				Spe Spe
PROGRAM S	SPECT	'RA		Spe
		==		Spe
An extens	sion	of LINEAR to 1	inearize ALl MF=5 spectra.	Spe
			utron induced, photon spectra.	Spe
			•	Spe
First rel	Lease	ed in 2010 - Ea	rlier below dates refer to LINEAR.	Spe
				Spe
VERSION 7	74-1	(MAY 1974)		Spe
VERSION 7	75-1	(APRIL 1975)		Spe
VERSION 7	76-2	(OCTOBER 1976)		Spe
VERSION 7	77-1	(JANUARY 1977)		Spe
		(JULY 1978)		Spe
		•	C-7600 AND CRAY-1 VERSION.	Spe
			, CDC AND CRAY VERSION.	Spe
		(DECEMBER 1980)	Spe
		(MARCH 1981)		Spe
		-	IMPROVED COMPUTER COMPATIBILITY.	Spe
VERSION 8	33-1	(JANUARY 1983)	*MAJOR RE-DESIGN.	Spe
			*PAGE SIZE INCREASED - 1002 TO 3006.	Spe
			*ELIMINATED COMPUTER DEPENDENT CODING.	-
			*NEW, MORE COMPATIBLE I/O UNIT NUMBER.	-
			*ADDED OPTION TO KEEP ALL ORIGINAL	Spe
			ENERGY POINTS FROM EVALUATION.	Spe
			*ADDED STANDARD ALLOWABLE ERROR OPTION	_
TERRITOR O		(OCHODED 1003)	(CURRENTLY 0.1 PER-CENT).	Spe
		•	IMPROVED BASED ON USER COMMENTS.	Spe
		(APRIL 1984)	IMPROVED BASED ON USER COMMENTS.	Spe
VERSION 6	94-2	(JUNE 1984)	*UPDATED FOR ENDF/B-VI FORMATS. *SPECIAL I/O ROUTINES TO GUARANTEE	Spe
			ACCURACY OF ENERGY.	Spe
			*DOUBLE PRECISION TREATMENT OF ENERGY	Spe
			(REQUIRED FOR NARROW RESONANCES).	Spe
VERSTON 8	35-1	(AUGUST 1985)	*FORTRAN-77/H VERSION	Spe
			*ENDF/B-VI FORMAT	Spe
			*DOUBLE PRECISION TREATMENT OF CROSS	Spe
		(0,	SECTION	Spe
VERSION 8	88-1	(JULY 1988)	*OPTIONINTERNALLY DEFINE ALL I/O	Spe
			FILE NAMES (SEE, SUBROUTINE FILEIO	Spe
			FOR DETAILS).	Spe
			*IMPROVED BASED ON USER COMMENTS.	Spe
VERSION 8	39-1	(JANUARY 1989)	*PSYCHOANALYZED BY PROGRAM FREUD TO	Spe
			INSURE PROGRAM WILL NOT DO ANYTHING	Spe
			CRAZY.	Spe
			*UPDATED TO USE NEW PROGRAM CONVERT	Spe
			KEYWORDS.	Spe
			*ADDED LIVERMORE CIVIC COMPILER	Spe
			CONVENTIONS.	Spe
VERSION 9	90-1	(JUNE 1990)	*EXTENDED TO LINEARIZE PHOTON	Spe
			INTERACTION DATA, MF=23 AND 27	Spe
			*ADDED FORTRAN SAVE OPTION	Spe
			*UPDATED BASED ON USER COMMENTS.	Spe
			*NEW MORE CONSISTENT ENERGY OUTPUT	Spe
			ROUTINE.	Spe
			*WARNINGINPUT PARAMETER FORMAT	Spe
			HAS BEEN CHANGEDSEE DESCRIPTION	Spe
			BELOW.	Spe
version 9	91-1	(JULY 1991)	*ADDED INTERPOLATION LAW 6 - ONLY USED	-
			FOR CHARGED PARTICLE CROSS SECTIONS	Spe
			WIND CONTRAME DEMEMBARTTTMTEC	Spe
		/	FOR COULOMB PENETRABILITIES. *ADDED NU-BAR (TOTAL, DELAYED, PROMPT)	_

		TO LINEARLY INTERPOLABLE	Spectra
	•	*INCREASED PAGE SIZE FROM 3006 TO 5010	Spectra
		POINTS.	Spectra
	•	*ALL ENERGIES INTERNALLY ROUNDED PRIOR	Spectra
		TO CALCULATIONS.	Spectra
	•	*COMPLETELY CONSISTENT I/O AND ROUNDING	Spectra
		ROUTINES - TO MINIMIZE COMPUTER	Spectra
		DEPENDENCE.	Spectra
VERSION 92-2	(JULY 1992)	*CORRECTED CONVERSION OF NU-BAR FROM	Spectra
		POLYNOMIAL TO TABULATED - COPY	Spectra
		SPONTANEOUS NU-BAR (BY DEFINITION	Spectra
		THE SPONTANEOUS NU-BAR IS NOT AN	Spectra
		ENERGY DEPENDENT QUANTITY).	Spectra
VERSION 93-1	(MARCH 1993)	*UPDATED FOR USE WITH LAHEY COMPILER	Spectra
		ON IBM-PCS.	Spectra
		*INCREASED PAGE SIZE FROM 5010 TO	Spectra
		30000 POINTS	Spectra
VERSION 94-1	(JANUARY 1994)	*VARIABLE ENDF/B DATA FILENAMES	Spectra
		TO ALLOW ACCESS TO FILE STRUCTURES	Spectra
		(WARNING - INPUT PARAMETER FORMAT	Spectra
		HAS BEEN CHANGED)	Spectra
		*CLOSE ALL FILES BEFORE TERMINATING	Spectra
		(SEE, SUBROUTINE ENDIT)	Spectra
VERSION 96-1	(JANUARY 1996)	*COMPLETE RE-WRITE	Spectra
		*IMPROVED COMPUTER INDEPENDENCE	Spectra
		*ALL DOUBLE PRECISION	Spectra
		*ON SCREEN OUTPUT	Spectra
		*UNIFORM TREATMENT OF ENDF/B I/O	Spectra
		*IMPROVED OUTPUT PRECISION	Spectra
		*DEFINED SCRATCH FILE NAMES	Spectra
		*ALWAYS INCLUDE THERMAL VALUE	Spectra
		*INCREASED PAGE SIZE FROM 30000 TO	Spectra
		60000 POINTS	Spectra
VERSION 99-1	(MARCH 1999)	*CORRECTED CHARACTER TO FLOATING	Spectra
	,	POINT READ FOR MORE DIGITS	Spectra
		*UPDATED TEST FOR ENDF/B FORMAT	Spectra
		VERSION BASED ON RECENT FORMAT CHANGE	
		*GENERAL IMPROVEMENTS BASED ON	Spectra
		USER FEEDBACK	Spectra
VERSION 99-2	(JUNE 1999)	*ASSUME ENDF/B-VI, NOT V, IF MISSING	Spectra
	(00000	MF=1, MT-451.	Spectra
VERS. 2000-1	(FEBRUARY 2000)*ADDED MF = 9 AND 10 LINEARIZATION	Spectra
	(*GENERAL IMPROVEMENTS BASED ON	Spectra
		USER FEEDBACK	Spectra
VERS. 2002-1	(MAY 2002)	*OPTIONAL INPUT PARAMETERS	Spectra
VERS. 2004-1	-	*GENERAL UPDATE BASED ON USER FEEDBACK	_
VERS. 2005-1	= = = = = = = = = = = = = = = = = = = =	*ALWAYS KEEP ORIGINAL TABULATED	Spectra
		NU-BAR POINTS.	Spectra
VERS. 2006-1	(FEB. 2006)	*CORRECTED INT=6 NEAR THRESHOLD	Spectra
	, ,	*NO SUBDIVIDE BELOW MINIMUM XCMIN	Spectra
VERS. 2007-1	(JAN. 2007)	*CHECKED AGAINST ALL ENDF/B-VII.	Spectra
	,,,	*INCREASED PAGE SIZE FROM 60,000 TO	Spectra
		600,000 POINTS	Spectra
VERS. 2010-1	(JUNE 2010)	*ADDED MF = 5 - MF = 6 STILL PLANNED.	Spectra
	,50112 2020/	*72 CHARACTER FILE NAMES.	Spectra
		*ONLY PROCESS MF=5 - SKIP ALL OTHERS	Spectra
		TO PREVENT CONFLICT WITH LINEAR	Spectra
		THINNING.	Spectra
VERS. 2012-1	(Aug 2012)	*Added MF=15, neutron induced photon	Spectra
· LI.W. 2012-1	(ug. 2012)	spectra.	Spectra
		*Added CODENAME	Spectra
		*32 and 64 bit Compatible	Spectra
		*Added ERROR stop	Spectra
		HAUGE BREOK SCOP	Specta

VERS. 2015-1 (Jan. 2015) *Extended OUT9.	Spectra
*Replaced ALL 3 way IF Statements.	Spectra
*Corrected MF=15 Data - it was adding	Spectra
SEND between sub-sections.	Spectra
*Deleted unused parts, e.g., NUBAR.	Spectra
VERS. 2017-1 (May 2017) *Increased page size to 3,000,000	Spectra
*Updated based on user feedback	Spectra
*Changed DGAMMA to REDGAMMA to avoid	Spectra
conflict with possble RESERVED NAME	Spectra
*All floating input parameters changed	Spectra
to character input + IN9 conversion.	Spectra
	Spectra
OWNED, MAINTAINED AND DISTRIBUTED BY	Spectra
	Spectra
THE NUCLEAR DATA SECTION	Spectra
INTERNATIONAL ATOMIC ENERGY AGENCY	Spectra
P.O. BOX 100	Spectra
A-1400, VIENNA, AUSTRIA	Spectra
EUROPE	Spectra
	Spectra
ORIGINALLY WRITTEN BY	Spectra
	Spectra
Dermott E. Cullen	Spectra
DDECEME COMMENT TATODAY TO	Spectra
PRESENT CONTACT INFORMATION	Spectra
Describb B. Calles	Spectra
Dermott E. Cullen	Spectra
1466 Hudson Way	Spectra
Livermore, CA 94550 U.S.A.	Spectra
	Spectra
Telephone 925-443-1911 E. Mail RedCullen1@Comcast.net	Spectra
Website RedCullen1.net/HOMEPAGE.NEW	Spectra Spectra
Website Redcuiteni.net/nomerage.wew	Spectra
AUTHORS MESSAGE	Spectra
	Spectra
THE REPORT DESCRIBED ABOVE IS THE LATEST PUBLISHED DOCUMENTATION	_
	Spectra
FOR THIS PROGRAM. HOWEVER. THE COMMENTS BELOW SHOULD BE CONSIDERED.	Spectra Spectra
FOR THIS PROGRAM. HOWEVER, THE COMMENTS BELOW SHOULD BE CONSIDERED THE LATEST DOCUMENTATION INCLUDING ALL RECENT IMPROVEMENTS. PLEASE	Spectra
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OPPOSED TO THE BINARY FORMAT) AND CAN HANDLE DATA IN ANY VERSION OF THE ENDF/B FORMAT (I.E., ENDF/B-I, II, III, IV, V OR VI FORMAT). Spectra

IT IS ASSUMED THAT THE DATA IS CORRECTLY CODED IN THE ENDF/B FORMAT AND NO ERROR CHECKING IS PERFORMED. IN PARTICULAR IT IS ASSUMED THAT THE MAT, MF AND MT ON EACH LINE IS CORRECT. SEQUENCE NUMBERS (COLUMNS 76-80) ARE IGNORED ON INPUT, BUT WILL BE CORRECTLY OUTPUT ON ALL LINES. THE FORMAT OF SECTION MF=1, MT=451 AND ALL SECTIONS OF MF=3 MUST BE CORRECT. THE PROGRAM COPIES ALL OTHER SECTION OF DATA AS HOLLERITH AND AS SUCH IS INSENSITIVE TO THE CORRECTNESS OR INCORRECTNESS OF ALL OTHER SECTIONS.

OUTPUT FORMAT

IN THIS VERSION OF LINEAR ALL ENERGIES WILL BE OUTPUT IN F (INSTEAD OF E) FORMAT IN ORDER TO ALLOW ENERGIES TO BE WRITTEN WITH UP TO 9 DIGITS OF ACCURACY. IN PREVIOUS VERSIONS THIS WAS AN OUTPUT OPTION. HOWEVER USE OF THIS OPTION TO COMPARE THE RESULTS OF ENERGIES WRITTEN IN THE NORMAL ENDF/B CONVENTION OF 6 DIGITS TO THE 9 DIGIT OUTPUT FROM THIS PROGRAM DEMONSTRATED THAT FAILURE TO USE THE 9 DIGIT OUTPUT CAN LEAD TO LARGE ERRORS IN THE DATA DUE TO TRUNCATION OF ENERGIES TO 6 DIGITS DURING OUTPUT.

CONTENTS OF OUTPUT

ENTIRE EVALUATIONS ARE OUTPUT, NOT JUST THE LINEARIZED DATA CROSS SECTIONS, E.G. ANGULAR AND ENERGY DISTRIBUTIONS ARE ALSO INCLUDED.

DOCUMENTATION

THE FACT THAT THIS PROGRAM HAS OPERATED ON THE DATA IS DOCUMENTED BY THE ADDITION OF 3 COMMENT LINES AT THE END OF EACH HOLLERITH SECTION IN THE FORM

******** PROGRAM SPECTRA (2017-1) *********** FOR ALL DATA GREATER THAN 1.00000-10 IN ABSOLUTE VALUE DATA LINEARIZED TO WITHIN AN ACCURACY OF 0.1 PER-CENT

THE ORDER OF SIMILAR COMMENTS (FROM RECENT, SIGMA1 AND GROUPIE) REPRESENTS A COMPLETE HISTORY OF ALL OPERATIONS PERFORMED ON THE DATA BY THESE PROGRAMS.

THESE COMMENT LINES ARE ONLY ADDED TO EXISTING HOLLERITH SECTIONS, Spectra I.E., THIS PROGRAM WILL NOT CREATE A HOLLERITH SECTION. THE FORMAT Spectra OF THE HOLLERITH SECTION IN ENDF/B-V DIFFERS FROM THE THAT OF EARLIER VERSIONS OF ENDF/B. BY READING AN EXISTING MF=1, MT=451 IT IS POSSIBLE FOR THIS PROGRAM TO DETERMINE WHICH VERSION OF THE ENDF/B FORMAT THE DATA IS IN. WITHOUT HAVING A SECTION OF MF=1, MT=451 PRESENT IT IS IMPOSSIBLE FOR THIS PROGRAM TO DETERMINE WHICH VERSION OF THE ENDF/B FORMAT THE DATA IS IN, AND AS SUCH IT IS IMPOSSIBLE FOR THE PROGRAM TO DETERMINE WHAT FORMAT SHOULD BE USED TO CREATE A HOLLERITH SECTION.

REACTION INDEX

THIS PROGRAM DOES NOT USE THE REACTION INDEX WHICH IS GIVEN IN SECTION MF=1, MT=451 OF EACH EVALUATION.

THIS PROGRAM DOES NOT UPDATE THE REACTION INDEX IN MF=1, MT=451. THIS CONVENTION HAS BEEN ADOPTED BECAUSE MOST USERS DO NOT REQUIRE A CORRECT REACTION INDEX FOR THEIR APPLICATIONS AND IT WAS Spectra NOT CONSIDERED WORTHWHILE TO INCLUDE THE OVERHEAD OF CONSTRUCTING Spectra

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A CORRECT REACTION INDEX IN THIS PROGRAM. HOWEVER, IF YOU REQUIRE Spectra A REACTION INDEX FOR YOUR APPLICATIONS, AFTER RUNNING THIS PROGRAM Spectra YOU MAY USE PROGRAM DICTIN TO CREATE A CORRECT REACTION INDEX. Spectra

SECTION SIZE

SINCE THIS PROGRAM USES A LOGICAL PAGING SYSTEM THERE IS NO LIMIT TO THE NUMBER OF POINTS IN ANY SECTION, E.G., THE TOTAL CROSS SECTION MAY BE REPRESENTED BY 200,000 DATA POINTS.

FOR ANY LINEARIZED SECTION THAT CONTAINS 60000 OR FEWER POINTS Spectra
THE ENTIRE OPERATION WILL BE PERFORMED IN CORE AND THE LINEARIZED Spectra
DATA WILL BE OUTPUT DIRECTLY TO THE ENDF/B FORMAT. FOR ANY SECTION Spectra
THAT CONTAINS MORE POINTS THE DATA WILL BE LINEARIZED A PAGE AT A
TIME (1 PAGE = 60000 POINTS) AND OUTPUT TO SCRATCH. AFTER THE SPECTRA
ENTIRE SECTION HAS BEEN LINEARIZED THE DATA WILL BE READ BACK FROM SPECTRA
SCRATCH AND OUTPUT TO THE ENDF/B FORMAT. Spectra

SELECTION OF DATA

THE PROGRAM SELECTS DATA TO BE LINEARIZED BASED EITHER ON EITHER MAT (ENDF/B MAT NO.) OR ZA AS WELL AS MF AND MT NUMBERS. THIS PROGRAM ALLOWS UP TO 100 MAT/MF/MT OR ZA/MF/MT RANGES TO BE SPECIFIED BY INPUT PARAMETERS. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN MAT ORDER, REGARDLESS OF THE CRITERIA USED TO RETRIEVE MATERIALS. IF RETRIEVAL IS BY MAT RANGE THE PROGRAM WILL TERMINATE WHEN A MAT IS FOUND THAT IS ABOVE ALL REQUESTED MAT RANGES. IF RETRIEVAL IS BY ZA RANGE THE PROGRAM WILL SEARCH THE ENTIRE ENDF/B TAPE.

PROGRAM OPERATION

EACH SECTION OF DATA IS CONSIDERED SEPARATELY. EACH SECTION OF Spectra ENDF/B DATA TO LINEARIZE IS REPRESENTED BY A TABLE OF ENERGY Spectra VS. CROSS SECTION AND ANY ONE OF FIVE ALLOWABLE INTERPOLATION LAWS SPECTRA BETWEEN ANY TWO TABULATED POINTS. THIS PROGRAM WILL REPLACE EACH SPECTRA SECTION OF DATA CROSS SECTIONS BY A NEW TABLE OF ENERGY VS. Spectra CROSS SECTION IN WHICH THE INTERPOLATION LAW IS ALWAYS LINEAR IN SPECTRA ENERGY AND CROSS SECTION BETWEEN ANY TWO TABULATED POINTS. Spectra

DATA IS READ AND LINEARIZED A PAGE AT A TIME (ONE PAGE CONTAINS 60000 DATA POINTS). IF THE FINAL LINEARIZED SECTION CONTAINS TWO PAGES OR LESS, DATA POINTS IT WILL BE ENTIRELY CORE RESIDENT AFTER IT HAS BEEN LINEARIZED AND WILL BE WRITTEN DIRECTLY FROM CORE TO THE OUTPUT TAPE. IF THE LINEARIZED SECTION IS LARGER THAN TWO PAGES, AFTER EACH PAGE IS LINEARIZED IT WILL BE WRITTEN TO SCRATCH. AFTER THE ENTIRE SECTION HAS BEEN LINEARIZED IT WILL BE READ BACK FROM SCRATCH, TWO PAGES AT A TIME, AND WRITTEN TO THE OUTPUT TAPE.

KEEP EVALUATED DATA POINTS

SOMETIMES IT IS CONVENIENT TO KEEP ALL ENERGY POINTS WHICH WERE PRESENT IN THE ORIGINAL EVALUATION AND TO MERELY SUPPLEMENT THESE POINTS WITH ADDITIONAL ENERGY POINTS IN ORDER TO LINEARIZE THE CROSS SECTIONS. FOR EXAMPLE, IT IS OFTEN CONVENIENT TO KEEP THE THERMAL VALUE (AT 0.0253 EV) OR THE VALUE AT 14.1 MEV.

THE CURRENT VERSION OF THIS PROGRAM WILL ALLOW THE USER TO KEEP ALL ORIGINAL EVALUATED DATA POINTS BY SPECIFYING 1 IN COLUMNS 34-44 OF THE FIRST INPUT LINE. THIS WILL TURN OFF THE BACKWARD THINNING (SEE UCRL-50400, VOL. 17, PART A FOR EXPLANATION) AND RESULT IN ALL ORIGINAL ENERGY POINTS BEING KEPT. CAUTION SHOULD

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BE EXERCISED IN USING THIS OPTION SINCE IT CAN RESULT IN A CONSIDERABLE INCREASE IN THE NUMBER OF DATA POINTS OUTPUT BY THIS CODE.

FOR ALL USERS WHO ARE NOT INTERESTED IN THIS OPTIONS NO CHANGES ARE REQUIRED IN THE INPUT TO THIS PROGRAM, I. E. IF COLUMNS 34-44 ARE BLANK (AS FOR ALL PREVIOUS VERSIONS OF THIS CODE) THE PROGRAM WILL OPERATE EXACTLY AS IT DID BEFORE.

ALLOWABLE ERROR

ALLOWABLE ERROR MUST ALWAYS BE SPECIFIED IN THE INPUT TO THIS PROGRAM AS A FRACTION, NOT A PER-CENT. FOR EXAMPLE, INPUT THE ALLOWABLE FRACTIONAL ERROR 0.001 IN ORDER TO OBTAIN DATA THAT IS ACCURATE TO WITHIN 0.1 PER-CENT.

THE CONVERSION OF THE DATA FROM THE GENERAL INTERPOLATION FORM TO Spectra LINARLY INTERPOLABLE FORM CANNOT BE PERFORMED EXACTLY. HOWEVER, IT Spectra CAN BE PERFORMED TO VIRTUALLY ANY REQUIRED ACCURACY AND MOST IMPORTANTLY CAN BE PERFORMED TO A TOLERANCE THAT IS SMALL COMPARED Spectra TO THE UNCERTAINTY IN THE CROSS SECTIONS THEMSELVES. AS SUCH THE CONVERSION OF CROSS SECTIONS TO LINEARLY INTERPOLABLE FORM CAN BE PERFORMED WITH ESSENTIALLY NO LOSE OF INFORMATION.

THE ALLOWABLE ERROR MAY BE ENERGY INDEPENDENT (CONSTANT) OR ENERGY Spectra DEPENDENT. THE ALLOWABLE ERROR IS DESCRIBED BY A TABULATED FUNCTION OF UP TO 20 (ENERGY, ERROR) PAIRS AND LINEAR INTERPOLATION Spectra BETWEEN TABULATED POINTS. IF ONLY ONE TABULATED POINT IS GIVEN THE Spectra ERROR WILL BE CONSIDERED CONSTANT OVER THE ENTIRE ENERGY RANGE. WITH THIS ENERGY DEPENDENT ERROR ONE MAY OPTIMIZE THE OUTPUT FOR ANY GIVEN APPLICATION BY USING A SMALL ERROR IN THE ENERGY RANGE OF INTEREST AND A LESS STRINGENT ERROR IN OTHER ENERGY RANGES.

DEFAULT ALLOWABLE ERROR

IN ORDER TO INSURE CONVERGENCE OF THE LINEARIZING ALGORITHM THE ALLOWABLE ERROR MUST BE POSITIVE. IF THE USER INPUTS AN ERROR THAT IS NOT POSITIVE IT WILL AUTOMATICALLY BE SET TO THE DEFAULT VALUE (CURRENTLY 0.001, CORRESPONDING TO 0.1 PER-CENT) AND INDICATED AS SUCH IN THE OUTPUT LISTING.

COULOMB PENETRABILITY (INTERPOLATION LAW = 6) _____ INTRODUCED FOR ENDF/B-VI. THIS IS DEFINED AS,

SIG(E) = C1*EXP(-C2/SQRT(E - T))

THIS PROGRAM ONLY CONSIDERS EXOTHERMIC REACTIONS - T = 0

SIG(E) = C1*EXP(-C2/SQRT(E))

WARNING...THIS INTERPOLATION LAW SHOULD ONLY BE USED FOR REACTIONS Spectra WHICH HAVE A POSITIVE Q-VALUE (EXOTHERMIC REACTIONS), SINCE HERE WE ONLY CONSIDER T = 0.0 IN THE FORMALISM. IN ALL OTHER CASES A WARNING MESSAGE WILL BE PRINTED.

INPUT FILES

UNIT DESCRIPTION

- 2 INPUT LINES (BCD 80 CHARACTERS/RECORD)
- 10 ORIGINAL ENDF/B DATA (BCD 80 CHARACTERS/RECORD)

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OUTPUT FILES Spectra _____ Spectra UNIT DESCRIPTION Spectra Spectra 3 OUTPUT REPORT (BCD - 120 CHARACTERS/RECORD) Spectra 11 FINAL ENDF/B DATA (BCD - 80 CHARACTERS/RECORD) Spectra Spectra SCRATCH FILES Spectra Spectra UNIT DESCRIPTION Spectra -----Spectra 12 SCRATCH FILE (BINARY - 180000 WORDS/RECORD Spectra Spectra OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINE FILEIO) Spectra Spectra UNIT FILE NAME Spectra ---- ------Spectra 2 SPECTRA. INP Spectra 3 SPECTRA.LST Spectra 10 ENDFB.IN Spectra ENDFB.OUT 11 Spectra (SCRATCH) Spectra Spectra Spectra INPUT PARAMETERS Spectra

FOR VERSIONS EARLIER THAN 90-1 THIS PROGRAM ONLY ALLOWED THE USER TO SPECIFY BY INPUT PARAMETERS WHICH MATERIALS (MAT) TO PROCESS. FOR EACH REQUESTED MATERIAL NEUTRON INTERACTION CROSS SECTIONS (MF=3) WOULD BE LINEARIZED AND THE REMAINDER OF THE MATERIAL WOULD BE COPIED.

FOR VERSIONS 90-1 AND LATER THIS PROGRAM WILL ALLOW THE USER TO TO SPECIFY BY INPUT PARAMETERS EXACTLY WHAT SECTIONS OF DATA TO PROCESS. FOR EACH SECTION OF DATA, SPECIFIED BY MAT, MF, MT RANGES, SECTIONS OF MF=3, 23 AND 27 WILL BE LINEARIZED AND ALL OTHER REQUESTED SECTIONS WILL BE COPIED. ALL SECTIONS WHICH ARE NOT EXPLICITLY REQUESTED WILL BE SKIPPED AND WILL NOT APPEAR ON ENDF/B FILE OUTPUT BY THIS PROGRAM.

WITH THIS NEW PROCEDURE YOU CAN MINIMIZE THE SIZE OF THE ENDF/B FILE OUTPUT BY THIS PROGRAM, E.G., IF YOU ONLY WANT NEUTRON CROSS SECTIONS FOR SUBSEQUENT PROCESSING YOU NEED ONLY REQUEST ONLY MF=3 DATA.

HOWEVER, YOU MUST UNDERSTAND THAT ONLY THOSE SECTIONS WHICH YOU EXPLICITLY REQUEST WILL APPEAR ON THE ENDF/B FILE OUTPUT BY THIS PROGRAM. FOR EXAMPLE, IF YOU WISH TO DOCUMENT EXACTLY HOW YOU LINEARIZED THE DATA BY INCLUDING COMMENTS IN MF=1, MT=451 THEN YOU MUST EXPLICITLY REQUEST THAT MF=1, MT=451 BE PROCESSED FOR EACH MATERIAL THAT YOU REQUEST. SIMILAR IF YOU WANT THE ENTIRE EVALUATION YOU MUST REQUEST ALL MF AND MT TO BE OUTPUT.

LINE COLS. DESCRIPTION

- 1-11 SELECTION CRITERIA (0=MAT, 1=ZA)
 - 12-22 MONITOR MODE SELECTOR
 - = 0 NORMAL OPERATION
 - = 1 MONITOR PROGRESS OF LINEARIZING OF THE DATA. EACH TIME A PAGE OF DATA POINTS IS WRITTEN TO THE SCRATCH FILE PRINT OUT THE TOTAL NUMBER OF Spectra POINTS ON SCRATCH AND THE LOWER AND UPPER ENERGY LIMITS OF THE PAGE (THIS OPTION MAY BE

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USED IN ORDER TO MONITOR THE EXECUTION SPEED Spectra OF LONG RUNNING JOBS). Spectra 23-33 MINIMUM CROSS SECTION OF INTEREST (BARNS). Spectra (IF 0.0 OR LESS IS INPUT THE PROGRAM WILL Spectra USE 1.0E-10). ENERGY INTERVALS WILL NOT BE Spectra SUB-DIVIDED IF THE ABSOLUTE VALUE OF THE CROSS Spectra SECTION WITHIN THE INTERVAL IS LESS THAN THIS VALUE. Spectra AN EXCEPTION TO THIS RULE IS NEAR THRESHOLDS ENERGY Spectra INTERVALS WILL BE SUB-DIVIDED UNTIL CONVERGENCE Spectra REGARDLESS OF THE MAGNITUDE OF THE CROSS SECTION. Spectra 34-44 KEEP ORIGINAL EVALUATED DATA POINTS. Spectra = 0 - NO.Spectra = 1 - YES - ADDITIONAL POINTS MAY BE ADDED IN ORDER Spectra TO LINEARIZE DATA, BUT ALL ORIGINAL Spectra DATA POINTS WILL BE INCLUDED IN THE Spectra RESULTS. Spectra 1-72 ENDF/B INPUT DATA FILENAME 2 Spectra (STANDARD OPTION = ENDFB.IN) Spectra 3 1-72 ENDF/B OUTPUT DATA FILENAME Spectra (STANDARD OPTION = ENDFB.OUT) Spectra 4-N 1- 6 LOWER MAT OR ZA LIMIT Spectra 7-8 LOWER MF LIMIT Spectra 9-11 LOWER MT LIMIT Spectra 12-17 UPPER MAT OR ZA LIMIT Spectra 18-19 UPPER MF LIMIT Spectra 20-22 UPPER MT LIMIT Spectra UP TO 100 RANGES MAY BE SPECIFIED, ONLY ONE RANGE Spectra PER LINE. THE LIST OF RANGES IS TERMINATED BY A Spectra BLANK LINE. IF THE UPPER MAT LIMIT OF ANY REQUEST Spectra IS LESS THAN THE LOW LIMIT IT WILL BE SET EQUAL TO Spectra THE LOWER LIMIT. IF THE UPPER LIMIT IS STILL ZERO Spectra IT WILL BE SET EQUAL TO 999999. IF THE UPPER MF OR Spectra MT LIMIT IS ZERO IT WILL BE SET TO 99 OR 999 Spectra RESPECTIVELY. Spectra VARY 1-11 ENERGY FOR ERROR LAW Spectra ALLOWABLE FRACTIONAL ERROR FOR ERROR LAW. Spectra THE ACCEPTABLE LINEARIZING ERROR MAY BE SPECIFIED TO Spectra BE EITHER ENERGY INDEPENDENT (DEFINED BY A SINGLE Spectra ERROR), OR ENERGY DEPENDENT (DEFINED BY UP TO 20 Spectra ENERGY, ERROR PAIRS). FOR THE ENERGY DEPENDENT CASE Spectra LINEAR INTERPOLATION WILL BE USED TO DEFINE THE ERROR Spectra AT ENERGIES BETWEEN THOSE AT WHICH IT IS TABULATED. Spectra IN ALL CASES THE ERROR LAW IS TERMINATED BY A BLANK Spectra LINE. IF ONLY ONE ENERGY, ERROR PAIR IS GIVEN THE Spectra THE LAW WILL BE CONSIDERED TO BE ENERGY INDEPENDENT. Spectra IF MORE THAN ONE PAIR IS GIVEN IT WILL BE CONSIDERED Spectra TO BE ENERGY DEPENDENT (NOTE, ENERGY INDEPENDENT Spectra FORM WILL RUN FASTER THAN THE EQUIVALENT ENERGY Spectra DEPENDENT FORM). FOR AN ENERGY DEPENDENT ERROR LAW Spectra ALL ENERGIES MUST BE ASCENDING ENERGY ORDER. FOR Spectra CONVERGENCE OF THE LINEARIZING ALGORITHM ALL ERRORS Spectra MUST BE POSITIVE. IF AN ALLOWABLE ERROR IS NOT Spectra POSITIVE IT WILL BE SET EQUAL TO THE STANDARD OPTION Spectra (CURRENTLY 0.001, CORRESPONDING TO 0.1 PER-CENT). Spectra IF THE FIRST ERROR LINE IS BLANK IT WILL TERMINATE Spectra THE ERROR LAW AND THE ERROR WILL BE TREATED AS Spectra ENERGY INDEPENDENT, EQUAL TO THE STANDARD OPTION Spectra (CURRENTLY 0.1 PER-CENT). (SEE EXAMPLE INPUT 4). Spectra Spectra EXAMPLE INPUT NO. 1 Spectra Spectra RETRIEVE DATA BY ZA IN ORDER TO FIND ALL URANIUM ISOTOPES AND Spectra THORIUM 232. RETRIEVE ALL NEUTRON INTERACTION CROSS SECTIONS Spectra

(MF=3). ALL ENERGY INTERVALS IN WHICH THE CROSS SECTION IS	Spectra
AT LEAST 1 MICRO-BARN (1.0E-06 BARNS) WILL BE SUBDIVIDED.	Spectra
BACKWARD THINNING WILL BE PERFORMED. FROM 0 TO 100 EV LINEARIZE	Spectra
TO WITHIN 0.1 PER-CENT ACCURACY. FROM 100 EV TO 1 KEV VARY	Spectra
ACCURACY BETWEEN 0.1 AND 1.0 PER-CENT. ABOVE 1 KEV USE 1	Spectra
	_
PER-CENT ACCURACY.	Spectra
THE TAXABLE OF THE OFFICE OF T	Spectra
EXPLICITLY SPECIFY THE STANDARD FILENAMES.	Spectra
	Spectra
IN THIS CASE THE FOLLOWING 11 INPUT LINES ARE REQUIRED	Spectra
	Spectra
1 0 1.00000- 6 0	Spectra
ENDFB.IN	Spectra
ENDFB.OUT	Spectra
92000 3 0 92999 3999	Spectra
90232 3 0 0 3 0 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999) Spectra
(END OF REQUEST LIST)	Spectra
0.00000+ 0 1.00000-03	Spectra
1.00000+ 2 1.00000-03	Spectra
1.00000+ 3 1.00000-02	Spectra
1.00000+ 9 1.00000-02	Spectra
(END OF ERROR LAW)	Spectra
·	Spectra
EXAMPLE INPUT NO. 2	Spectra
	Spectra
SAME AS THE ABOVE CASE, EXCEPT LINEARIZE ALL DATA TO WITHIN THE	Spectra
STANDARD ACCURACY (CURRENTLY 0.1 PER-CENT). IN ORDER TO USE THE	Spectra
STANDARD ACCURACY YOU NEED NOT SPECIFY ANY ERROR LAW AT ALL. IN	Spectra
THIS CASE INCLUDE THE HOLLERITH SECTION, MF=1, MT=451, FOR EACH	Spectra
MATERIAL.	Spectra
MATERIAL.	-
THE DESIGNATION OF THE STIENNING DIAMS. THE DOCUMENTS	Spectra
LEAVE THE DEFINITION OF THE FILENAMES BLANK - THE PROGRAM WILL	Spectra
THEN USE STANDARD FILENAMES.	Spectra
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THE WHITE CASE WHE COLLECTIVE A THRUM TIMES ARE RECUIRED	Spectra
IN THIS CASE THE FOLLOWING 9 INPUT LINES ARE REQUIRED	Spectra
	Spectra Spectra
1 0 1.00000- 6 0	Spectra Spectra Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN)	Spectra Spectra Spectra Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT)	Spectra Spectra Spectra Spectra Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451	Spectra Spectra Spectra Spectra Spectra Spectra Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451 92000 3 0 92999 3999	Spectra Spectra Spectra Spectra Spectra Spectra Spectra Spectra
1 0 1.00000-6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451 92000 3 0 92999 3999 90232 1451 0 1451	Spectra Spectra Spectra Spectra Spectra Spectra Spectra Spectra Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451 92000 3 0 92999 3999 90232 1451 0 1451 90232 3 0 0 3 0 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999	Spectra Spectra Spectra Spectra Spectra Spectra Spectra Spectra Spectra
1 0 1.00000-6 0	Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451 92000 3 0 92999 3999 90232 1451 0 1451 90232 3 0 0 3 0 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999	Spectra
1 0 1.00000-6 0	Spectra
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1 0 1.00000-6 0	Spectra
1 0 1.00000-6 0	Spectra
1 0 1.00000- 6 0	Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451 92000 3 0 92999 3999 90232 1451 0 1451 90232 3 0 0 3 0 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999 (END OF REQUEST LIST) (0.1 PER-CENT ERROR, END OF ERROR LAW) EXAMPLE INPUT NO. 3 LINEARIZE ALL MATERIALS ON AN ENDF/B TAPE TO WITHIN AN ACCURACY OF 0.5 PER-CENT (0.005 AS A FRACTION). IN THIS CASE YOU NEED NOT SPECIFY THE MAT, MF, MT RANGES. READ THE ENDF/B DATA FROM \ENDFB6\ZA092238 AND WRITE THE ENDF/B DATA TO \ENDFB6\LINEAR\ZA092238. IN THIS CASE THE FOLLOWING 6 INPUT LINES ARE REQUIRED	Spectra
1 0 1.00000- 6 0	Spectra
1 0 1.00000- 6 0	Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451 92000 3 0 92999 3999 90232 1451 0 1451 90232 3 0 0 3 0 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999 (END OF REQUEST LIST) (0.1 PER-CENT ERROR, END OF ERROR LAW) EXAMPLE INPUT NO. 3	Spectra
1 0 1.00000- 6 0	Spectra
1 0 1.00000- 6 0 (USE DEFAULT FILENAME = ENDFB.IN) (USE DEFAULT FILENAME = ENDFB.OUT) 92000 1451 92999 1451 92000 3 0 92999 3999 90232 1451 0 1451 90232 3 0 0 3 0 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999 (END OF REQUEST LIST) (0.1 PER-CENT ERROR, END OF ERROR LAW) EXAMPLE INPUT NO. 3	Spectra

Spectra NOTE THAT IN THIS CASE IF THE INPUT HAD SPECIFIED AN EQUIVALENT Spectra ENERGY DEPENDENT ERROR LAW BY GIVING A NUMBER OF ENERGY POINTS Spectra AT EACH OF WHICH THE ERROR IS 0.5 PER-CENT THE PROGRAM WOULD TAKE Spectra LONGER TO RUN (I.E., ONLY USE AN ENERGY DEPENDENT ERROR LAW WHEN Spectra IT IS NECESSARY). Spectra Spectra EXAMPLE INPUT NO. 4 Spectra Spectra IN ORDER TO LINEARIZE ALL MATERIALS ON AN ENDF/B TAPE TO THE Spectra STANDARD OPTION OF 0.1 PER-CENT IT IS ADEQUATE TO INPUT A SET Spectra OF COMPLETELY BLANK LINES WHICH WILL AUTOMATICALLY INVOKE ALL Spectra OF THE STANDARD OPTIONS. Spectra Spectra LEAVE THE DEFINITION OF THE FILENAMES BLANK - THE PROGRAM WILL Spectra THEN USE STANDARD FILENAMES. Spectra Spectra IN THIS CASE THE FOLLOWING THREE INPUT LINES ARE REQUIRED Spectra Spectra (MAT, 1.0E-10 BARNS, THIN) Spectra (USE DEFAULT FILENAME = ENDFB.IN) Spectra (USE DEFAULT FILENAME = ENDFB.OUT) Spectra (RETRIEVE ALL DATA, END REQUEST LIST) Spectra (0.1 PER-CENT ERROR, END OF ERROR LAW) Spectra

Spectra

====== Spectra