

```

===== Spectra
PROGRAM SPECTRA Spectra
===== Spectra
An extension of LINEAR to linearize ALL MF=5 spectra. Spectra
05/28/2012 - Added MF=15 neutron induced, photon spectra. Spectra

First released in 2010 - Earlier below dates refer to LINEAR. 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 OPTION 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 ROUNDING Spectra
                                ROUTINES - TO MINIMIZE COMPUTER Spectra
                                DEPENDENCE. Spectra

```


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 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)

OUTPUT FILES

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

SCRATCH FILES

UNIT	DESCRIPTION
12	SCRATCH FILE (BINARY - 180000 WORDS/RECORD)

OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINE FILEIO)

UNIT	FILE NAME
2	SPECTRA.INP
3	SPECTRA.LST
10	ENDFB.IN
11	ENDFB.OUT
12	(SCRATCH)

INPUT PARAMETERS

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 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  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 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

EXAMPLE INPUT NO. 1 Spectra
----- Spectra
RETRIEVE DATA BY ZA IN ORDER TO FIND ALL URANIUM ISOTOPES AND Spectra

```


EXAMPLE INPUT NO. 4

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.

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

IN THIS CASE THE FOLLOWING THREE INPUT LINES ARE REQUIRED

(MAT, 1.0E-10 BARNS, THIN)
(USE DEFAULT FILENAME = ENDFB.IN)
(USE DEFAULT FILENAME = ENDFB.OUT)
(RETRIEVE ALL DATA, END REQUEST LIST)
(0.1 PER-CENT ERROR, END OF ERROR LAW)

Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra
Spectra

=====