1	1	1	11
3281.889000	114.204100	219.924000	1	1	1	11
3481.973000	114.209600	1876.38600	1	1	1	11

3734.771000	114.206200	727.898200	1	1	1	11
4135.211000	114.200200	488.227100	1	1	1	11
4231.531000	114.200600	2168.07800	1	1	1	11
4445.840000	114.200900	4862.07600	1	1	1	11
4757.601000	114.200000	139.909700	1	1	1	11
4780.107000	114.200200	1489.77900	1	1	1	11
4810.150000	114.201400	10285.3900	1	1	1	11
4954.901000	114.200000	169.879600	1	1	1	11
5161.900000	114.200200	2083.38400	1	1	1	11
5201.900000	114.201200	1352.8100	1	1	1	11
5202.501000	114.200000	239.823800	1	1	1	11
5223.319000	114.200600	10190.7300	1	1	1	11
6096.803000	114.200100	1791.34900	1	1	1	11
6109.803000	114.200100	2088.27800	1	1	1	11
6280.710000	114.200300	6780.55200	1	1	1	11
6444.703000	114.200100	2287.32300	1	1	1	11
6621.105000	114.200200	3767.16400	1	1	1	11
6854.201000	114.200000	1196.92600	1	1	1	11
6996.301000	114.200000	997.944500	1	1	1	11
7102.111000	100.000000	1132.60200	1	1	1	11
9015.633000	100.000000	98653.6500	1	1	1	11
804.240000	114.200000	8.80000000	1	1	1	12
940.980000	114.200000	2.45000000	1	1	1	12
1286.500000	114.200000	4.30000000	1	1	1	12
1405.800000	114.200000	3.20000000	1	1	1	13
1567.400000	114.200000	4.35000000	1	1	1	13
2039.000000	114.200000	9.90000000	1	1	1	12
2250.700000	114.200000	14.50000000	1	1	1	13
2285.100000	114.200000	31.00000000	1	1	1	13
2352.600000	114.200000	8.00000000	1	1	1	12
2723.200000	114.200000	27.00000000	1	1	1	12
2804.400000	114.200000	8.00000000	1	1	1	13
3048.200000	114.200000	8.50000000	1	1	1	13
3100.000000	114.200000	9.50000000	1	1	1	13
3209.000000	114.200000	26.50000000	1	1	1	13
3416.800000	114.200000	23.50000000	1	1	1	13
3460.500000	114.200000	28.00000000	1	1	1	13
3520.300000	114.200000	71.00000000	1	1	1	12
3521.000000	114.200000	17.00000000	1	1	1	12
3911.600000	114.200000	17.00000000	1	1	1	13
4095.800000	114.200000	30.00000000	1	1	1	13
4348.800000	114.200000	25.50000000	1	1	1	13
4389.200000	114.200000	14.50000000	1	1	1	13
4916.000000	114.200000	42.00000000	1	1	1	13

1.000000000

Channel RadII in key-word format

RadII= 7.500000, 7.500000

Flags=0,-1

Uncertainties= 0.037500, 7.500000

Group= 1

Chan= 1,

Flags=0,-1

RadII= 7.400000, 7.400000

Flags=0,-1

Uncertainties= 0.037000, 7.400000

Group= 2

Chan= 1,

Flags=0,-1

RadII= 7.4561270 7.4561270

Flags=0,-1

Uncertainties= 0.040000, 7.4561270

Group= 3

Chan= 1,

Flags=0,-1

RadII= 7.469997, 7.469997

Flags=0,-1

Uncertainties= 0.040000, 7.469997

Group= 4

Chan= 1,

Flags=0,-1

RadII= 5.900000, 5.900000

Flags=0,-1

Uncertainties= 0.030000, 5.900000

Group= 6

Chan= 1,

Flags=0,-1

Group= 7

Chan= 1,

Flags=0,-1

Group= 8

Chan= 1,

Flags=0,-1

RadII= 7.497540, 7.497540	Flags=0,-1
Uncertainties= 0.040000, 7.497540	
Group= 9	Chan= 1, Flags=0,-1
RadII= 7.511213, 7.511213	Flags=0,-1
Uncertainties= 0.036553, 7.511213	
Group= 11	Chan= 1, Flags=0,-1
Group= 12	Chan= 1, Flags=0,-1
Group= 13	Chan= 1, Flags=0,-1

Prior uncertainties in key-word format

Relative uncertainty = 0.10

Group=1

Ch=1

Relative uncertainty = 0.10

Group=2

Ch=1

Relative uncertainty = 0.10

Group=3

Ch=1

Relative uncertainty = 0.10

Group=4

Ch=1

Relative uncertainty = 0.10

Group=5

Ch=1

Relative uncertainty = 0.10

Group=6

Ch=1

Relative uncertainty = 0.10

Group=7

Ch=1

Relative uncertainty = 0.10

Group=8

Ch=1

Relative uncertainty = 0.10

Group=9

Ch=1

Relative uncertainty = 0.10

Group=10

Ch=1

Relative uncertainty = 0.10

Group=11

Ch=1

Relative uncertainty = 0.10

Group=12

Ch=1

Relative uncertainty = 0.10

Group=13

Ch=1

Relative uncertainty = 0.10

pp=capture

Relative uncertainty = 0.10

pp=energy

Figure 43. SAMMY parameter file containing particle-pair definition for seven Dy isotopes and related quantum number information. Resonance parameters (energy and resonance widths) are also shown.

```

Multiplication factor to increase uncertainty from 1.e-5 up 0.5 eV
factors defined below for each isotope.
spin group = 1
0.0      1.
0.5      1.

spin group = 2
0.0      1.
0.5      1.

spin group = 3
0.0      1.
0.5      1.

spin group = 4
0.0      3.
10.      3.

spin group = 5
0.0      3.
10.      3.

spin group = 6
0.0      5.
10.      5.

spin group = 7
0.0      1.
0.5      1.

spin group = 8
0.0      1.
0.5      1.

spin group = 9
-3000.0  2.
0.0      2.

spin group = 10
-3000.0  2.
0.0      2.

spin group = 11
-3000.0  10.
0.        10.

spin group = 12
0.0      1.
0.5      1.

spin group = 13
0.0      1.
0.5      1.

```

Figure 44. SAMMY energy dependent uncertainty file containing the multiplication factors applied on a specific energy range defined by the energy levels of a specific isotope.