

=====
Groupie

Groupie

PROGRAM GROUPIE

Groupie

=====

Groupie

VERSION 76-1 (NOVEMBER 1976)

Groupie

VERSION 79-1 (OCTOBER 1979) CDC-7600 AND CRAY-1 VERSION.

Groupie

VERSION 80-1 (MAY 1980) IBM, CDC AND CRAY VERSION

Groupie

VERSION 81-1 (JANUARY 1981) EXTENSION TO 3000 GROUPS

Groupie

VERSION 81-2 (MARCH 1981) IMPROVED SPEED

Groupie

VERSION 81-3 (AUGUST 1981) BUILT-IN 1/E WEIGHTING SPECTRUM

Groupie

VERSION 82-1 (JANUARY 1982) IMPROVED COMPUTER COMPATIBILITY

Groupie

VERSION 83-1 (JANUARY 1983) *MAJOR RE-DESIGN.

Groupie

*ELIMINATED COMPUTER DEPENDENT CODING.

Groupie

*NEW, MORE COMPATIBLE I/O UNIT NUMBERS.

Groupie

*NEW MULTI-BAND LIBRARY BINARY FORMAT.

Groupie

VERSION 83-2 (OCTOBER 1983) ADDED OPTION TO ALLOW SIGMA-0 TO BE

Groupie

DEFINED EITHER AS MULTIPLES OF

Groupie

UNSHIELDED TOTAL CROSS SECTION IN EACH

Groupie

GROUP, OR POWERS OF 10 IN ALL GROUPS.

Groupie

VERSION 84-1 (APRIL 1984) ADDED MORE BUILT IN MULTIGROUP ENERGY

Groupie

STRUCTURES.

Groupie

VERSION 85-1 (APRIL 1985) *UPDATED FOR ENDF/B-VI FORMATS.

Groupie

*SPECIAL I/O ROUTINES TO GUARANTEE

Groupie

ACCURACY OF ENERGY.

Groupie

*DOUBLE PRECISION TREATMENT OF ENERGY

Groupie

(REQUIRED FOR NARROW RESONANCES).

Groupie

*MINIMUM TOTAL CROSS SECTION TREATMENT

Groupie

VERSION 85-2 (AUGUST 1985) *FORTRAN-77/H VERSION
Groupie
VERSION 86-1 (JANUARY 1986) *ENDF/B-VI FORMAT
Groupie
VERSION 86-2 (JUNE 1986) *BUILT-IN MAXWELLIAN, 1/E AND FISSION
Groupie
WEIGHTING SPECTRUM.
Groupie
VERSION 88-1 (JULY 1988) *OPTION...INTERNALLY DEFINE ALL I/O
Groupie
FILE NAMES (SEE, SUBROUTINES FILIO1
Groupie
FILIO2 FOR DETAILS).
Groupie
*IMPROVED BASED ON USER COMMENTS.
Groupie
VERSION 89-1 (JANUARY 1989) *PSYCHOANALYZED BY PROGRAM FREUD TO
Groupie
INSURE PROGRAM WILL NOT DO ANYTHING
Groupie
CRAZY.
Groupie
*UPDATED TO USE NEW PROGRAM CONVERT
Groupie
KEYWORDS.
Groupie
*ADDED LIVERMORE CIVIC COMPILER
Groupie
CONVENTIONS.
Groupie
VERSION 91-1 (JUNE 1991) *INCREASED PAGE SIZE FROM 1002 TO 5010
Groupie
POINTS
Groupie
*UPDATED BASED ON USER COMMENTS
Groupie
*ADDED FORTRAN SAVE OPTION
Groupie
*COMPLETELY CONSISTENT ROUTINE TO READ
Groupie
FLOATING POINT NUMBERS.
Groupie
VERSION 92-1 (JANUARY 1992) *ADDED RESONANCE INTEGRAL CALCULATION -
Groupie
UNSHIELDED AND/OR SHIELDED - FOR
Groupie
DETAILS SEE BELOW
Groupie
*INCREASED NUMBER OF ENERGY POINTS
Groupie
IN BUILT-IN SPECTRA - TO IMPROVE
Groupie
ACCURACY.
Groupie

Groupie		*ALLOW SELECTION OF ZA/MF/MT OR
Groupie		MAT/MF/MT RANGES - ALL DATA NOT
Groupie		SELECTED IS SKIPPED ON INPUT AND
Groupie		NOT WRITTEN AS OUTPUT.
Groupie		*COMPLETELY CONSISTENT I/O ROUTINES -
Groupie		TO MINIMIZE COMPUTER DEPENDENCE.
Groupie		*NOTE, CHANGES IN INPUT PARAMETER
Groupie		FORMAT - FOR ZA/MF/MT OR MAT/MF/MT
Groupie		RANGES.
Groupie	VERSION 92-2 (JUNE 1992)	*MULTIBAND PARAMETERS OUTOUT AS
Groupie		CHARACTER (RATHER THAN BINARY) FILE.
Groupie	VERSION 93-1 (APRIL 1993)	*INCREASED PAGE SIZE FROM 5010 TO
Groupie		30000 POINTS
Groupie		*ELIMINATED COMPUTER DEPENDENCE.
Groupie	VERSION 94-1 (JANUARY 1994)	*VARIABLE ENDF/B DATA FILENAMES
Groupie		TO ALLOW ACCESS TO FILE STRUCTURES
Groupie		(WARNING - INPUT PARAMETER FORMAT
Groupie		HAS BEEN CHANGED)
Groupie		*CLOSE ALL FILES BEFORE TERMINATING
Groupie		(SEE, SUBROUTINE ENDIT)
Groupie	VERSION 95-1 (JANUARY 1994)	*CORRECTED MAXWELLIAN WEIGHTING
Groupie		*CHANGING WEIGHTING SPECTRUM FROM
Groupie		0.1 TO 0.001 % UNCERTAINTY
Groupie	VERSION 96-1 (JANUARY 1996)	*COMPLETE RE-WRITE
Groupie		*IMPROVED COMPUTER INDEPENDEENCE
Groupie		*ALL DOUBLE PRECISION
Groupie		*ON SCREEN OUTPUT

Groupie		*UNIFORM TREATMENT OF ENDF/B I/O
Groupie		*IMPROVED OUTPUT PRECISION
Groupie		*DEFINED SCRATCH FILE NAMES
Groupie		*UP TO 1000 GROUP MULTI-BAND
Groupie		CALCULATION (PREVIOUSLY 175)
Groupie		*MAXIMUM NUMBER OF GROUPS REDUCED
Groupie		FROM 3,000 TO 1,000
Groupie		*UP TO 1000 MATERIALS
Groupie		(PREVIOUSLY 100)
Groupie		*CORRECTED USE OF MAXWELLIAN +
Groupie		1/E + FISSION SPECTRUM
Groupie		*ONLY 2 BAND VERSION DISTRIBUTED
Groupie		(CONTACT AUTHOR FOR DETAILS)
Groupie		*DEFINED SCRATCH FILE NAMES
Groupie	VERSION 99-1 (MARCH 1999)	*CORRECTED CHARACTER TO FLOATING
Groupie		POINT READ FOR MORE DIGITS
Groupie		*UPDATED TEST FOR ENDF/B FORMAT
Groupie		VERSION BASED ON RECENT FORMAT CHANGE
Groupie		*GENERAL IMPROVEMENTS BASED ON
Groupie		USER FEEDBACK
Groupie	VERSION 99-2 (JUNE 1999)	*ASSUME ENDF/B-VI, NOT V, IF MISSING
Groupie		MF=1, MT-451.
Groupie	VERS. 2000-1 (FEBRUARY 2000)	*ADDED MF=10, ACTIVATION CROSS SECTION
Groupie		PROCESSING.
Groupie		*GENERAL IMPROVEMENTS BASED ON
Groupie		USER FEEDBACK
Groupie	VERS. 2002-1 (FEBRUARY 2002)	*ADDED TART 700 GROUP STRUCTURE
Groupie		

Groupie		*ADDED VARIABLE SIGMA0 INPUT OPTION
Groupie	(MAY 2002)	*OPTIONAL INPUT PARAMETERS
Groupie	(NOV. 2002)	*ADDED SAND-II EXTENDED DOWN TO
Groupie		1.0D-5 EV.
Groupie	(JUNE 2003)	*CORRECTED SAND-II 620 AND 640 GROUP
Groupie		ENERGY BOUNDARIES DEFINITIONS.
Groupie	VERS. 2004-1 (SEPT. 2004)	*INCREASED PAGE SIZE FROM 30000 TO
Groupie		120000 POINTS
Groupie		*ADDED "OTHER" AS ADDITIONAL REACTION
Groupie		TO IMPROVE MULTI-BAND FITTING
Groupie		*ADDED ITERATION FOR "BEST" PARTIAL
Groupie		PARAMETERS.
Groupie		*DO NOT SKIP LOW TOTAL ENERGY RANGES
Groupie		WHEN DEFINING AVERAGE CROSS SECTIONS -
Groupie		THIS MAKES OUTPUT COMPATIBLE WITH
Groupie		ANY STANDARD AVERAGING PROCEDURE
Groupie	VERS. 2005-1 (JAN. 2005)	*ADDED OPTION TO CHANGE TEMPERATURE OF
Groupie		BUILT-IN STANDARD SPECTRUM.
Groupie	VERS. 2007-1 (JAN. 2007)	*CHECKED AGAINST ALL ENDF/B-VII.
Groupie		*INCREASED PAGE SIZE FROM 120,000 TO
Groupie		600,000 POINTS
Groupie	VERS. 2008-1 (JAN. 2008)	*72 CHARACTER FILE NAMES.
Groupie		*GENERAL UPDATES
Groupie	VERS. 2010-1 (Apr. 2010)	*INCREASED WEIGHTING SPECTRUM TO 30,000
Groupie		FROM 3,000 ENERGY POINTS.
Groupie		*ADDED OUTPUT TO PLOT/COMPARE SHIELDED
Groupie		AND UNSHIELDED CROSS SECTIONS.

VERS. 2011-1 (June 2011)	*Corrected TART 700 groups to extend up
Groupie	to 1 GeV (1,000 MeV) - previously it
Groupie	was ERRONEOUSLY cutoff at 20 MeV.
Groupie	
VERS. 2011-2 (Nov. 2011)	*Corrected TART 616 groups lowest
Groupie	energy from 1.0D-4 eV to 1.0D-5 eV.
Groupie	
Groupie	*Added TART 666 to 200 MeV (for TENDL).
Groupie	
Groupie	*Optional high energy cross section
Groupie	extension above tabulated energy range
Groupie	(either = 0 = standard, or constant)
Groupie	WARNING - ENDF/B standard convention
Groupie	is that the cross section = 0 where it
Groupie	is not explicitly defined - extension
Groupie	= 0 is standard, constant is NOT, so
Groupie	constant extension is NOT RECOMMENDED.
Groupie	
VERS. 2012-1 (Aug. 2012)	*Added CODENAME
Groupie	
Groupie	*32 and 64 bit Compatible
Groupie	
Groupie	*Added ERROR stop.
VERS. 2013-1 (Nov. 2013)	*Extended OUT9.
Groupie	
Groupie	*Uses OUT9, not OUT10 for energies.
VERS. 2015-1 (Jan. 2015)	*Corrected SPECTM - handle ALL included
Groupie	group structures, i.e., even those
Groupie	that start above thermal range by
Groupie	ALWAYS constructing weighting spectrum
Groupie	to be AT LEAST 1.0D-5 eV to 20 MeV.
Groupie	
Groupie	*Extended OUT9
Groupie	
Groupie	*Replaced ALL 3 way IF Statements.
Groupie	
Groupie	*Generalized TART Group Structures.

Groupie *Generalized SAND-II Group Structures.
Groupie *Extended SAND-II to 60, 150, 200 MeV.
Groupie **VERS. 2015-2 (Mar. 2015)** *Deleted 1P from formats reading input
Groupie parameters, causing incorrect scaling
Groupie *Changed ALL data to "D" instead of
Groupie "E" to insure it is REAL*8 and avoid
Groupie Truncation ERRORS.

Groupie
Groupie 2015-2 Acknowledgment

Groupie
Groupie =====

Groupie I thank Chuck Whitmer (TerraPower,WA) and Andrej Trkov (NDS,IAEA)
Groupie for reporting the errors that led to the 2015-2 Improvements in
Groupie this code.

Groupie
Groupie I thank Jean-Christophe Sublet (UKAEA) for contributing MAC
Groupie executables and Bojan Zefran (IJS, Slovenia) for contributing
Groupie LINUX (32 or 63 bit) executables. And most of all I must thank
Groupie Andrej Trkov (NDS, IAEA) for overseeing the entire PREPRO project
Groupie at IAEA, Vienna. This was a truly International team who worked
Groupie together to produce PREPRO 2015-2.

Groupie

Groupie OWNED, MAINTAINED AND DISTRIBUTED BY
Groupie

Groupie -----

Groupie THE NUCLEAR DATA SECTION
Groupie INTERNATIONAL ATOMIC ENERGY AGENCY

Groupie P.O. BOX 100
Groupie A-1400, VIENNA, AUSTRIA

Groupie

EUROPE
Groupie

Groupie
ORIGINALLY WRITTEN BY
Groupie

Groupie
Dermott E. Cullen
Groupie

Groupie
PRESENT CONTACT INFORMATION
Groupie

Groupie
Dermott E. Cullen
Groupie
1466 Hudson Way
Groupie
Livermore, CA 94550
Groupie
U.S.A.
Groupie
Telephone 925-443-1911
Groupie
E. Mail RedCullen1@Comcast.net
Groupie
Website <http://home.comcast.net/~redcullen1>
Groupie

Groupie
AUTHORS MESSAGE
Groupie

Groupie
THE REPORT DESCRIBED ABOVE IS THE LATEST PUBLISHED DOCUMENTATION
Groupie
FOR THIS PROGRAM. HOWEVER, THE COMMENTS BELOW SHOULD BE CONSIDERED
Groupie
THE LATEST DOCUMENTATION INCLUDING ALL RECENT IMPROVEMENTS. PLEASE
Groupie
READ ALL OF THESE COMMENTS BEFORE IMPLEMENTATION, PARTICULARLY
Groupie
THE COMMENTS CONCERNING MACHINE DEPENDENT CODING.
Groupie

Groupie
AT THE PRESENT TIME WE ARE ATTEMPTING TO DEVELOP A SET OF COMPUTER
Groupie
INDEPENDENT PROGRAMS THAT CAN EASILY BE IMPLEMENTED ON ANY ONE
Groupie
OF A WIDE VARIETY OF COMPUTERS. IN ORDER TO ASSIST IN THIS PROJECT
Groupie

IT WOULD BE APPRECIATED IF YOU WOULD NOTIFY THE AUTHOR OF ANY
Groupie
COMPILER DIAGNOSTICS, OPERATING PROBLEMS OR SUGGESTIONS ON HOW TO
Groupie
IMPROVE THIS PROGRAM. HOPEFULLY, IN THIS WAY FUTURE VERSIONS OF
Groupie
THIS PROGRAM WILL BE COMPLETELY COMPATIBLE FOR USE ON YOUR
Groupie
COMPUTER.

Groupie

Groupie
PURPOSE

Groupie

Groupie
THIS PROGRAM IS DESIGNED TO CALCULATE ANY COMBINATION OF
Groupie
THE FOLLOWING QUANTITIES FROM LINEARLY INTERPOLABLE TABULATED
Groupie
CROSS SECTIONS IN THE ENDF/B FORMAT

Groupie

Groupie
(1) UNSHIELDED GROUP AVERAGED CROSS SECTIONS

Groupie
(2) BONDARENKO SELF-SHIELDED GROUP AVERAGED CROSS SECTIONS

Groupie
(3) MULTI-BAND PARAMETERS

Groupie

Groupie
IN THE FOLLOWING FOR SIMPLICITY THE ENDF/B TERMINOLOGY--ENDF/B
Groupie
TAPE--WILL BE USED. IN FACT THE ACTUAL MEDIUM MAY BE TAPE, CARDS,
Groupie
DISK OR ANY OTHER MEDIUM.

Groupie

Groupie
ENDF/B FORMAT

Groupie

Groupie
THIS PROGRAM ONLY USES THE ENDF/B BCD OR CARD IMAGE FORMAT (AS
Groupie
OPPOSED TO THE BINARY FORMAT) AND CAN HANDLE DATA IN ANY VERSION
Groupie
OF THE ENDF/B FORMAT (I.E., ENDF/B-I, II, III, IV OR V FORMAT).

Groupie

Groupie
IT IS ASSUMED THAT THE DATA IS CORRECTLY CODED IN THE ENDF/B

Groupie

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

Groupie
ALL FILE 3 CROSS SECTIONS THAT ARE USED BY THIS PROGRAM MUST BE
Groupie
LINEARLY INTERPOLABLE IN ENERGY AND CROSS SECTION (ENDF/B
Groupie
INTERPOLATION LAW 2). FILE 3 BACKGROUND CROSS SECTIONS MAY BE MADE
Groupie
LINEARLY INTERPOLABLE USING PROGRAM LINEAR (UCRL-50400, VOL. 17,
Groupie
PART A). THE RESONANCE CONTRIBUTION MAY BE ADDED TO THE BACKGROUND
Groupie
CROSS SECTIONS USING PROGRAM RECENT (UCRL-50400, VOL. 17, PART B).
Groupie
IF THIS PROGRAM FINDS THAT THE FILE 3 CROSS SECTIONS ARE NOT
Groupie
LINEARLY INTERPOLABLE THIS PROGRAM WILL TERMINATE EXECUTION.
Groupie

Groupie
CONTENTS OF OUTPUT
Groupie

Groupie
IF ENDF/B FORMATTED OUTPUT IS REQUESTED ENTIRE EVALUATIONS ARE
Groupie
OUTPUT, NOT JUST THE MULTI-GROUPED FILE 3 CROSS SECTIONS, E.G.
Groupie
ANGULAR AND ENERGY DISTRIBUTIONS ARE ALSO INCLUDED.
Groupie

Groupie
DOCUMENTATION
Groupie

Groupie
THE FACT THAT THIS PROGRAM HAS OPERATED ON THE DATA IS DOCUMENTED
Groupie
BY THE ADDITION OF THREE COMMENT CARDS AT THE END OF EACH
Groupie

HOLLERITH SECTION TO DESCRIBE THE GROUP STRUCTURE AND WEIGHTING
Groupie
SPECTRUM, E.G.
Groupie

Groupie
***** PROGRAM GROUPIE (2015-2) *****
Groupie
UNSHIELDED GROUP AVERAGES USING 69 GROUPS (WIMS)
Groupie
MAXWELLIAN, 1/E AND FISSION WEIGHTING SPECTRUM
Groupie

Groupie
THE ORDER OF ALL SIMILAR COMMENTS (FROM LINEAR, RECENT AND SIGMA1)
Groupie
REPRESENTS A COMPLETE HISTORY OF ALL OPERATIONS PERFORMED ON
Groupie
THE DATA.
Groupie

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

Groupie
REACTION INDEX
Groupie

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

THIS PROGRAM DOES NOT UPDATE THE REACTION INDEX IN MF=1, MT=451.
Groupie
THIS CONVENTION HAS BEEN ADOPTED BECAUSE MOST USERS DO NOT
Groupie
REQUIRE A CORRECT REACTION INDEX FOR THEIR APPLICATIONS AND IT WAS
Groupie
NOT CONSIDERED WORTHWHILE TO INCLUDE THE OVERHEAD OF CONSTRUCTING
Groupie
A CORRECT REACTION INDEX IN THIS PROGRAM. HOWEVER, IF YOU REQUIRE
Groupie
A REACTION INDEX FOR YOUR APPLICATIONS, AFTER RUNNING THIS PROGRAM
Groupie
YOU MAY USE PROGRAM DICTIN TO CREATE A CORRECT REACTION INDEX.
Groupie

Groupie
SECTION SIZE

Groupie

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

Groupie
SELECTION OF DATA

Groupie

Groupie
THE PROGRAM SELECTS MATERIALS TO BE PROCESSED BASED EITHER ON
Groupie
MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR
Groupie
ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE
Groupie
ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS
Groupie
USED TO SELECT MATERIALS, AND WILL TERMINATE WHEN A MAT OR ZA
Groupie
IS FOUND THAT IS ABOVE THE RANGE OF ALL REQUESTS.
Groupie

Groupie
ENERGY ORDER AND UNITS

Groupie

Groupie
ALL ENERGIES (FOR CROSS SECTIONS, WEIGHTING SPECTRUM OR GROUP
Groupie
BOUNDARIES) MUST BE IN UNITS OF EV AND MUST BE IN ASCENDING
Groupie

NUMERICAL ORDER.
Groupie

Groupie
ENERGY GRID
Groupie

Groupie
ALTHOUGH ALL REACTIONS MUST TO LINEARLY INTERPOLABLE, THEY DO NOT
Groupie
ALL HAVE TO USE THE SAME ENERGY GRID. EACH REACTION CAN BE GIVEN
Groupie
BY AN INDEPENDENT ENERGY GRID. THIS PROGRAM WILL PROCEED FROM
Groupie
THE LOWEST TO HIGHEST ENERGY SELECTING EACH ENERGY INTERVAL OVER
Groupie
WHICH ALL DATA, FOR ANY GIVEN CALCULATION, ARE ALL LINEARLY
Groupie
INTERPOLABLE.
Groupie

Groupie
GROUP STRUCTURE
Groupie

Groupie
THIS PROGRAM IS DESIGNED TO USE AN ARBITRARY ENERGY GROUP
Groupie
STRUCTURE WHERE THE ENERGIES ARE IN EV AND ARE IN INCREASING
Groupie
ENERGY ORDER. THE MAXIMUM NUMBER OF GROUPS IS 1000.
Groupie

Groupie
THE USER MAY INPUT AN ARBITRARY GROUP STRUCTURE OR THE USER MAY
Groupie
USE USE ONE OF THE SEVEN BUILT-IN GROUP STRUCTURES.
Groupie
(0) 175 GROUP (TART STRUCTURE)
Groupie
(1) 50 GROUP (ORNL STRUCTURE)
Groupie
(2) 126 GROUP (ORNL STRUCTURE)
Groupie
(3) 171 GROUP (ORNL STRUCTURE)
Groupie
(4) 620 GROUP (SAND-II STRUCTURE, UP TO 18 MEV)
Groupie
(5) 640 GROUP (SAND-II STRUCTURE, UP TO 20 MEV)
Groupie
(6) 69 GROUP (WIMS STRUCTURE)
Groupie
(7) 68 GROUP (GAM-I STRUCTURE)
Groupie

(8) 99 GROUP (GAM-II STRUCTURE)
 Groupie
 (9) 54 GROUP (MUFT STRUCTURE)
 Groupie
 (10) 28 GROUP (ABBN STRUCTURE)
 Groupie
 (11) 616 GROUP (TART STRUCTURE TO 20 MeV)
 Groupie
 (12) 700 GROUP (TART STRUCTURE TO 1 GEV)
 Groupie
 (13) 665 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 18 MEV)
 Groupie
 (14) 685 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 20 MEV)
 Groupie
 (15) 666 GROUP (TART STRUCTURE TO 200 MeV)
 Groupie
 (16) 725 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 60 MEV)
 Groupie
 (17) 755 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 150 MEV)
 Groupie
 (18) 765 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 200 MEV)
 Groupie

Groupie
 GROUP AVERAGES
 Groupie

Groupie
 THIS PROGRAM DEFINES GROUP AVERAGED CROSS SECTIONS AS...
 Groupie

Groupie

$$\text{AVERAGE} = \frac{\text{(INTEGRAL E1 TO E2) (SIGMA(E) * S(E) * WT(E) * DE)}}{\text{(INTEGRAL E1 TO E2) (S(E) * WT(E) * DE)}}$$
 Groupie

Groupie
 WHERE...
 Groupie

Groupie
 AVERAGE = GROUP AVERAGED CROSS SECTION
 Groupie
 E1, E2 = ENERGY LIMITS OF THE GROUP
 Groupie
 SIGMA(E) = ENERGY DEPENDENT CROSS SECTION FOR ANY GIVEN REACTION
 Groupie
 S(E) = ENERGY DEPENDENT WEIGHTING SPECTRUM
 Groupie
 WT(E) = ENERGY DEPENDENT SELF-SHIELDING FACTOR.
 Groupie

Groupie

ENERGY DEPENDENT WEIGHTING SPECTRUM

Groupie -----

Groupie THE ENERGY DEPENDENT WEIGHTING SPECTRUM IS GIVEN BY AN ARBITRARY

Groupie TABULATED LINERLY INTERPOLABLE FUNCTION WHICH CAN BE DESCRIBED

Groupie BY AN ARBITRARY NUMBER OF POINTS. THIS ALLOWS THE USER TO

Groupie SPECIFY ANY DESIRED WEIGHTING SPECTRUM TO ANY GIVEN DEGREE OF

Groupie ACCURACY. REMEMBER THAT THE PROGRAM WILL ASSUME THAT THE SPECTRUM

Groupie IS LINEARLY INTERPOLABLE BETWEEN TABULATED POINTS. THEREFORE THE

Groupie USER SHOULD USE ENOUGH POINTS TO INSURE AN ADEQUATE REPRESENTATION

Groupie OF THE SPECTRUM BETWEEN TABULATED DATA POINTS.

Groupie

Groupie THE PRESENT VERSION OF THE CODE HAS THREE BULIT-IN WEIGHTING

Groupie SPECTRA,

Groupie

Groupie (1) CONSTANT

Groupie (2) 1/E

Groupie (3) MAXWELLIAN = $E \cdot \exp(-E/KT) / KT$ (0.0 TO $4 \cdot KT$)

Groupie 1/E = $C1/E$ ($4 \cdot KT$ TO 67 KEV)

Groupie FISSION = $C2 \cdot \exp(-E/WA) \cdot \sinh(\sqrt{E \cdot WB})$ (ABOVE 67 KEV)

Groupie

Groupie $KT = 0.253 \text{ EV (293 KELVIN)}$

Groupie $WA = 9.65D+5$

Groupie $WB = 2.29D-6$

Groupie $C1, C2 = \text{DEFINED TO MAKE SPECTRUM CONTINUOUS}$

Groupie

Groupie FISSION SPECTRUM CONSTANTS FROM

Groupie A.F.HENRY, NUCLEAR REACTOR ANALYSIS, P. 11, MIT PRESS (1975)

Groupie

Groupie
 UNSHIELDED GROUP AVERAGES
 Groupie

 Groupie
 FOR UNSHIELDED AVERAGES THE SELF-SHIELDING FACTOR (WT(E)) IS SET
 Groupie
 TO UNITY. THIS PROGRAM ALLOWS UP TO 1000 GROUPS.
 Groupie

Groupie
 SELF-SHIELDED GROUP AVERAGES
 Groupie

 Groupie
 IF SELF-SHIELDED AVERAGES AND/OR MULTI-BAND PARAMETERS ARE
 Groupie
 CALCULATED THIS PROGRAM ALLOWS UP TO 1000 GROUPS. SELF-SHIELDED
 Groupie
 AVERAGES AND/OR MULTI-BAND PARAMETERS ARE CALCULATED FOR THE
 Groupie
 TOTAL, ELASTIC, CAPTURE AND FISSION.
 Groupie

Groupie
 FOR THE TOTAL, ELASTIC, CAPTURE AND FISSION THE PROGRAM USES A
 Groupie
 WEIGHTING FUNCTION THAT IS A PRODUCT OF THE ENERGY DEPENDENT
 Groupie
 WEIGHTING SPECTRUM TIMES A BONDERENKO TYPE SELF-SHIELDING FACTOR.
 Groupie

Groupie

$$WT(E) = S(E) / (TOTAL(E) + SIGMA0) ** N$$

 Groupie

Groupie
 WHERE...
 Groupie

Groupie
 S(E) - ENERGY DEPENDENT WEIGHTING SPECTRUM (DEFINED BY
 Groupie
 TABULATED VALUES AND LINEAR INTERPOLATION BETWEEN
 Groupie
 TABULATED VALUES).
 Groupie
 TOTAL(E) - ENERGY DEPENDENT TOTAL CROSS SECTION FOR ONE MATERIAL
 Groupie
 (DEFINED BY TABULATED VALUES AND LINEAR INTERPOLATION
 Groupie
 BETWEEN TABULATED VALUES).
 Groupie

SIGMA0 - CROSS SECTION TO REPRESENT THE EFFECT OF ALL OTHER
 Groupie
 MATERIALS AND LEAKAGE (DEFINED WITHIN EACH GROUP TO BE
 Groupie
 A MULTIPLE OF THE UNSHIELDED TOTAL CROSS SECTION WITHIN
 Groupie
 THAT GROUP OR POWERS OF 10 - INPUT OPTION).
 Groupie
 N - A POSITIVE INTEGER (0, 1, 2 OR 3).
 Groupie
 Groupie
 THE PROGRAM WILL USE ONE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E)
 Groupie
 AND 25 DIFFERENT BONDERENKO TYPE SELF-SHIELDING FACTORS (25 SIGMA0
 Groupie
 AND N COMBINATIONS) TO DEFINE 25 DIFFERENT AVERAGE CROSS SECTIONS,
 Groupie
 FOR EACH REACTION, WITHIN EACH GROUP.
 Groupie
 Groupie
 THE 25 WEIGHTING FUNCTIONS USED ARE....
 Groupie
 (1) - UNSHIELDED CROSS SECTIONS (N=0)
 Groupie
 (2-22)- PARTIALLY SHIELDED CROSS SECTIONS (N=1 ,VARIOUS SIGMA0)
 Groupie
 THE VALUES OF SIGMA0 USED WILL BE EITHER,
 Groupie
 (A) THE VALUES OF SIGMA0 THAT ARE USED VARY FROM 1024
 Groupie
 TIMES THE UNSHIELDED TOTAL CROSS SECTIONS IN STEPS OF 1/2
 Groupie
 DOWN TO 1/1024 TIMES THE UNSHIELDED TOTAL CROSS SECTION
 Groupie
 (A RANGE OF OVER 1 MILLION, CENTERED ON THE UNSHIELDED
 Groupie
 TOTAL CROSS SECTION WITHIN EACH GROUP).
 Groupie
 (B) THE SAME CONSTANT VALUES OF SIGMA0 IN EACH GROUP. THE
 Groupie
 VALUES OF SIGMA0 USED INCLUDE 40000, 20000, 10000, 7000,
 Groupie
 4000, 2000, 1000, 700, 400, 200, 100, 70, 40, 20, 10, 7,
 Groupie
 4, 2, 1, 0.7, 0.4 (A RANGE OF 100,000 SPANNING MORE THAN
 Groupie
 THE RANGE OF SIGMA0 VALUES THAT MAY BE ENCOUNTERED IN
 Groupie
 ACTUAL APPLICATIONS)
 Groupie
 (23) - TOTALLY SHIELDED FLUX WEIGHTED CROSS SECTION
 Groupie

(N=1, SIGMA0=0)
Groupie
(24) - TOTALLY SHIELDED CURRENT WEIGHTED CROSS SECTION
Groupie
(N=2, SIGMA0=0)
Groupie
(25) - TOTALLY SHIELDED COSINE SQUARED WEIGHTED CROSS SECTION
Groupie
(N=3, SIGMA0=0)
Groupie

Groupie
FOR ALL OTHER REACTIONS (EXCEPT TOTAL, ELASTIC, CAPTURE AND
Groupie
FISSION) THE PROGRAM WILL USE THE ENERGY DEPENDENT WEIGHTING
Groupie
SPECTRUM S(E) TO DEFINE THE UNSHIELDED (BONDERENKO N=0)
Groupie
AVERAGED CROSS SECTION WITHIN EACH GROUP.
Groupie

Groupie
CALCULATION OF RESONANCE INTEGRALS
Groupie

Groupie
IN A PURE ELASTIC ISOTROPICALLY SCATTERING MATERIAL WITH A
Groupie
CONSTANT CROSS SECTION THE SPECTRUM WILL BE 1/E AND THERE WILL
Groupie
BE NO SELF-SHIELDING.
Groupie

Groupie
IN THIS CASE IF THE CROSS SECTION VARIES WITH ENERGY THE
Groupie
SPECTRUM WILL STILL BE 1/E AND THE SELF-SHIELDING FACTOR WILL
Groupie
BE EXACTLY 1/SIG-TOT(E) - WHERE SIG-TOT(E) = SIG-EL(E), SINCE
Groupie
THERE IS ONLY SCATTERING.
Groupie

Groupie
IF WE HAVE AN INFINITELY DILUTE AMOUNT OF A MATERIAL UNIFORMLY
Groupie
MIXED WITH A PURE ELASTIC ISOTROPICALLY SCATTERING MATERIAL WITH
Groupie
A CONSTANT CROSS SECTION THE STANDARD DEFINITION OF THE RESONANCE
Groupie
INTEGRAL CAN BE USED TO DEFINE REACTION RATES FOR EACH REACTION.
Groupie

Groupie

THE RESONANCE INTEGRAL IS DEFINED AS,
Groupie

Groupie
$$RI = (\text{INTEGRAL } E1 \text{ TO } E2) (\text{SIGMA}(E) * S(E) * WT(E) * DE)$$

Groupie

Groupie
WHERE NORMALLY,

Groupie
$$S(E) = 1/E$$

Groupie
$$WT(E) = 1 \quad - \text{ NO SELF-SHIELDING}$$

Groupie

Groupie
FROM THE ABOVE DEFINITION OF GROUP AVERAGED CROSS SECTIONS THE

Groupie
RESONANCE INTEGRAL IS,

Groupie

Groupie
$$RI = \text{AVERAGE} * (\text{INTEGRAL } E1 \text{ TO } E2) (S(E) * WT(E) * DE)$$

Groupie

Groupie
FOR A 1/E SPECTRUM AND NO SELF-SHIELDING THIS REDUCES TO,

Groupie

Groupie
$$RI = \text{AVERAGE} * \text{LOG}(E2/E1)$$

Groupie

Groupie
IN ANY OTHER SITUATION, INCLUDING ABSORPTION AND/OR ENERGY

Groupie
DEPENDENT CROSS SECTIONS, THE SPECTRUM WILL NOT BE 1/E -

Groupie
ABSORPTION WILL TEND TO DECREASE THE SPECTRUM PROGRESSIVELY

Groupie
MORE AT LOWER ENERGIES - ENERGY DEPENDENCE OF THE CROSS SECTION

Groupie
WILL LEAD TO SELF-SHIELDING.

Groupie

Groupie
HERE WE WILL NOT ATTEMPT TO PERFORM A DETAILED SPECTRUM

Groupie
CALCULATION TO ACCOUNT FOR ABSORPTION.

Groupie

Groupie
HOWEVER, WE WILL EXTEND THE DEFINITION OF THE RESONANCE INTEGRAL

Groupie

TO ACCOUNT FOR SELF-SHIELDING EFFECTS BY ALLOWING FOR INCLUSION
Groupie
OF SELF-SHIELDING EFFECTS IN THE DEFINITION OF GROUP AVERAGES
Groupie
AND THEN DEFINING THE RESONANCE INTEGRAL AS,
Groupie

Groupie
RI = AVERAGE* LOG(E2/E1)
Groupie

Groupie
IN ORDER TO CALCULATE RESONANCE INTEGRALS YOU MUST FOLLOW THESE
Groupie
STEPS,
Groupie

- Groupie
- 1) SELECT A 1/E SPECTRUM - ON FIRST LINE OF INPUT PARAMETERS.
Groupie
 - 2) SELECT THE ENERGY BOUNDARIES - NORMALLY ONLY 1 GROUP FROM
Groupie
0.5 EV UP TO 20 MEV - HOWEVER, YOU ARE FREE TO SELECT ANY
Groupie
ENERGY RANGE THAT YOU WISH - YOU MAY EVEN SELECT MORE THAN
Groupie
1 GROUP MERELY BY SPECIFYING MORE THAN 1 GROUP AS INPUT -
Groupie
THIS CAN BE USED TO DEFINE THE CONTRIBUTIONS TO THE RESONANCE
Groupie
INTEGRAL FROM INDIVIDUAL ENERGY RANGES.
Groupie
 - 3) SELECT THIS OPTION FOR THE UNSHIELDED AND/OR SHIELDED OUTPUT
Groupie
LISTING - ON THE SECOND LINE OF INPUT PARAMETERS.
Groupie

Groupie
WHEN THIS OPTION IS USED THE PROGRAM WILL CALCULATE GROUP AVERAGED
Groupie
CROSS SECTIONS - AS DEFINED ABOVE - PRIOR TO OUTPUT THE RESULTS
Groupie
WILL MERELY BE MULTIPLIED BY THE WIDTH OF THE GROUP ASSUMING YOU
Groupie
HAVE SELECTED A 1/E SPECTRUM - THERE IS NO CHECK ON THIS - THE
Groupie
PROGRAM MERELY MULTIPLIES THE GROUP AVERAGED CROSS SECTIONS BY,
Groupie

Groupie
LOG(E2/E1) - WHERE E2 AND E1 ARE THE GROUP ENERGY BOUNDARIES.
Groupie

Groupie

WARNING - IT IS UP TO YOU TO INSURE THAT YOU FOLLOW EXACTLY THE
STEPS OUTLINED ABOVE IF YOU WISH TO OBTAIN MEANINGFUL
RESULTS.

NOTE - OUTPUT IN THE ENDF/B FORMAT IS ALWAYS GROUP AVERAGED CROSS
SECTIONS, REGARDLESS OF WHETHER YOU ASK FOR AVERAGED CROSS
SECTIONS OR RESONANCE INTEGRALS - THIS IS BECAUSE DATA IN
THE ENDF/B FORMAT IS EXPLICITLY DEFINED TO BE CROSS
SECTIONS.

RESONANCE INTEGRAL OUTPUT CAN ONLY BE OBTAINED IN THE
LISTING FORMATS.

MINIMUM TOTAL CROSS SECTION TREATMENT

SINCE THE BONDARENKO SELF-SHIELDING DEPENDS ON 1/TOTAL CROSS
SECTION, THE ALGORITHM WILL BECOME NUMERICALLY UNSTABLE IF THE
TOTAL CROSS SECTION IS NEGATIVE (AS OCCURS IN MANY ENDF/B
EVALUATIONS). IF THE TOTAL IS LESS THAN SOME MINIMUM ALLOWABLE
VALUE (DEFINE BY OKMIN, PRESENTLY 1 MILLI-BARN) AN ERROR MESSAGE
WILL BE PRINTED AND FOR THE SELF-SHIELDING CALCULATION ALL ENERGY
INTERVALS IN WHICH THE TOTAL IS LESS THAN THE MINIMUM WILL BE
IGNORED.

NOTE, FOR THE UNSHIELDED CALCULATIONS ALL CROSS SECTIONS WILL BE
CONSIDERED WHETHER THEY ARE POSITIVE OR NEGATIVE. THEREFORE IF
THE TOTAL CROSS SECTION IS NEGATIVE OR LESS THAN THE MINIMUM

VALUE THERE MAY BE AN INCONSISTENCY BETWEEN THE UNSHIELDED AND
Groupie
THE SELF-SHIELDED CROSS SECTIONS. IF THE TOTAL CROSS SECTION IS
Groupie
NEGATIVE AND SELF-SHIELDED CROSS SECTIONS ARE CALCULATED THE
Groupie
PROGRAM WILL PRINT AN ERROR MESSAGE INDICATING THAT THE SELF-
Groupie
SHIELDED RESULTS ARE UNRELIABLE AND SHOULD NOT BE USED. THEREFORE
Groupie
IN THIS CASE THE PROGRAM WILL NOT ATTEMPT TO MODIFY THE UNSHIELDED
Groupie
RESULTS TO ELIMINATE THE EFFECT OF NEGATIVE CROSS SECTIONS, SINCE
Groupie
THE UNSHIELDED RESULTS ARE THE ONLY ONES WHICH TRULY REFLECT THE
Groupie
ACTUAL INPUT.

Groupie

Groupie
RESOLVED RESONANCE REGION

Groupie

Groupie

IN THE RESOLVED RESONANCE REGION (ACTUALLY EVERYWHERE BUT IN THE
Groupie
UNRESOLVED RESONANCE REGION) THE CROSS SECTIONS OUTPUT BY LINEAR-
Groupie
RECENT-SIGMA1 WILL BE ACTUAL ENERGY DEPENDENT CROSS SECTIONS AND
Groupie
THE CALCULATIONS BY THIS PROGRAM WILL YIELD ACTUAL SHIELDED AND
Groupie
UNRESOLVED CROSS SECTIONS.

Groupie

Groupie
UNRESOLVED RESONANCE REGION

Groupie

Groupie

IN THE UNRESOLVED RESONANCE REGION PROGRAM RECENT USES THE
Groupie
UNRESOLVED RESONANCE PARAMETERS TO CALCULATE INFINITELY DILUTE
Groupie
AVERAGE CROSS SECTIONS. THIS PROGRAM WILL MERELY READ THIS
Groupie
INFINITELY DILUTE DATA AS IF IT WERE ENERGY DEPENDENT DATA AND
Groupie
GROUP AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT
Groupie
UNRESOLVED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT
Groupie
IT WILL NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS.

Groupie

Groupie

ACCURACY OF RESULTS

Groupie

Groupie

ALL INTEGRALS ARE PERFORMED ANALYTICALLY. THEREFORE NO ERROR IS

Groupie

INTRODUCED DUE TO THE USE OF TRAPAZOIDAL OR OTHER INTEGRATION

Groupie

SCHEME. THE TOTAL ERROR THAT CAN BE ASSIGNED TO THE RESULTING

Groupie

AVERAGES IS JUST THAT DUE TO THE ERROR IN THE CROSS SECTIONS

Groupie

AND ENERGY DEPENDENT WEIGHTING SPECTRUM. GENERALLY SINCE THE

Groupie

THE ENERGY DEPENDENT WEIGHTING SPECTRUM APPEARS IN BOTH THE

Groupie

NUMERATOR AND THE DENOMINATOR THE AVERAGES RAPIDLY BECOME

Groupie

INSENSITIVE TO THE WEIGHTING SPECTRUM AS MORE GROUPS ARE USED.

Groupie

SINCE THE WEIGHTING SPECTRUM IS LOADED IN THE PAGING SYSTEM THE

Groupie

USER CAN DESCRIBE THE SPECTRUM TO ANY REQUIRED ACCURACY USING

Groupie

ANY NUMBER OF ENERGY VS. SPECTRUM PAIRS.

Groupie

Groupie

MULTI-BAND PARAMETERS

Groupie

Groupie

MULTI-BAND PARAMETERS ARE CALCULATED FOR THE TOTAL, ELASTIC,

Groupie

CAPTURE AND FISSION REACTIONS. WITH THE NUMBER OF GROUPS THAT

Groupie

ARE NORMALLY USED (SEE BUILT IN GROUP STRUCTURES) ALL OTHER

Groupie

REACTIONS RESULT IN A NEGLIGABLE AMOUNT OF SELF-SHIELDING. AS

Groupie

SUCH THEIR EQUIVALENT BAND CROSS SECTION WILL MERELY BE THEIR

Groupie

UNSHIELDED VALUE WITHIN EACH BAND.

Groupie

Groupie

FOR ANY GIVEN EVALUATION, WITHIN ANY GIVEN GROUP THIS PROGRAM

Groupie

WILL GENERATE THE MINIMUM NUMBER OF BANDS REQUIRED WITHIN THAT

Groupie

GROUP. AS OUTPUT TO THE COMPUTER READABLE DISK FILE THE BAND

Groupie

PARAMETERS FOR EACH EVALUATION WILL BE FORMATTED TO HAVE THE
Groupie SAME NUMBER OF BANDS IN ALL GROUPS (WITH ZERO WEIGHT FOR SOME
Groupie BANDS WITHIN ANY GROUP). THE USER MAY DECIDE TO HAVE OUTPUT
Groupie EITHER WITH THE MINIMUM NUMBER OF BANDS REQUIRED FOR EACH
Groupie EVALUATION (E.G. 2 BANDS FOR HYDROGEN AND 4 BANDS FOR U-233) OR
Groupie THE SAME NUMBER OF BANDS FOR ALL EVALUATIONS (E.G. 4 BANDS FOR
Groupie BOTH HYDROGEN AND U-233).

Groupie
Groupie FOR 2 OR FEWER BANDS THE PROGRAM USES AN ANALYTIC EXPRESSION
Groupie TO DEFINE ALL MULTI-BAND PARAMETERS. FOR MORE THAN 2 BANDS THE
Groupie PROGRAM PERFORMS A NON-LINEAR FIT TO SELECT THE MULTI-BAND
Groupie PARAMETERS THAT MINIMIZE THE MAXIMUM FRACTIONAL ERROR AT ANY
Groupie POINT ALONG THE ENTIRE SELF-SHIELDING CURVE. THE NUMBER OF BANDS
Groupie REQUIRED WITHIN ANY GIVEN GROUP IS DEFINED BY INSURING THAT THE
Groupie MULTI-BAND PARAMETERS CAN BE USED TO ACCURATELY DEFINE SELF-
Groupie SHIELDED CROSS SECTIONS ALONG THE ENTIRE SELF-SHIELDING CURVE
Groupie FROM SIGMA0 = 0 TO INFINITY. THE USER MAY DEFINE THE ACCURACY
Groupie REQUIRED.

Groupie
Groupie ENDF/B FORMATTED UNSHIELDED AVERAGES
Groupie -----
Groupie UNSHIELDED MULTI-GROUP AVERAGED CROSS SECTIONS FOR ALL REACTIONS
Groupie MAY BE OBTAINED IN THE ENDF/B FORTRAN IN EITHER HISTOGRAM
Groupie (INTERPOLATION LAW 1) OR LINEARLY INTERPOLABLE (INTERPOLATION
Groupie LAW 2) FORM. SEE INPUT BELOW FOR DETAILS.

Groupie
Groupie MIXTURES OF MATERIALS AND RESONANCE OVERLAP
Groupie

Groupie
THE SELF-SHIELDED CROSS SECTIONS FOR THE INDIVIDUAL CONSTITUENTS
Groupie
OF ANY MIXTURE CAN BE CALCULATED BY THIS PROGRAM BY REALIZING THAT
Groupie
THIS PROGRAM ESSENTIALLY ONLY USES THE TOTAL CROSS SECTION AS A
Groupie
WEIGHTING FUNCTION TO ACCOUNT FOR SELF-SHIELDING EFFECTS. FOR A
Groupie
MIXTURE IT IS THEREFORE ONLY NECESSARY TO USE THE TOTAL CROSS
Groupie
SECTION FOR THE MIXTURE IN PLACE OF THE ACTUAL TOTAL CROSS SECTION
Groupie
FOR EACH CONSTITUENT AND TO RUN THIS PROGRAM. THIS CAN BE DONE BY
Groupie
FIRST RUNNING PROGRAM MIXER TO CALCULATE THE ENERGY DEPENDENT
Groupie
TOTAL CROSS SECTION FOR ANY COMPOSITE MIXTURE. NEXT, SUBSTITUTE
Groupie
THIS COMPOSITE TOTAL CROSS SECTION FOR THE ACTUAL TOTAL CROSS
Groupie
SECTION OF EACH CONSTITUENT (IN EACH ENDF/B FORMATTED EVALUATION).
Groupie
FINALLY, RUN THIS PROGRAM TO CALCULATE THE SELF-SHIELDED CROSS
Groupie
SECTION FOR EACH CONSTITUENT, PROPERLY ACCOUNTING FOR RESONANCE
Groupie
OVERLAP BETWEEN THE RESONANCES OF ALL OF THE CONSTITUENTS OF THE
Groupie
MIXTURE. DURING THE SAME RUN THESE SELF-SHIELDED CROSS SECTIONS
Groupie
CAN IN TURN BE USED TO CALCULATE FULLY CORRELATED MULT-BAND
Groupie

Groupie
MULTI-BAND PARAMETER OUTPUT FORMAT

Groupie

Groupie
FOR VERSIONS 92-2 AND LATER VERSIONS THE MULTI-BAND PARAMETERS
Groupie
ARE OUTPUT IN A SIMPLE CHARACTER FORMAT, THAT CAN BE TRANSFERRED
Groupie
AND USED ON VIRTUALLY ANY COMPUTER.
Groupie

Groupie
THE BINARY FORMAT USED IN EARLIER VERSIONS OF THIS CODE IS NO
Groupie
LONGER USED.
Groupie

Groupie

CONTACT THE AUTHOR IF YOU WOULD LIKE TO RECEIVE A SIMPLE PROGRAM
 Groupie
 TO READ THE CHARACTER FORMATTED MULTI-BAND PARAMETER FILE AND
 Groupie
 CREATE A BINARY, RANDOM ACCESS FILE FOR USE ON VIRTUALLY ANY
 Groupie
 COMPUTER.

Groupie
 THE FORMAT OF THE CHARACTER FILE IS,
 Groupie

RECORD	COLUMNS	FORMAT	DESCRIPTION
1	1-72	18A4	LIBRARY DESCRIPTION (AS READ)
2	1-11	I11	MATERIAL ZA
	12-22	I11	NUMBER GROUPS
	23-33	I11	NUMBER OF BANDS
	34-44	D11.4	TEMPERATURE (KELVIN)
	45-55	1X,10A1	HOLLERITH DESCRIPTION OF ZA
3	1-11	D11.4	ENERGY (EV) - GROUP BOUNDARY.
	12-22	D11.4	TOTAL (FIRST BAND)
	23-33	D11.4	ELASTIC
	34-44	D11.4	CAPTURE
	35-55	D11.4	FISSION
4	1-11	-----	BLANK
	12-22	D11.4	TOTAL (SECOND BAND)
	23-33	D11.4	ELASTIC
	34-44	D11.4	CAPTURE
	35-55	D11.4	FISSION

Groupie
 LINES 3 AND 4 ARE REPEATED FOR EACH GROUP. THE LAST LINE FOR EACH
 Groupie
 MATERIAL (ZA) IS,
 Groupie

Groupie
N 1-11 D11.4 ENERGY (EV) - UPPER ENERGY LIMIT OF
Groupie
LAST GROUP.

Groupie

Groupie
FOR EXAMPLE, A 175 GROUP, 2 BAND FILE, FOR EACH MATERIAL WILL
Groupie
CONTAIN 352 LINES = 1 HEADER LINE, 175 * 2 LINES OF PARAMETERS,
Groupie
AND 1 FINAL LINE WITH THE UPPER ENERGY LIMIT
Groupie
OF THE LAST GROUP.

Groupie

Groupie
INPUT FILES

Groupie

Groupie

UNIT DESCRIPTION

Groupie

---- -

Groupie

2 INPUT DATA (BCD - 80 CHARACTERS/RECORD)

Groupie

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

Groupie

Groupie
OUTPUT FILES

Groupie

Groupie

UNIT DESCRIPTION

Groupie

---- -

Groupie

31 MULTI-BAND PARAMETERS CHARACTER FILE - OPTIONAL

Groupie

(BCD - 80 CHARACTERS/RECORD)

Groupie

32 SELF-SHIELDED CROSS SECTION LISTING - OPTIONAL

Groupie

(BCD - 120 CHARACTERS/RECORD)

Groupie

33 MULTI-BAND PARAMETER LISTING - OPTIONAL

Groupie

(BCD - 120 CHARACTERS/RECORD)

Groupie

34 UNSHIELDED CROSS SECTION LISTING - OPTION

Groupie

```

                (BCD - 120 CHARACTERS/RECORD)
Groupie
  3  OUTPUT REPORT (BCD - 80 CHARACTERS/RECORD)
Groupie
 11  MULTI-GROUP ENDF/B DATA - OPTIONAL
Groupie
        (BCD - 80 CHARACTERS/RECORD)
Groupie

Groupie
      SCRATCH FILES
Groupie
-----
Groupie
      UNIT  FILENAME  DESCRIPTION
Groupie
-----  -----  -----
Groupie
   8  ENERGY DEPENDENT WEIGHTING SPECTRUM
Groupie
        (BINARY - 40080 WORDS/BLOCK)
Groupie
   9  TOTAL CROSS SECTION
Groupie
        (BINARY - 40080 WORDS/BLOCK)
Groupie
  12  ELASTIC CROSS SECTION - ONLY FOR SELF-SHIELDING CALCULATION
Groupie
        (BINARY - 40080 WORDS/BLOCK)
Groupie
  13  CAPTURE CROSS SECTION - ONLY FOR SELF-SHIELDING CALCULATION
Groupie
        (BINARY - 40080 WORDS/BLOCK)
Groupie
  14  FISSION CROSS SECTION - ONLY FOR SELF-SHIELDING CALCULATION
Groupie
        (BINARY - 40080 WORDS/BLOCK)
Groupie

Groupie
      OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINES FILIO1 AND FILIO2)
Groupie
-----
Groupie
      UNIT  FILE NAME
Groupie
-----  -----
Groupie
   2  GROUPIE.INP
Groupie
   3  GROUPIE.LST
Groupie
   8  (SCRATCH)
Groupie

```

Groupie 9 (SCRATCH)
 Groupie 10 ENDFB.IN
 Groupie 11 ENDFB.OUT
 Groupie 12 (SCRATCH)
 Groupie 13 (SCRATCH)
 Groupie 14 (SCRATCH)
 Groupie 31 MULTBAND.TAB
 Groupie 32 SHIELD.LST
 Groupie 33 MULTBAND.LST
 Groupie 34 UNSHIELD.LST
 Groupie

Groupie I/O UNITS USED
 Groupie

Groupie -----

Groupie UNITS 2, 3 8, 9 AND 10 WILL ALWAYS BE USED.

Groupie UNITS 31 THROUGH 34 AND 11 ARE OPTIONALLY USED DEPENDING ON THE
 Groupie OUTPUT REQUESTED.

Groupie UNITS 12, 13 AND 14 WILL ONLY BE USED IF SELF-SHIELDED OR
 Groupie MULTIBAND OUTPUT IS REQUESTED.

Groupie

Groupie INPUT CARDS
 Groupie

Groupie -----

CARD	COLS.	FORMAT	DESCRIPTION
------	-------	--------	-------------

----	-----	-----	-----
------	-------	-------	-------

1	1-11	I11	SELECTION CRITERIA (0=MAT, 1=ZA)
---	------	-----	----------------------------------

1	12-22	I11	NUMBER OF GROUPS.
---	-------	-----	-------------------

Groupie =.GT.0 - ARBITRARY GROUP BOUNDARIES ARE READ

Groupie FROM INPUT FILE (N GROUPS REQUIRE

Groupie

N+1 GROUP BOUNDARIES). CURRENT
 PROGRAM MAXIMUM IS 1000 GROUPS.
 BUILT-IN OPTIONS INCLUDE....

Groupie	= 0	- TART	175 GROUPS
Groupie	= -1	- ORNL	50 GROUPS
Groupie	= -2	- ORNL	126 GROUPS
Groupie	= -3	- ORNL	171 GROUPS
Groupie	= -4	- SAND-II	620 (665) GROUPS TO 18 MEV
Groupie	= -5	- SAND-II	640 (685) GROUPS TO 20 MEV
Groupie	= -6	- WIMS	69 GROUPS
Groupie	= -7	- GAM-I	68 GROUPS
Groupie	= -8	- GAM-II	99 GROUPS
Groupie	= -9	- MUFT	54 GROUPS
Groupie	== -10	- ABBN	28 GROUPS
Groupie	== -11	- TART	616 GROUPS TO 20 MEV
Groupie	== -12	- TART	700 GROUPS TO 1 GEV
Groupie	== -13	- SAND-II	665 GROUPS TO 18 MEV
Groupie	== -14	- SAND-II	685 GROUPS TO 20 MEV
Groupie	== -15	- TART	666 GROUPS TO 200 MEV
Groupie	== -16	- SAND-II	725 GROUPS TO 60 MEV
Groupie	== -17	- SAND-II	755 GROUPS TO 150 MEV
Groupie	== -18	- SAND-II	765 GROUPS TO 200 MEV

Groupie 1 23-33 I11 MULTI-BAND SELECTOR

Groupie	= 0	- NO MULTI-BAND CALCULATIONS
Groupie	= 1	- 2 BAND. CONSERVE AV(TOT), AV(1/TOT) AND AV(1/TOT**2)
Groupie	= 2	- 2 BAND. CONSERVE AV(TOT), AV(1/TOT)

AND $AV(1/(TOT+SIGMA0))$ WHERE
 SIGMA0 = AV(TOT) IN EACH GROUP
 = 3-5- MULTI-BAND FIT. CONSERVE AV(TOT) AND
 MINIMIZE FRACTIONAL ERROR FOR ENTIRE
 SELF-SHIELDING CURVE (SIGMA0 = 0 TO
 INFINITY)
 IF THE SELECTOR IS POSITIVE (1 TO 5) THE
 MINIMUM NUMBER OF BANDS WILL BE OUTPUT FOR
 EACH ISOTOPE INDEPENDENTLY. IF THE SELECTOR
 IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF
 BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR
 ALL ISOTOPES.
 1 34-44 I11 NUMBER OF POINTS USED TO DESCRIBE ENERGY
 DEPENDENT WEIGHTING SPECTRUM S(E).
 = -2 - MAXWELLIAN - UP TO 0.1 EV
 1/E - 0.1 EV TO 67 KEV
 FISSION - ABOVE 67 KEV
 05/01/20-----ADDED OPTION TO ALLOW TEMPERATURE OF THE
 MAXWELLIAN TO BE CHANGED - SEE INPUT LINE 4,
 COLUMNS 55 - 66.
 = -1 - 1/E
 = 0 OR 1- ENERGY INDEPENDENT (SO CALLED FLAT
 WEIGHTING SPECTRUM).
 = .GT.1 - READ THIS MANY POINTS FROM INPUT
 TO DESCRIBE WEIGHTING SPECTRUM.
 NO LIMIT TO THE NUMBER OF POINTS
 USED TO DESCRIBE WEIGHTING.

1 45-55 D11.4 MULTI-BAND CONVERGENCE CRITERIA.
 Groupie ONLY USED FOR 3 OR MORE BANDS. THE NUMBER OF
 Groupie BANDS IN EACH GROUPS IS SELECTED TO INSURE
 Groupie THAT THE ENTIRE SELF-SHIELDING CURVE CAN BE
 Groupie REPRODUCED TO WITHIN THIS FRACTIONAL ERROR.
 Groupie = .LT. 0.0001 - USE STANDARD 0.001
 Groupie (0.1 PER-CENT)
 Groupie = .GE. 0.0001 - USE AS CONVERGENCE CRITERIA
 1 56-66 I11 SIGMA-0 DEFINITION SELECTOR.
 Groupie < 0 - 21 VALUES OF SIGMA0 ARE READ INPUT AND
 Groupie INTERPRETED AS FIXED VALUES = SAME AS
 Groupie = 1 DESCRIPTION BELOW
 Groupie INPUT VALUES MUST ALL BE,
 Groupie 1) GREATER THAN 0
 Groupie 2) IN DESCENDING VALUE ORDER
 Groupie = 0 - SIGMA-0 WILL BE DEFINED AS A MULTIPLE
 Groupie OF THE UNSHIELDED TOTAL CROSS SECTION
 Groupie IN EACH GROUP (VALUES OF 1/1024 TO
 Groupie 1024 IN STEPS OF A FACTOR OF 2 WILL
 Groupie BE USED AS THE MULTIPLIER).
 Groupie = 1 - SIGMA-0 WILL BE DEFINED AS THE SAME
 Groupie NUMBER OF BARNS IN EACH GROUP (VALUES
 Groupie 40000 TO 0.4 BARNS WILL BE USED. WITHIN
 Groupie EACH DECADE VALUES OF 10, 7, 4, 2, 1
 Groupie BARNS WILL BE USED).
 1 67-70 I4 High energy extension = definition of cross
 Groupie section above highest tabulated energy.
 Groupie

Groupie = 0 = cross section = 0 (standard ENDF/B)
 Groupie = 1 = cross section = constant (equal to
 value at highest tabulated energy).
 Groupie 2-4 1-66 6D11.4 IF SIGMA-0 DEFINITION SELECTOR < 0, THE NEXT
 Groupie 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0,
 Groupie 6 PER LINE.
 Groupie 2 1-72 A72 ENDF/B INPUT DATA FILENAME
 Groupie (STANDARD OPTION = ENDFB.IN)
 Groupie 3 1-72 A72 ENDF/B OUTPUT DATA FILENAME
 Groupie (STANDARD OPTION = ENDFB.OUT)
 Groupie
 Groupie THE FOURTH INPUT CARD IS USED TO SELECT ALL DESIRED OUTPUT MODES.
 Groupie EACH OUTPUT DEVICE MAY BE TURNED OFF (0) OR ON (1). THEREFORE
 Groupie THEREFORE EACH OF THE FOLLOWING INPUT PARAMETERS MAY BE EITHER
 Groupie ZERO TO INDICATE NO OUTPUT OR NON-ZERO TO INDICATE OUTPUT.
 Groupie
 Groupie 4 1-11 I11 SELF-SHIELDED CROSS SECTION LISTING
 Groupie = 1 - CROSS SECTIONS
 Groupie = 2 - RESONANCE INTEGRALS
 Groupie 4 12-22 I11 MULTI-BAND PARAMETER LISTING
 Groupie 4 23-33 I11 MULTI-BAND PARAMETERS COMPUTER READABLE
 Groupie 4 34-44 I11 UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT
 Groupie = 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1)
 Groupie = 2 - LINEAR-LINEAR (INTERPOLATION LAW 2)
 Groupie 4 45-55 I11 UNSHIELDED CROSS SECTIONS LISTING
 Groupie = 1 - CROSS SECTIONS
 Groupie = 2 - RESONANCE INTEGRALS
 Groupie

05/01/20 - ADDED THE BELOW OPTION

Groupie
4 56-66 D11.4 IF THE STANDARD BUILT-IN SPECTRA IS USED,
Groupie
Groupie INPUT LINE 1, COLUMNS 34-44 = 2, THIS FIELD
Groupie CAN BE USED TO OPTIONALLY CHANGE TEMPERATURE
Groupie OF THE MAXWELLIAN.
Groupie INPUT IS IN EV (0.0253 EV = ROOM TEMPERATURE)
Groupie = 0 - USE DEFAULT 0.0253 EV, ROOM TEMPERATURE
Groupie > 0 - USE THIS AS THE TEMPERATURE
Groupie
Groupie RESTRICTION - TEMPERATURE CANNOT EXCEED
Groupie 1000 EV.

Groupie
5 1-80 18A4 LIBRARY IDENTIFICATION. ANY TEXT THAT THE
Groupie
Groupie USER WISHES TO IDENTIFY THE MULTI-BAND
Groupie PARAMETERS. THIS LIBRARY IDENTIFICATION IS
Groupie WRITTEN INTO THE COMPUTER READABLE MULTI-BAND
Groupie DATA FILE.

Groupie
6-N 1- 6 I6 LOWER MAT OR ZA LIMIT
Groupie
Groupie 7- 8 I2 LOWER MF LIMIT
Groupie
Groupie 9-11 I3 LOWER MT LIMIT
Groupie
Groupie 12-17 I11 UPPER MAT OR ZA LIMIT
Groupie
Groupie 18-19 I2 UPPER MF LIMIT
Groupie
Groupie 20-22 I3 UPPER MT LIMIT
Groupie
Groupie UP TO 100 RANGES MAY BE SPECIFIED, ONE RANGE
Groupie
Groupie PER LINE. THE LIST OF RANGES IS TERMINATED
Groupie
Groupie BY A BLANK CARD. IF THE UPPER MAT OR ZA
Groupie
Groupie LIMIT IS LESS THAN THE LOWER LIMIT THE UPPER

Groupie
Groupie
Groupie
Groupie
Groupie
Groupie
Groupie

IS SET EQUAL TO THE LOWER LIMIT. IF THE UPPER MF OR MT LIMIT IS ZERO IT WILL BE SET EQUAL TO ITS MAXIMUM VALUE, 99 OR 999, RESPECTIVELY IF THE FIRST REQUEST LINE IS BLANK IT WILL TERMINATE THE LIST OF REQUESTS AND CAUSE ALL DATA TO BE RETRIEVED (SEE EXAMPLE INPUT).

Groupie
VARY 1-66 6D11.4
Groupie
Groupie

ENERGY GROUP BOUNDARIES. ONLY REQUIRED IF THE NUMBER OF GROUPS INDICATED ON THE FIRST INPUT CARD IS POSITIVE. ALL ENERGIES MUST BE IN ASCENDING ENERGY IN EV. THE PRESENT LIMITS ARE 1 TO 1000 GROUPS. FOR N GROUPS N+1 BOUNDARIES WILL BE READ FROM THE INPUT FILE, E.G. IF THE FIRST INPUT CARD INDICATES 20 GROUPS, 21 ENERGY BOUNDARIES WILL BE READ FROM THE INPUT FILE.

Groupie
VARY 1-66 6D11.4
Groupie
Groupie

ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY REQUIRED IF THE NUMBER OF POINTS INDICATED ON FIRST CARD IS MORE THAN ONE. DATA IS GIVEN IN (ENERGY, WEIGHT) PAIRS, UP TO 3 PAIRS PER CARD, USING ANY NUMBER OF CARDS REQUIRED. ENERGIES MUST BE IN ASCENDING ORDER IN EV. THE SPECTRUM VALUES MUST BE NON-NEGATIVE. THE ENERGY RANGE OF SPECTRUM MUST AT LEAST SPAN THE ENERGY RANGE OF THE ENERGY GROUPS. SINCE SPECTRUM IS STORED IN

Groupie PAGING SYSTEM THERE IS NO LIMIT TO NUMBER
Groupie OF POINTS THAT CAN BE USED TO DESCRIBE THE
Groupie WEIGHTING SPECTRUM.

Groupie
Groupie EXAMPLE INPUT NO. 1

Groupie -----

Groupie REQUEST DATA BY MAT AND PROCESS ALL DATA (ALL MAT BETWEEN 1 AND
Groupie 9999). USE THE TART 175 GROUP STRUCTURE, GENERATE 2 BAND
Groupie PARAMETERS (THE FOR ALL ISOTOPES) TO 0.1 PER-CENT ACCURACY
Groupie IN THE SELF-SHIELDING CURVE. OUTPUT ALL LISTING, COMPUTER
Groupie READABLE AND ENDF/B FORMAT GROUP AVERAGES.

Groupie
Groupie EXPLICITLY SPECIFY THE STANDARD FILENAMES.

Groupie
Groupie THE FOLLOWING 7 INPUT LINES ARE REQUIRED.

Groupie
Groupie 0 0 -2 0 1.00000-03 0

Groupie ENDFB.IN

Groupie ENDFB.OUT

Groupie 1 1 1 1 1 1

Groupie TART 175 GROUP, 2 BAND LIBRARY TO 0.1 PER-CENT ACCURACY

Groupie 1 1 1 9999 0 0

Groupie (BLANK CARD TERMINATES REQUEST LIST)

Groupie
Groupie EXAMPLE INPUT NO. 2

Groupie -----

Groupie
Groupie THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ

(USE STANDARD FILENAME = ENDFB.OUT)

Groupie

0 0 0 0 1

Groupie

RESONANCE INTEGRAL CALCULATION (FROM 0.5 EV TO 1 MEV)

Groupie

(RETRIEVE ALL DATA, TERMINATE REQUEST LIST)

Groupie

5.00000-01 1.00000+06

Groupie

Groupie

=====

Groupie