

=====		SPECTRA
PROGRAM SPECTRA		SPECTRA
=====		SPECTRA
An extension of LINEAR to linearize ALL MF=5 spectra.		SPECTRA
2012/05/28 - Added MF=15 neutron induced, photon spectra.		SPECTRA
2021/01/26 - MF=6 Still NOT Implemented		SPECTRA
		SPECTRA
First released in 2010 - Earlier below dates refer to LINEAR.		SPECTRA
		SPECTRA
VERSION 74-1 (MAY 1974)		SPECTRA
VERSION 75-1 (APRIL 1975)		SPECTRA
VERSION 76-2 (OCTOBER 1976)		SPECTRA
VERSION 77-1 (JANUARY 1977)		SPECTRA
VERSION 78-1 (JULY 1978)		SPECTRA
VERSION 79-1 (JULY 1979) CDC-7600 AND CRAY-1 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.	SPECTRA
	*NEW, MORE COMPATIBLE I/O UNIT NUMBER.	SPECTRA
	*ADDED OPTION TO KEEP ALL ORIGINAL	SPECTRA
	ENERGY POINTS FROM EVALUATION.	SPECTRA
	*ADDED STANDARD ALLOWABLE ERROR OPTIONS	SPECTRA
	(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	SPECTRA
VERSION 88-1 (JULY 1988) *OPTION...INTERNALLY 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	SPECTRA
	ROUTINE.	SPECTRA
	*WARNING...INPUT PARAMETER FORMAT	SPECTRA
	HAS BEEN CHANGED...SEE 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
	*INCREASED PAGE SIZE FROM 3006 TO 5010	SPECTRA
	POINTS.	SPECTRA
	*ALL ENERGIES INTERNALLY ROUNDED PRIOR	SPECTRA
	TO CALCULATIONS.	SPECTRA
	*COMPLETELY CONSISTENT I/O AND ROUNDINGS	SPECTRA
	ROUTINES - TO MINIMIZE COMPUTER	SPECTRA

	DEPENDENCE.	SPECTRA
VERSION 92-2 (JULY 1992)	*CORRECTED CONVERSION OF NU-BAR FROM POLYNOMIAL TO TABULATED - COPY	SPECTRA
	SPONTANEOUS NU-BAR (BY DEFINITION THE SPONTANEOUS NU-BAR IS NOT AN ENERGY DEPENDENT QUANTITY).	SPECTRA
VERSION 93-1 (MARCH 1993)	*UPDATED FOR USE WITH LAHEY COMPILER ON IBM-PCS.	SPECTRA
	*INCREASED PAGE SIZE FROM 5010 TO 30000 POINTS	SPECTRA
VERSION 94-1 (JANUARY 1994)	*VARIABLE ENDF/B DATA FILENAMES TO ALLOW ACCESS TO FILE STRUCTURES (WARNING - INPUT PARAMETER FORMAT HAS BEEN CHANGED)	SPECTRA
	*CLOSE ALL FILES BEFORE TERMINATING (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 60000 POINTS	SPECTRA
VERSION 99-1 (MARCH 1999)	*CORRECTED CHARACTER TO FLOATING POINT READ FOR MORE DIGITS	SPECTRA
	*UPDATED TEST FOR ENDF/B FORMAT VERSION BASED ON RECENT FORMAT CHANGES	SPECTRA
	*GENERAL IMPROVEMENTS BASED ON USER FEEDBACK	SPECTRA
VERSION 99-2 (JUNE 1999)	*ASSUME ENDF/B-VI, NOT V, IF MISSING MF=1, MT-451.	SPECTRA
VERS. 2000-1 (FEBRUARY 2000)	*ADDED MF = 9 AND 10 LINEARIZATION	SPECTRA
	*GENERAL IMPROVEMENTS BASED ON USER FEEDBACK	SPECTRA
VERS. 2002-1 (MAY 2002)	*OPTIONAL INPUT PARAMETERS	SPECTRA
VERS. 2004-1 (JAN. 2004)	*GENERAL UPDATE BASED ON USER FEEDBACKS	SPECTRA
VERS. 2005-1 (JAN. 2005)	*ALWAYS KEEP ORIGINAL TABULATED 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 600,000 POINTS	SPECTRA
VERS. 2010-1 (JUNE 2010)	*ADDED MF = 5 - MF = 6 STILL PLANNED.	SPECTRA
	*72 CHARACTER FILE NAMES.	SPECTRA
	*ONLY PROCESS MF=5 - SKIP ALL OTHERS TO PREVENT CONFLICT WITH LINEAR THINNING.	SPECTRA
VERS. 2012-1 (Aug. 2012)	*Added MF=15, neutron induced photon spectra.	SPECTRA
	*Added CODENAME	SPECTRA
	*32 and 64 bit Compatible	SPECTRA
	*Added ERROR stop	SPECTRA
VERS. 2015-1 (Jan. 2015)	*Extended OUT9.	SPECTRA
	*Replaced ALL 3 way IF Statements.	SPECTRA
	*Corrected MF=15 Data - it was adding 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 conflict with possible RESERVED NAME	SPECTRA
	*All floating input parameters changed to character input + IN9 conversion.	SPECTRA
VERS. 2018-1 (Jan. 2018)	*On-line output for ALL ENDERROR	SPECTRA
VERS. 2019-1 (June 2019)	*Additional Interpolation Law Tests	SPECTRA
	*Check Maximum Tabulated Energy to insure it is the same for all MTs -	SPECTRA

if not, print WARNING messages. SPECTRA
*Corrected END Histogram linearized - SPECTRA
Previously deleted last point - ERRORS SPECTRA
to assume this has Y=0 - now keep SPECTRA
point, but insure Y = 0. SPECTRA
VERS. 2020-1 (Mar. 2020) *Added Target Isomer State SPECTRA
VERS. 2021-1 (Jan. 2021) *Updated for FORTRAN 2018 SPECTRA
VERS. 2023-1 (Feb. 2023) *Decreased page size from 3,000,000 SPECTRA
to 120,000 SPECTRA

OWNED, MAINTAINED AND DISTRIBUTED BY

THE NUCLEAR DATA SECTION
INTERNATIONAL ATOMIC ENERGY AGENCY
P.O. BOX 100
A-1400, VIENNA, AUSTRIA
EUROPE

ORIGINALLY WRITTEN BY

Dermott E. Cullen

PRESENT CONTACT INFORMATION

Dermott E. Cullen
1466 Hudson Way
Livermore, CA 94550
U.S.A.
Telephone 925-443-1911
E. Mail RedCullen1@Comcast.net
Website RedCullen1.net/HOMEPAGE.NEW

AUTHORS MESSAGE

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

AT THE PRESENT TIME WE ARE ATTEMPTING TO DEVELOP A SET OF COMPUTERS SPECTRA
INDEPENDENT PROGRAMS THAT CAN EASILY BE IMPLEMENTED ON ANY ONE SPECTRA
OF A WIDE VARIETY OF COMPUTERS. IN ORDER TO ASSIST IN THIS PROJECTS SPECTRA
IT WOULD BE APPRECIATED 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

PURPOSE

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

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

ENDF/B FORMAT

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

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
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
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
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
***** PROGRAM SPECTRA (2023-1) *****	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
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
THESE COMMENT LINES ARE ONLY ADDED TO EXISTING HOLLERITH SECTIONS,	SPECTRA
I.E., THIS PROGRAM WILL NOT CREATE A HOLLERITH SECTION. THE FORMATS	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
THE ENDF/B FORMAT THE DATA IS IN. WITHOUT HAVING A SECTION OF	SPECTRA
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
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
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 PROGRAMS	SPECTRA
YOU MAY USE PROGRAM DICTIN TO CREATE A CORRECT REACTION INDEX.	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
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 SECTIONS	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

FUNCTION OF UP TO 20 (ENERGY,ERROR) PAIRS AND LINEAR INTERPOLATIONSPECTRA
BETWEEN TABULATED POINTS. IF ONLY ONE TABULATED POINT IS GIVEN THESPECTRA
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

DEFAULT ALLOWABLE ERROR

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

COULOMB PENETRABILITY (INTERPOLATION LAW = 6)

INTRODUCED FOR ENDF/B-VI. THIS IS DEFINED AS, SPECTRA

$$\text{SIG}(E) = C1 * \text{EXP}(-C2 / \text{SQRT}(E - T))$$

THIS PROGRAM ONLY CONSIDERS EXOTHERMIC REACTIONS - T = 0 SPECTRA

$$\text{SIG}(E) = C1 * \text{EXP}(-C2 / \text{SQRT}(E))$$

WARNING...THIS INTERPOLATION LAW SHOULD ONLY BE USED FOR REACTIONS SPECTRA
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

UNIT DESCRIPTION SPECTRA
----- SPECTRA

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

OUTPUT FILES

UNIT DESCRIPTION SPECTRA
----- SPECTRA

- 3 OUTPUT REPORT (BCD - 120 CHARACTERS/RECORD) SPECTRA
- 11 FINAL ENDF/B DATA (BCD - 80 CHARACTERS/RECORD) SPECTRA

SCRATCH FILES

UNIT DESCRIPTION SPECTRA
----- SPECTRA

- 12 SCRATCH FILE (BINARY - 180000 WORDS/RECORD) SPECTRA

OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINE FILEIO)

UNIT FILE NAME SPECTRA
----- SPECTRA

- 2 SPECTRA.INP SPECTRA
- 3 SPECTRA.LST SPECTRA
- 10 ENDFB.IN SPECTRA
- 11 ENDFB.OUT SPECTRA
- 12 (SCRATCH) SPECTRA

INPUT PARAMETERS

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

FOR VERSIONS 90-1 AND LATER THIS PROGRAM WILL ALLOW THE USER TO SPECTRA
TO SPECIFY BY INPUT PARAMETERS EXACTLY WHAT SECTIONS OF DATA SPECTRA

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	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.	SPECTRA
		EACH TIME A PAGE OF DATA POINTS IS WRITTEN TO	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).	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
2	1-72	ENDF/B INPUT DATA FILENAME	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
	12-22	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 ERRORS	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

EXAMPLE INPUT NO. 1

 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

EXPLICITLY SPECIFY THE STANDARD FILENAMES. SPECTRA

IN THIS CASE THE FOLLOWING 11 INPUT LINES ARE REQUIRED 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

EXAMPLE INPUT NO. 2

 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

LEAVE THE DEFINITION OF THE FILENAMES BLANK - THE PROGRAM WILL SPECTRA
 THEN USE STANDARD FILENAMES. SPECTRA

IN THIS CASE THE FOLLOWING 9 INPUT LINES ARE REQUIRED 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 0 3 0 (UPPER LIMIT AUTOMATICALLY SET TO 90232 3999) SPECTRA
 (END OF REQUEST LIST) SPECTRA
 (0.1 PER-CENT ERROR, END OF ERROR LAW) SPECTRA

EXAMPLE INPUT NO. 3

 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
READ THE ENDF/B DATA FROM \ENDFB6\ZA092238 AND WRITE THE ENDF/B DATA TO \ENDFB6\LINEAR\ZA092238.	SPECTRA
IN THIS CASE THE FOLLOWING 6 INPUT LINES ARE REQUIRED	SPECTRA
(MAT, 1.0E-10 BARNs, THIN)	SPECTRA
\ENDFB6\ZA092238	SPECTRA
\ENDFB6\LINEAR\ZA092238	SPECTRA
5.00000-03 (RETRIEVE ALL DATA, END REQUEST LIST)	SPECTRA
(END OF ERROR LAW)	SPECTRA
NOTE THAT IN THIS CASE IF THE INPUT HAD SPECIFIED AN EQUIVALENT ENERGY DEPENDENT ERROR LAW BY GIVING A NUMBER OF ENERGY POINTS AT EACH OF WHICH THE ERROR IS 0.5 PER-CENT THE PROGRAM WOULD TAKE LONGER TO RUN (I.E., ONLY USE AN ENERGY DEPENDENT ERROR LAW WHEN IT IS NECESSARY).	SPECTRA
EXAMPLE INPUT NO. 4	SPECTRA
-----	SPECTRA
IN ORDER TO LINEARIZE ALL MATERIALS ON AN ENDF/B TAPE TO THE STANDARD OPTION OF 0.1 PER-CENT IT IS ADEQUATE TO INPUT A SET OF COMPLETELY BLANK LINES WHICH WILL AUTOMATICALLY INVOKE ALL OF THE STANDARD OPTIONS.	SPECTRA
LEAVE THE DEFINITION OF THE FILENAMES BLANK - THE PROGRAM WILL THEN USE STANDARD FILENAMES.	SPECTRA
IN THIS CASE THE FOLLOWING THREE INPUT LINES ARE REQUIRED	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