PROGRAM SPECTRA An extension of LINEAR to linearize ALl MF=5 spectra. 2012/05/28 - Added MF=15 neutron induced, photon spectra. 2019/01/03 - MF=6 Still NOT Implemented First released in 2010 - Earlier below dates refer to LINEAR. VERSION 74-1 (MAY 1974) VERSION 75-1 (APRIL 1975) VERSION 76-2 (OCTOBER 1976) VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978) VERSION 79-1 (JULY 1979) CDC-7600 AND CRAY-1 VERSION.	Spectra Spectra Spectra Spectra Spectra Spectra
An extension of LINEAR to linearize ALl MF=5 spectra. 2012/05/28 - Added MF=15 neutron induced, photon spectra. 2019/01/03 - MF=6 Still NOT Implemented First released in 2010 - Earlier below dates refer to LINEAR. VERSION 74-1 (MAY 1974) VERSION 75-1 (APRIL 1975) VERSION 76-2 (OCTOBER 1976) VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978)	Spectra Spectra Spectra Spectra Spectra
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VERSION 74-1 (MAY 1974) VERSION 75-1 (APRIL 1975) VERSION 76-2 (OCTOBER 1976) VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978)	Spectra
VERSION 74-1 (MAY 1974) VERSION 75-1 (APRIL 1975) VERSION 76-2 (OCTOBER 1976) VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978)	-
VERSION 75-1 (APRIL 1975) VERSION 76-2 (OCTOBER 1976) VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978)	
VERSION 75-1 (APRIL 1975) VERSION 76-2 (OCTOBER 1976) VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978)	Spectra
VERSION 76-2 (OCTOBER 1976) VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978)	Spectra
VERSION 77-1 (JANUARY 1977) VERSION 78-1 (JULY 1978)	Spectra
	Spectra
(27)	Spectra
VERSION /9-1 (BOLI 1979) CDC-7000 AND CRAI-I VERSION.	Spectra
VERSION 80-1 (MAY 1980) IBM, CDC AND CRAY VERSION.	Spectra
VERSION 80-2 (DECEMBER 1980)	Spectra
VERSION 81-1 (MARCH 1981)	Spectra
VERSION 82-1 (JANUARY 1982) IMPROVED COMPUTER COMPATIBILITY.	Spectra
VERSION 83-1 (JANUARY 1983) *MAJOR RE-DESIGN.	Spectra
*PAGE SIZE INCREASED - 1002 TO 3006.	Spectra
*ELIMINATED COMPUTER DEPENDENT CODING.	-
*NEW, MORE COMPATIBLE I/O UNIT NUMBER. *ADDED OPTION TO KEEP ALL ORIGINAL	-
ENERGY POINTS FROM EVALUATION.	Spectra
*ADDED STANDARD ALLOWABLE ERROR OPTION	
(CURRENTLY 0.1 PER-CENT).	Spectra
VERSION 83-2 (OCTOBER 1983) IMPROVED BASED ON USER COMMENTS.	Spectra
VERSION 84-1 (APRIL 1984) IMPROVED BASED ON USER COMMENTS.	Spectra
VERSION 84-2 (JUNE 1984) *UPDATED FOR ENDF/B-VI FORMATS.	Spectra
*SPECIAL I/O ROUTINES TO GUARANTEE	Spectra
ACCURACY OF ENERGY.	Spectra
*DOUBLE PRECISION TREATMENT OF ENERGY	Spectra
(REQUIRED FOR NARROW RESONANCES).	Spectra
VERSION 85-1 (AUGUST 1985) *FORTRAN-77/H VERSION	Spectra
VERSION 86-1 (JANUARY 1986) *ENDF/B-VI FORMAT	Spectra
VERSION 87-1 (JANUARY 1987) *DOUBLE PRECISION TREATMENT OF CROSS	Spectra
SECTION VERSION 88-1 (JULY 1988) *OPTIONINTERNALLY DEFINE ALL I/O	Spectra
FILE NAMES (SEE, SUBROUTINE FILEIO	Spectra
FOR DETAILS).	Spectra
*IMPROVED BASED ON USER COMMENTS.	Spectra
VERSION 89-1 (JANUARY 1989) * PSYCHOANALYZED BY PROGRAM FREUD TO	Spectra
INSURE PROGRAM WILL NOT DO ANYTHING	Spectra
CRAZY.	Spectra
*UPDATED TO USE NEW PROGRAM CONVERT	Spectra
KEYWORDS.	Spectra
*ADDED LIVERMORE CIVIC COMPILER	Spectra
CONVENTIONS.	Spectra
VERSION 90-1 (JUNE 1990) *EXTENDED TO LINEARIZE PHOTON	Spectra
INTERACTION DATA, MF=23 AND 27	Spectra
*ADDED FORTRAN SAVE OPTION	Spectra
*UPDATED BASED ON USER COMMENTS.	Spectra
*NEW MORE CONSISTENT ENERGY OUTPUT ROUTINE.	Spectra
*WARNINGINPUT PARAMETER FORMAT	Spectra
HAS BEEN CHANGEDSEE DESCRIPTION	Spectra
BELOW.	Spectra
VERSION 91-1 (JULY 1991) *ADDED INTERPOLATION LAW 6 - ONLY USED	Spectra
FOR CHARGED PARTICLE CROSS SECTIONS	Spectra
FOR COULOMB PENETRABILITIES.	Spectra
VERSION 92-1 (JANUARY 1992)*ADDED NU-BAR (TOTAL, DELAYED, PROMPT)	Spectra
POLYNOMIAL OR TABULATED ALL CONVERTED	Spectra
TO LINEARLY INTERPOLABLE	Spectra
	Spectra
*INCREASED PAGE SIZE FROM 3006 TO 5010	Spectra
POINTS.	Spectra
POINTS. *ALL ENERGIES INTERNALLY ROUNDED PRIOR	-
POINTS. *ALL ENERGIES INTERNALLY ROUNDED PRIOR TO CALCULATIONS.	Spectra
POINTS. *ALL ENERGIES INTERNALLY ROUNDED PRIOR	Spectra

				DEPENDENCE .	Spectra
VERSI	ON 92-2	(JULY 1		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
THEOT	07 02 1		10021 4	ENERGY DEPENDENT QUANTITY).	Spectra
VERSIO	ON 93-1	(MARCH)	1993) *	UPDATED FOR USE WITH LAHEY COMPILER ON IBM-PCS.	Spectra Spectra
			*	INCREASED PAGE SIZE FROM 5010 TO	Spectra
				30000 POINTS	Spectra
VERSIO	ON 94-1	(JANUAR	Y 1994)*	VARIABLE ENDF/B DATA FILENAMES	Spectra
				TO ALLOW ACCESS TO FILE STRUCTURES	Spectra
				(WARNING - INPUT PARAMETER FORMAT	Spectra
			*	HAS BEEN CHANGED) CLOSE ALL FILES BEFORE TERMINATING	Spectra Spectra
				(SEE, SUBROUTINE ENDIT)	Spectra
VERSI	ON 96-1	(JANUAR	Y 1996)	*COMPLETE RE-WRITE	Spectra
				*IMPROVED COMPUTER INDEPENDENCE	Spectra
				*ALL DOUBLE PRECISION	Spectra
				*ON SCREEN OUTPUT	Spectra
				*UNIFORM TREATMENT OF ENDF/B I/O *IMPROVED OUTPUT PRECISION	Spectra Spectra
				*DEFINED SCRATCH FILE NAMES	Spectra
				*ALWAYS INCLUDE THERMAL VALUE	Spectra
				*INCREASED PAGE SIZE FROM 30000 TO	Spectra
				60000 POINTS	Spectra
VERSI	ON 99-1	(MARCH)	1999)	*CORRECTED CHARACTER TO FLOATING	Spectra
				POINT READ FOR MORE DIGITS *UPDATED TEST FOR ENDF/B FORMAT	Spectra
				VERSION BASED ON RECENT FORMAT CHANGE	Spectra Spectra
				*GENERAL IMPROVEMENTS BASED ON	Spectra
				USER FEEDBACK	Spectra
VERSIO	ON 99-2	(JUNE 19	999)	*ASSUME ENDF/B-VI, NOT V, IF MISSING	Spectra
	0000 1	/====		MF=1, MT-451.	Spectra
VERS.	2000-1	(FEBRUAI	RY 2000)	*ADDED MF = 9 AND 10 LINEARIZATION *GENERAL IMPROVEMENTS BASED ON	Spectra
				USER FEEDBACK	Spectra Spectra
VERS.	2002-1	(MAY 20	02)	*OPTIONAL INPUT PARAMETERS	Spectra
VERS.	2004-1	(JAN. 20	004)	*GENERAL UPDATE BASED ON USER FEEDBACK	Spectra
VERS.	2005-1	(JAN. 20	005)	*ALWAYS KEEP ORIGINAL TABULATED	Spectra
VEDO	2006-1	(EED 0)	0061	NU-BAR POINTS. *CORRECTED INT=6 NEAR THRESHOLD	Spectra
VERS.	2006-1	(FEB. 20	008)	*NO SUBDIVIDE BELOW MINIMUM XCMIN	Spectra Spectra
VERS.	2007-1	(JAN. 20	007)	*CHECKED AGAINST ALL ENDF/B-VII.	Spectra
		•		*INCREASED PAGE SIZE FROM 60,000 TO	Spectra
				600,000 POINTS	Spectra
VERS.	2010-1	(JUNE 2)	010)	*ADDED MF = 5 - MF = 6 STILL PLANNED.	Spectra
				*72 CHARACTER FILE NAMES. *ONLY PROCESS MF=5 - SKIP ALL OTHERS	Spectra
				TO PREVENT CONFLICT WITH LINEAR	Spectra Spectra
				THINNING.	Spectra
VERS.	2012-1	(Aug. 20	012)	*Added MF=15, neutron induced photon	Spectra
				spectra.	Spectra
				*Added CODENAME	Spectra
				*32 and 64 bit Compatible *Added ERROR stop	Spectra Spectra
VERS.	2015-1	(Jan. 20	015)	*Extended OUT9.	Spectra
		,	,	*Replaced ALL 3 way IF Statements.	Spectra
				*Corrected MF=15 Data - it was adding	Spectra
				SEND between sub-sections.	Spectra
VEDO	2017 1	(Mar- 0)	017)	*Deleted unused parts, e.g., NUBAR.	Spectra
VERS.	2017-1	(may 20	017)	*Increased page size to 3,000,000 *Updated based on user feedback	Spectra Spectra
				*Changed DGAMMA to REDGAMMA to avoid	Spectra
				conflict with possble RESERVED NAME	Spectra
				*All floating input parameters changed	
	0010 -	(-	01.01	to character input + IN9 conversion.	Spectra
		(Jan. 20		*On-line output for ALL ENDERROR	Spectra
VERO.	2019-1	(June 20	019)	*Additional Interpolation Law Tests *Check Maximum Tabulated Energy to	Spectra Spectra
				insure it is the same for all MTs -	Spectra

if not, print WARNING messages. Spectra *Corrected END Histogram linearized -Spectra Previously deleted last point - ERROR Spectra to assume this has Y=0 - now keep Spectra point, but insure Y = 0. 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 Spectra PRESENT CONTACT INFORMATION Spectra Spectra Dermott E. Cullen Spectra 1466 Hudson Way Spectra Livermore, CA 94550 Spectra U.S.A. Spectra Telephone 925-443-1911 Spectra E. Mail RedCullen1@Comcast.net Spectra Website RedCullen1.net/HOMEPAGE.NEW Spectra 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 THE LATEST DOCUMENTATION INCLUDING ALL RECENT IMPROVEMENTS. PLEASE Spectra READ ALL OF THESE COMMENTS BEFORE IMPLEMENTATION. Spectra Spectra AT THE PRESENT TIME WE ARE ATTEMPTING TO DEVELOP A SET OF COMPUTER Spectra INDEPENDENT PROGRAMS THAT CAN EASILY BE IMPLEMENTED ON ANY ONE Spectra OF A WIDE VARIETY OF COMPUTERS. IN ORDER TO ASSIST IN THIS PROJECT Spectra IT WOULD BE APPECIATED IF YOU WOULD NOTIFY THE AUTHOR OF ANY Spectra COMPILER DIAGNOSTICS, OPERATING PROBLEMS OR SUGGESTIONS ON HOW TO Spectra IMPROVE THIS PROGRAM. HOPEFULLY, IN THIS WAY FUTURE VERSIONS OF Spectra THIS PROGRAM WILL BE COMPLETELY COMPATIBLE FOR USE ON YOUR Spectra COMPUTER. Spectra Spectra PURPOSE Spectra Spectra THIS PROGRAM IS DESIGNED TO CONVERT ENDF/B FILE 3, 23 AND 27 DATA Spectra TO LINEAR-LINEAR INTERPOLABLE FORM. ANY SECTION THAT IS ALREADY Spectra LINEAR-LINEAR INTERPOLABLE WILL BE THINNED. Spectra Spectra IN THE FOLLOWING DISCUSSION FOR SIMPLICITY THE ENDF/B TERMINOLOGY Spectra ---ENDF/B TAPE---WILL BE USED. IN FACT THE ACTUAL MEDIUM MAY BE Spectra TAPE, CARDS, DISK OR ANY OTHER MEDIUM. Spectra Spectra ENDF/B FORMAT Spectra _____ Spectra THIS PROGRAM ONLY USES THE ENDF/B BCD OR CARD IMAGE FORMAT (AS Spectra OPPOSED TO THE BINARY FORMAT) AND CAN HANDLE DATA IN ANY VERSION Spectra OF THE ENDF/B FORMAT (I.E., ENDF/B-I, II, III, IV, V OR VI FORMAT). Spectra Spectra IT IS ASSUMED THAT THE DATA IS CORRECTLY CODED IN THE ENDF/B Spectra FORMAT AND NO ERROR CHECKING IS PERFORMED. IN PARTICULAR IT IS Spectra ASSUMED THAT THE MAT, MF AND MT ON EACH LINE IS CORRECT. SEQUENCE Spectra NUMBERS (COLUMNS 76-80) ARE IGNORED ON INPUT, BUT WILL BE Spectra CORRECTLY OUTPUT ON ALL LINES. THE FORMAT OF SECTION MF=1, MT=451 Spectra AND ALL SECTIONS OF MF=3 MUST BE CORRECT. THE PROGRAM COPIES ALL Spectra OTHER SECTION OF DATA AS HOLLERITH AND AS SUCH IS INSENSITIVE TO Spectra THE CORRECTNESS OR INCORRECTNESS OF ALL OTHER SECTIONS. Spectra Spectra OUTPUT FORMAT Spectra Spectra -----

IN THIS VERSION OF LINEAR ALL ENERGIES WILL BE OUTPUT IN Spectra F (INSTEAD OF E) FORMAT IN ORDER TO ALLOW ENERGIES TO BE WRITTEN Spectra WITH UP TO 9 DIGITS OF ACCURACY. IN PREVIOUS VERSIONS THIS WAS AN Spectra OUTPUT OPTION. HOWEVER USE OF THIS OPTION TO COMPARE THE RESULTS Spectra OF ENERGIES WRITTEN IN THE NORMAL ENDF/B CONVENTION OF 6 DIGITS Spectra TO THE 9 DIGIT OUTPUT FROM THIS PROGRAM DEMONSTRATED THAT FAILURE Spectra TO USE THE 9 DIGIT OUTPUT CAN LEAD TO LARGE ERRORS IN THE DATA Spectra DUE TO TRUNCATION OF ENERGIES TO 6 DIGITS DURING OUTPUT. Spectra Spectra CONTENTS OF OUTPUT Spectra _____ Spectra ENTIRE EVALUATIONS ARE OUTPUT, NOT JUST THE LINEARIZED DATA Spectra CROSS SECTIONS, E.G. ANGULAR AND ENERGY DISTRIBUTIONS ARE ALSO Spectra INCLUDED. Spectra Spectra DOCUMENTATION Spectra Spectra THE FACT THAT THIS PROGRAM HAS OPERATED ON THE DATA IS DOCUMENTED Spectra BY THE ADDITION OF 3 COMMENT LINES AT THE END OF EACH HOLLERITH Spectra SECTION IN THE FORM Spectra Spectra Spectra FOR ALL DATA GREATER THAN 1.00000-10 IN ABSOLUTE VALUE Spectra DATA LINEARIZED TO WITHIN AN ACCURACY OF 0.1 PER-CENT Spectra Spectra THE ORDER OF SIMILAR COMMENTS (FROM RECENT, SIGMA1 AND GROUPIE) Spectra REPRESENTS A COMPLETE HISTORY OF ALL OPERATIONS PERFORMED ON Spectra THE DATA BY THESE PROGRAMS. Spectra Spectra 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 Spectra EARLIER VERSIONS OF ENDF/B. BY READING AN EXISTING MF=1, MT=451 Spectra IT IS POSSIBLE FOR THIS PROGRAM TO DETERMINE WHICH VERSION OF Spectra Spectra 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 Spectra DETERMINE WHICH VERSION OF THE ENDF/B FORMAT THE DATA IS IN, AND Spectra AS SUCH IT IS IMPOSSIBLE FOR THE PROGRAM TO DETERMINE WHAT FORMAT Spectra SHOULD BE USED TO CREATE A HOLLERITH SECTION. Spectra Spectra REACTION INDEX Spectra Spectra THIS PROGRAM DOES NOT USE THE REACTION INDEX WHICH IS GIVEN IN Spectra SECTION MF=1, MT=451 OF EACH EVALUATION. Spectra Spectra THIS PROGRAM DOES NOT UPDATE THE REACTION INDEX IN MF=1, MT=451. Spectra THIS CONVENTION HAS BEEN ADOPTED BECAUSE MOST USERS DO NOT Spectra REQUIRE A CORRECT REACTION INDEX FOR THEIR APPLICATIONS AND IT WAS Spectra NOT CONSIDERED WORTHWHILE TO INCLUDE THE OVERHEAD OF CONSTRUCTING Spectra 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 Spectra SECTION SIZE Spectra Spectra SINCE THIS PROGRAM USES A LOGICAL PAGING SYSTEM THERE IS NO LIMIT Spectra TO THE NUMBER OF POINTS IN ANY SECTION, E.G., THE TOTAL CROSS Spectra SECTION MAY BE REPRESENTED BY 200,000 DATA POINTS. Spectra Spectra 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 Spectra 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 Spectra SELECTION OF DATA Spectra Spectra THE PROGRAM SELECTS DATA TO BE LINEARIZED BASED EITHER ON EITHER Spectra MAT (ENDF/B MAT NO.) OR ZA AS WELL AS MF AND MT NUMBERS. THIS Spectra

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. Spectra

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

KEEP EVALUATED DATA POINTS

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

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 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 Spectra IMPORTANTLY CAN BE PERFORMED TO A TOLERANCE THAT IS SMALL COMPARED Spectra TO THE UNCERTAINTY IN THE CROSS SECTIONS THEMSELVES. AS SUCH THE Spectra CONVERSION OF CROSS SECTIONS TO LINEARLY INTERPOLABLE FORM CAN BE Spectra PERFORMED WITH ESSENTIALLY NO LOSE OF INFORMATION. Spectra

THE ALLOWABLE ERROR MAY BE ENERGY INDEPENDENT (CONSTANT) OR ENERGY Spectra DEPENDENT. THE ALLOWABLE ERROR IS DESCRIBED BY A TABULATED Spectra 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. Spectra WITH THIS ENERGY DEPENDENT ERROR ONE MAY OPTIMIZE THE OUTPUT FOR Spectra

ANY GIVEN APPLICATION BY USING A SMALL ERROR IN THE ENERGY RANGE	Spectra
OF INTEREST AND A LESS STRINGENT ERROR IN OTHER ENERGY RANGES.	Spectra Spectra
DEFAULT ALLOWABLE ERROR	Spectra
	Spectra
IN ORDER TO INSURE CONVERGENCE OF THE LINEARIZING ALGORITHM THE	Spectra
ALLOWABLE ERROR MUST BE POSITIVE. IF THE USER INPUTS AN ERROR THAT IS NOT POSITIVE IT WILL AUTOMATICALLY BE SET TO THE DEFAULT	Spectra Spectra
VALUE (CURRENTLY 0.001, CORRESPONDING TO 0.1 PER-CENT) AND	Spectra
INDICATED AS SUCH IN THE OUTPUT LISTING.	Spectra
	Spectra
COULOMB PENETRABILITY (INTERPOLATION LAW = 6)	Spectra
	Spectra
INTRODUCED FOR ENDF/B-VI. THIS IS DEFINED AS,	Spectra
$SIG(E) = C1 \times EXP(-C2/SQRT(E - T))$	Spectra Spectra
$Sig(E) = Ci^{*}Ext(Cz) SQRI(E - 1))$	Spectra
THIS PROGRAM ONLY CONSIDERS EXOTHERMIC REACTIONS - $T = 0$	Spectra
	Spectra
$SIG(E) = C1 \times EXP(-C2/SQRT(E))$	Spectra
	Spectra
WARNINGTHIS INTERPOLATION LAW SHOULD ONLY BE USED FOR REACTIONS	-
WHICH HAVE A POSITIVE Q-VALUE (EXOTHERMIC REACTIONS),	Spectra
SINCE HERE WE ONLY CONSIDER $T = 0.0$ in the Formalism.	Spectra
IN ALL OTHER CASES A WARNING MESSAGE WILL BE PRINTED.	Spectra
INPUT FILES	Spectra Spectra
	Spectra
UNIT DESCRIPTION	Spectra
	Spectra
2 INPUT LINES (BCD - 80 CHARACTERS/RECORD)	Spectra
10 ORIGINAL ENDF/B DATA (BCD - 80 CHARACTERS/RECORD)	Spectra
	Spectra
OUTPUT FILES	Spectra
	Spectra
UNIT DESCRIPTION	Spectra
3 OUTPUT REPORT (BCD - 120 CHARACTERS/RECORD)	Spectra 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
ODETANNE CENTRADD ETTE NAMES (SEE SUDDOUETNE ETTETO)	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
11 ENDFB.OUT	Spectra
12 (SCRATCH)	Spectra
	Spectra
INPUT PARAMETERS	Spectra Spectra
	Spectra
FOR VERSIONS EARLIER THAN 90-1 THIS PROGRAM ONLY ALLOWED THE USER	Spectra
TO SPECIFY BY INPUT PARAMETERS WHICH MATERIALS (MAT) TO PROCESS.	Spectra
FOR EACH REQUESTED MATERIAL NEUTRON INTERACTION CROSS SECTIONS	Spectra
(MF=3) WOULD BE LINEARIZED AND THE REMAINDER OF THE MATERIAL	Spectra
WOULD BE COPIED.	Spectra
	Spectra
FOR VERSIONS 90-1 AND LATER THIS PROGRAM WILL ALLOW THE USER TO	Spectra
TO SPECIFY BY INPUT PARAMETERS EXACTLY WHAT SECTIONS OF DATA TO PROCESS. FOR EACH SECTION OF DATA, SPECIFIED BY MAT, MF, MT	Spectra
RANGES, SECTIONS OF MF=3, 23 AND 27 WILL BE LINEARIZED AND ALL	Spectra Spectra
OTHER REQUESTED SECTIONS WILL BE COPIED. ALL SECTIONS WHICH ARE	Spectra
NOT EXPLICITLY REQUESTED WILL BE SKIPPED AND WILL NOT APPEAR ON	Spectra
-	-

ENDF/1	B FILE (OUTPUT BY THIS PROGRAM.	Spectra
WITH :	THIS NEW	W PROCEDURE YOU CAN MINIMIZE THE SIZE OF THE ENDF/B	Spectra Spectra
		BY THIS PROGRAM, E.G., IF YOU ONLY WANT NEUTRON	Spectra
CROSS	SECTION	NS FOR SUBSEQUENT PROCESSING YOU NEED ONLY REQUEST	Spectra
ONLY I	MF=3 DAT	TA.	Spectra
HOMEN		MUST UNDERSTAND THAT ONLY THOSE SECTIONS WHICH YOU	Spectra
		EQUEST WILL APPEAR ON THE ENDF/B FILE OUTPUT BY	Spectra Spectra
		. FOR EXAMPLE, IF YOU WISH TO DOCUMENT EXACTLY	Spectra
		ARIZED THE DATA BY INCLUDING COMMENTS IN MF=1, MT=451	Spectra
		T EXPLICITLY REQUEST THAT MF=1, MT=451 BE PROCESSED	Spectra
		ERIAL THAT YOU REQUEST. SIMILAR IF YOU WANT THE	Spectra
ENTIR	E EVALUA	ATION YOU MUST REQUEST ALL MF AND MT TO BE OUTPUT.	Spectra Spectra
LINE	COLS.	DESCRIPTION	Spectra
			Spectra
1	1-11	SELECTION CRITERIA (0=MAT, 1=ZA)	Spectra
	12-22	MONITOR MODE SELECTOR	Spectra
		= 0 - NORMAL OPERATION	Spectra
		= 1 - MONITOR PROGRESS OF LINEARIZING OF THE DATA. EACH TIME A PAGE OF DATA POINTS IS WRITTEN TO	Spectra Spectra
		THE SCRATCH FILE PRINT OUT THE TOTAL NUMBER OF	Spectra
		POINTS ON SCRATCH AND THE LOWER AND UPPER	Spectra
		ENERGY LIMITS OF THE PAGE (THIS OPTION MAY BE	Spectra
		USED IN ORDER TO MONITOR THE EXECUTION SPEED	Spectra
	~~ ~~	OF LONG RUNNING JOBS).	Spectra
	23-33	MINIMUM CROSS SECTION OF INTEREST (BARNS). (IF 0.0 OR LESS IS INPUT THE PROGRAM WILL	Spectra 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
	34-44	REGARDLESS OF THE MAGNITUDE OF THE CROSS SECTION. KEEP ORIGINAL EVALUATED DATA POINTS.	Spectra Spectra
	51 11	= 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
2	1-72	RESULTS. ENDF/B INPUT DATA FILENAME	Spectra Spectra
2	1 12	(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 9-11	LOWER MF LIMIT LOWER MT LIMIT	Spectra
	9-11 12-17	UPPER MAT OR ZA LIMIT	Spectra Spectra
		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 IS LESS THAN THE LOW LIMIT IT WILL BE SET EQUAL TO	Spectra Spectra
		THE LOWER LIMIT. IF THE UPPER LIMIT IS STILL ZERO	Spectra
		IT WILL BE SET EQUAL TO 9999999. IF THE UPPER MF OR	Spectra
		MT LIMIT IS ZERO IT WILL BE SET TO 99 OR 999	Spectra
177.017	1-11	RESPECTIVELY. ENERGY FOR ERROR LAW	Spectra
VARI	12-22	ENERGY FOR ERROR LAW ALLOWABLE FRACTIONAL ERROR FOR ERROR LAW.	Spectra 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 AT ENERGIES BETWEEN THOSE AT WHICH IT IS TABULATED.	-
		AT ENERGIES BETWEEN THOSE AT WHICH IT IS TABULATED. IN ALL CASES THE ERROR LAW IS TERMINATED BY A BLANK	Spectra 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 EOUIVALENT 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 Spectra EXPLICITLY SPECIFY THE STANDARD FILENAMES. Spectra Spectra IN THIS CASE THE FOLLOWING 11 INPUT LINES ARE REQUIRED Spectra Spectra 0 1.00000- 6 0 Spectra 1 ENDER IN Spectra ENDFB.OUT Spectra 92000 3 0 92999 3999 Spectra 90232 3 0 030 (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 Spectra LEAVE THE DEFINITION OF THE FILENAMES BLANK - THE PROGRAM WILL Spectra THEN USE STANDARD FILENAMES. Spectra Spectra IN THIS CASE THE FOLLOWING 9 INPUT LINES ARE REQUIRED Spectra Spectra 1 0 1.00000- 6 0 Spectra (USE DEFAULT FILENAME = ENDFB.IN) Spectra (USE DEFAULT FILENAME = ENDFB.OUT) Spectra 92000 1451 92999 1451 Spectra 92000 3 0 92999 3999 Spectra 90232 1451 0 1451 Spectra 90232 3 0 030 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999) Spectra (END OF REQUEST LIST) Spectra (0.1 PER-CENT ERROR, END OF ERROR LAW) Spectra Spectra EXAMPLE INPUT NO. 3 Spectra Spectra _____ LINEARIZE ALL MATERIALS ON AN ENDF/B TAPE TO WITHIN AN ACCURACY Spectra OF 0.5 PER-CENT (0.005 AS A FRACTION). IN THIS CASE YOU NEED NOT Spectra SPECIFY THE MAT, MF, MT RANGES. Spectra Spectra READ THE ENDF/B DATA FROM \ENDFB6\ZA092238 AND WRITE THE ENDF/B Spectra DATA TO \ENDFB6\LINEAR\ZA092238. Spectra

	Spectra
IN THIS CASE THE FOLLOWING 6 INPUT LINES ARE REQUIRED	Spectra
	Spectra
(MAT, 1.0E-10 BARNS, THIN)	Spectra
\ENDFB6\ZA092238	Spectra
\ENDFB6\LINEAR\ZA092238	Spectra
(RETRIEVE ALL DATA, END REQUEST LIST)	Spectra
5.00000-03	Spectra
(END OF ERROR LAW)	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