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PROGRAM GROUPIE
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VERSION 76-1 (NOVEMBER 1976)
VERSION 79-1 (OCTOBER 1979) CDC-7600 AND CRAY-1 VERSION.
VERSION 80-1 (MAY 1980) IBM, CDC AND CRAY VERSION
VERSION 81-1 (JANUARY 1981) EXTENSION TO 3000 GROUPS
VERSION 81-2 (MARCH 1981) IMPROVED SPEED
VERSION 81-3 (AUGUST 1981) BUILT-IN 1/E WEIGHTING SPECTRUM
VERSION 82-1 (JANUARY 1982) IMPROVED COMPUTER COMPATIBILITY
VERSION 83-1 (JANUARY 1983) *MAJOR RE-DESIGN.
*ELIMINATED COMPUTER DEPENDENT CODING.
*NEW, MORE COMPATIBLE I/O UNIT NUMBERS.
*NEW MULTI-BAND LIBRARY BINARY FORMAT.
VERSION 83-2 (OCTOBER 1983) ADDED OPTION TO ALLOW SIGMA-0 TO BE
DEFINED EITHER AS MULTIPLES OF
UNSHIELDED TOTAL CROSS SECTION IN EACH
GROUP, OR POWERS OF 10 IN ALL GROUPS.
VERSION 84-1 (APRIL 1984) ADDED MORE BUILT IN MULTIGROUP ENERGY
STRUCTURES.
VERSION 85-1 (APRIL 1985) *UPDATED FOR ENDF/B-VI FORMATS.
*SPECIAL I/O ROUTINES TO GUARANTEE
ACCURACY OF ENERGY.
*DOUBLE PRECISION TREATMENT OF ENERGY
(REQUIRED FOR NARROW RESONANCES).
*MINIMUM TOTAL CROSS SECTION TREATMENT
VERSION 85-2 (AUGUST 1985) *FORTRAN-77/H VERSION
VERSION 86-1 (JANUARY 1986) *ENDF/B-VI FORMAT
VERSION 86-2 (JUNE 1986) *BUILT-IN MAXWELLIAN, 1/E AND FISSION
WEIGHTING SPECTRUM.
VERSION 88-1 (JULY 1988) *OPTION...INTERNALLY DEFINE ALL I/O
FILE NAMES (SEE, SUBROUTINES FILIO1
FILIO2 FOR DETAILS).
*IMPROVED BASED ON USER COMMENTS.
VERSION 89-1 (JANUARY 1989) *PSYCHOANALYZED BY PROGRAM FREUD TO
INSURE PROGRAM WILL NOT DO ANYTHING
CRAZY.
*UPDATED TO USE NEW PROGRAM CONVERT
KEYWORDS.
*ADDED LIVERMORE CIVIC COMPILER
CONVENTIONS.
VERSION 91-1 (JUNE 1991) *INCREASED PAGE SIZE FROM 1002 TO 5010
POINTS
*UPDATED BASED ON USER COMMENTS
*ADDED FORTRAN SAVE OPTION
*COMPLETELY CONSISTENT ROUTINE TO READ
FLOATING POINT NUMBERS.
VERSION 92-1 (JANUARY 1992) *ADDED RESONANCE INTEGRAL CALCULATION -
UNSHIELDED AND/OR SHIELDED - FOR
DETAILS SEE BELOW
*INCREASED NUMBER OF ENERGY POINTS
IN BUILT-IN SPECTRA - TO IMPROVE
ACCURACY.
*ALLOW SELECTION OF ZA/MF/MT OR
MAT/MF/MT RANGES - ALL DATA NOT
SELECTED IS SKIPPED ON INPUT AND
NOT WRITTEN AS OUTPUT.
*COMPLETELY CONSISTENT I/O ROUTINES -
TO MINIMIZE COMPUTER DEPENDENCE.
*NOTE, CHANGES IN INPUT PARAMETER
FORMAT - FOR ZA/MF/MT OR MAT/MF/MT
RANGES.
VERSION 92-2 (JUNE 1992) *MULTIBAND PARAMETERS OUTOUT AS
CHARACTER (RATHER THAN BINARY) FILE.
VERSION 93-1 (APRIL 1993) *INCREASED PAGE SIZE FROM 5010 TO
30000 POINTS
*ELIMINATED COMPUTER DEPENDENCE.
VERSION 94-1 (JANUARY 1994) *VARIABLE ENDF/B DATA FILENAMES
TO ALLOW ACCESS TO FILE STRUCTURES
(WARNING - INPUT PARAMETER FORMAT

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	HAS BEEN CHANGED)	GROUPIE
	*CLOSE ALL FILES BEFORE TERMINATING (SEE, SUBROUTINE ENDIT)	GROUPIE
VERSION 95-1 (JANUARY 1994)	*CORRECTED MAXWELLIAN WEIGHTING	GROUPIE
	*CHANGING WEIGHTING SPECTRUM FROM 0.1 TO 0.001 % UNCERTAINTY	GROUPIE
VERSION 96-1 (JANUARY 1996)	*COMPLETE RE-WRITE	GROUPIE
	*IMPROVED COMPUTER INDEPENDENCE	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 CALCULATION (PREVIOUSLY 175)	GROUPIE
	*MAXIMUM NUMBER OF GROUPS REDUCED FROM 3,000 TO 1,000	GROUPIE
	*UP TO 1000 MATERIALS (PREVIOUSLY 100)	GROUPIE
	*CORRECTED USE OF MAXWELLIAN + 1/E + FISSION SPECTRUM	GROUPIE
	*ONLY 2 BAND VERSION DISTRIBUTED (CONTACT AUTHOR FOR DETAILS)	GROUPIE
	*DEFINED SCRATCH FILE NAMES	GROUPIE
VERSION 99-1 (MARCH 1999)	*CORRECTED CHARACTER TO FLOATING POINT READ FOR MORE DIGITS	GROUPIE
	*UPDATED TEST FOR ENDF/B FORMAT	GROUPIE
	VERSION BASED ON RECENT FORMAT CHANGE	GROUPIE
	*GENERAL IMPROVEMENTS BASED ON USER FEEDBACK	GROUPIE
VERSION 99-2 (JUNE 1999)	*ASSUME ENDF/B-VI, NOT V, IF MISSING MF=1, MT-451.	GROUPIE
VERS. 2000-1 (FEBRUARY 2000)	*ADDED MF=10, ACTIVATION CROSS SECTION PROCESSING.	GROUPIE
	*GENERAL IMPROVEMENTS BASED ON USER FEEDBACK	GROUPIE
VERS. 2002-1 (FEBRUARY 2002)	*ADDED TART 700 GROUP STRUCTURE	GROUPIE
	*ADDED VARIABLE SIGMA0 INPUT OPTION	GROUPIE
(MAY 2002)	*OPTIONAL INPUT PARAMETERS	GROUPIE
(NOV. 2002)	*ADDED SAND-II EXTENDED DOWN TO 1.0D-5 EV.	GROUPIE
(JUNE 2003)	*CORRECTED SAND-II 620 AND 640 GROUP ENERGY BOUNDARIES DEFINITIONS.	GROUPIE
VERS. 2004-1 (SEPT. 2004)	*INCREASED PAGE SIZE FROM 30000 TO 120000 POINTS	GROUPIE
	*ADDED "OTHER" AS ADDITIONAL REACTION TO IMPROVE MULTI-BAND FITTING	GROUPIE
	*ADDED ITERATION FOR "BEST" PARTIAL PARAMETERS.	GROUPIE
	*DO NOT SKIP LOW TOTAL ENERGY RANGES WHEN DEFINING AVERAGE CROSS SECTIONS	GROUPIE
	THIS MAKES OUTPUT COMPATIBLE WITH ANY STANDARD AVERAGING PROCEDURE	GROUPIE
VERS. 2005-1 (JAN. 2005)	*ADDED OPTION TO CHANGE TEMPERATURE OF BUILT-IN STANDARD SPECTRUM.	GROUPIE
VERS. 2007-1 (JAN. 2007)	*CHECKED AGAINST ALL ENDF/B-VII.	GROUPIE
	*INCREASED PAGE SIZE FROM 120,000 TO 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 FROM 3,000 ENERGY POINTS.	GROUPIE
	*ADDED OUTPUT TO PLOT/COMPARE SHIELDED AND UNSHIELDED CROSS SECTIONS.	GROUPIE
VERS. 2011-1 (June 2011)	*Corrected TART 700 groups to extend up to 1 GeV (1,000 MeV) - previously it was ERRONEOUSLY cutoff at 20 MeV.	GROUPIE
VERS. 2011-2 (Nov. 2011)	*Corrected TART 616 groups lowest energy from 1.0D-4 eV to 1.0D-5 eV.	GROUPIE
	*Added TART 666 to 200 MeV (for TENDL).	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
*32 and 64 bit Compatible GROUPIE
*Added ERROR stop. GROUPIE

VERS. 2013-1 (Nov. 2013) *Extended OUT9. GROUPIE
*Uses OUTG, not OUT10 for energies. GROUPIE

VERS. 2015-1 (Jan. 2015) *Corrected SPECTM - handle ALL included GROUPIE
group structures, i.e., even those GROUPIE
that start above thremal range by GROUPIE
ALWAYS constructing weigthning spectrum GROUPIE
to be AT LEAST 1.0D-5 eV to 20 MeV. GROUPIE
*Extended OUTG GROUPIE
*Replaced ALL 3 way IF Statements. 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

VERS. 2015-3 (July 2015) *Insure no 10 digit output - not GROUPIE
needed for multi-group and this makes GROUPIE
listings simpler. GROUPIE
*Corrected High Energy Extension = GROUPIE
Can effect highest energy group. GROUPIE

VERS. 2016-1 (July 2016) *Added UKAEA 1102 Group Structure. GROUPIE
*Increased storage to accommodate GROUPIE
much larger group structures = GROUPIE
up to 20,000 Groups. GROUPIE
*Added output listing of the complete GROUPIE
input parameters for URRFIT, including GROUPIE
the NJOY parameters LSSF and ICOMP. GROUPIE
*Changed multiple IF statements to GROUPIE
accommodate compiler optimizer GROUPIE
*Cosmetic changes based on FREUD GROUPIE
psychoanalysis. GROUPIE
*Updated multi-band treatment to GROUPIE
explitly handle small shielding GROUPIE
limit - without this update the small GROUPIE
limit becomes numerically unstable. GROUPIE

VERS. 2017-1 (May 2017) *Increased max. points to 3,000,000. GROUPIE
*METHODEB was incorrecctly named GROUPIE
METHOD in one routine = corrected. GROUPIE
*Default multi-band is method #2 = GROUPIE
conserve <x>, <1/(x+<x>>), <1/x>. GROUPIE
*Definition of built-in group structure GROUPIE
using SUBROUTINE GROPE is identical GROUPIE
for GROUPIE and VIRGIN. GROUPIE
*All floating input parameters changed GROUPIE
to character input + IN9 conversion. GROUPIE
*Output report identfies MF now that GROUPIE
this code does more than just MF=3. GROUPIE
*Added NRO = energy dependent scatter GROUPIE
radius to copying FILE2 parameters GROUPIE
to define unresolved energy range. GROUPIE
*Corrected energy dependent scatter GROUPIE
for all resonance types (see, above GROUPIE
comments) = for multi-band output GROUPIE

VERS. 2018-1 (Jan. 2018) *Added on-line output for ALL ENDERROR GROUPIE

VERS. 2019-1 (June 2019) *Major re-write to re-order output to GROUPIE
include Unresolved Resonance Region GROUPIE
self-shielding. GROUPIE
*Added Unresolved self-shielding by GROUPIE
Extrapolating cross section moments GROUPIE

IF THIS PROGRAM FINDS THAT THE FILE 3 CROSS SECTIONS ARE NOT LINEARLY INTERPOLABLE THIS PROGRAM WILL TERMINATE EXECUTION.	GROUPIE
	GROUPIE
	GROUPIE
CONTENTS OF OUTPUT	GROUPIE
-----	GROUPIE
IF ENDF/B FORMATTED OUTPUT IS REQUESTED ENTIRE EVALUATIONS ARE OUTPUT, NOT JUST THE MULTI-GROUPED FILE 3 CROSS SECTIONS, E.G. ANGULAR AND ENERGY DISTRIBUTIONS ARE ALSO INCLUDED.	GROUPIE
	GROUPIE
	GROUPIE
DOCUMENTATION	GROUPIE
-----	GROUPIE
THE FACT THAT THIS PROGRAM HAS OPERATED ON THE DATA IS DOCUMENTED BY THE ADDITION OF THREE COMMENT CARDS AT THE END OF EACH HOLLERITH SECTION TO DESCRIBE THE GROUP STRUCTURE AND WEIGHTING SPECTRUM, E.G.	GROUPIE
	GROUPIE
	GROUPIE
***** PROGRAM GROUPIE (2021-1) *****	GROUPIE
UNSHIELDED GROUP AVERAGES USING 69 GROUPS (WIMS)	GROUPIE
MAXWELLIAN, 1/E, FISSION TO CONSTANT WEIGHTING SPECTRUM	GROUPIE
	GROUPIE
THE ORDER OF ALL SIMILAR COMMENTS (FROM LINEAR, RECENT AND SIGMA1) REPRESENTS A COMPLETE HISTORY OF ALL OPERATIONS PERFORMED ON THE DATA.	GROUPIE
	GROUPIE
	GROUPIE
THESE COMMENT CARDS ARE ONLY ADDED TO EXISTING HOLLERITH SECTIONS, I.E., THIS PROGRAM WILL NOT CREATE A HOLLERITH SECTION. THE FORMAT OF THE HOLLERITH SECTION IN ENDF/B-V DIFFERS FROM THE THAT OF EARLIER VERSIONS OF ENDF/B. BY READING AN EXISTING MF=1, MT=451 IT IS POSSIBLE FOR THIS PROGRAM TO DETERMINE WHICH VERSION OF THE ENDF/B FORMAT THE DATA IS IN. WITHOUT HAVING A SECTION OF MF=1, MT=451 PRESENT IT IS IMPOSSIBLE FOR THIS PROGRAM TO DETERMINE WHICH VERSION OF THE ENDF/B FORMAT THE DATA IS IN, AND AS SUCH IT IS IMPOSSIBLE FOR THE PROGRAM TO DETERMINE WHAT FORMAT SHOULD BE USED TO CREATE A HOLLERITH SECTION.	GROUPIE
	GROUPIE
	GROUPIE
REACTION INDEX	GROUPIE
-----	GROUPIE
THIS PROGRAM DOES NOT USE THE REACTION INDEX WHICH IS GIVEN IN SECTION MF=1, MT=451 OF EACH EVALUATION.	GROUPIE
	GROUPIE
	GROUPIE
THIS PROGRAM DOES NOT UPDATE THE REACTION INDEX IN MF=1, MT=451. THIS CONVENTION HAS BEEN ADOPTED BECAUSE MOST USERS DO NOT REQUIRE A CORRECT REACTION INDEX FOR THEIR APPLICATIONS AND IT WAS NOT CONSIDERED WORTHWHILE TO INCLUDE THE OVERHEAD OF CONSTRUCTING A CORRECT REACTION INDEX IN THIS PROGRAM. HOWEVER, IF YOU REQUIRE A REACTION INDEX FOR YOUR APPLICATIONS, AFTER RUNNING THIS PROGRAM YOU MAY USE PROGRAM DICTIN TO CREATE A CORRECT REACTION INDEX.	GROUPIE
	GROUPIE
	GROUPIE
SECTION SIZE	GROUPIE
-----	GROUPIE
SINCE THIS PROGRAM USES A LOGICAL PAGING SYSTEM THERE IS NO LIMIT TO THE NUMBER OF POINTS IN ANY SECTION, E.G., THE TOTAL CROSS SECTION MAY BE REPRESENTED BY 200,000 DATA POINTS.	GROUPIE
	GROUPIE
	GROUPIE
SELECTION OF DATA	GROUPIE
-----	GROUPIE
THE PROGRAM SELECTS MATERIALS TO BE PROCESSED BASED EITHER ON MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS USED TO SELECT MATERIALS, AND WILL TERMINATE WHEN A MAT OR ZA IS FOUND THAT IS ABOVE THE RANGE OF ALL REQUESTS.	GROUPIE
	GROUPIE
	GROUPIE
ENERGY ORDER AND UNITS	GROUPIE
-----	GROUPIE
ALL ENERGIES (FOR CROSS SECTIONS, WEIGHTING SPECTRUM OR GROUP BOUNDARIES) MUST BE IN UNITS OF EV AND MUST BE IN ASCENDING NUMERICAL ORDER.	GROUPIE
	GROUPIE
	GROUPIE
ENERGY GRID	GROUPIE
-----	GROUPIE

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

GROUP STRUCTURE

THIS PROGRAM IS DESIGNED TO USE AN ARBITRARY ENERGY GROUP STRUCTURE WHERE THE ENERGIES ARE IN EV AND ARE IN INCREASING ENERGY ORDER. THE MAXIMUM NUMBER OF GROUPS IS 20,000.

THE USER MAY INPUT AN ARBITRARY GROUP STRUCTURE OR THE USER MAY USE ONE OF THE BUILT-IN GROUP STRUCTURES.

- (0) 175 GROUP (TART STRUCTURE)
- (1) 50 GROUP (ORNL STRUCTURE)
- (2) 126 GROUP (ORNL STRUCTURE)
- (3) 171 GROUP (ORNL STRUCTURE)
- (4) 620 GROUP (SAND-II STRUCTURE, UP TO 18 MEV)
- (5) 640 GROUP (SAND-II STRUCTURE, UP TO 20 MEV)
- (6) 69 GROUP (WIMS STRUCTURE)
- (7) 68 GROUP (GAM-I STRUCTURE)
- (8) 99 GROUP (GAM-II STRUCTURE)
- (9) 54 GROUP (MUFT STRUCTURE)
- (10) 28 GROUP (ABBN STRUCTURE)
- (11) 616 GROUP (TART STRUCTURE TO 20 MeV)
- (12) 700 GROUP (TART STRUCTURE TO 1 GEV)
- (13) 665 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 18 MEV)
- (14) 685 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 20 MEV)
- (15) 666 GROUP (TART STRUCTURE TO 200 MeV)
- (16) 725 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 60 MEV)
- (17) 755 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 150 MEV)
- (18) 765 GROUP (SAND-II STRUCTURE, 1.0D-5 eV, UP TO 200 MEV)
- (19) 1102 GROUP (UKAEA STRUCTURE, 1.0D-5 eV, UP TO 1 GeV)

GROUP AVERAGES

THIS PROGRAM DEFINES GROUP AVERAGED CROSS SECTIONS AS...

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

WHERE...

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

ENERGY DEPENDENT WEIGHTING SPECTRUM

THE ENERGY DEPENDENT WEIGHTING SPECTRUM IS GIVEN BY AN ARBITRARY TABULATED LINEARLY INTERPOLABLE FUNCTION WHICH CAN BE DESCRIBED BY AN ARBITRARY NUMBER OF POINTS. THIS ALLOWS THE USER TO SPECIFY ANY DESIRED WEIGHTING SPECTRUM TO ANY GIVEN DEGREE OF ACCURACY. REMEMBER THAT THE PROGRAM WILL ASSUME THAT THE SPECTRUM IS LINEARLY INTERPOLABLE BETWEEN TABULATED POINTS. THEREFORE THE USER SHOULD USE ENOUGH POINTS TO INSURE AN ADEQUATE REPRESENTATION OF THE SPECTRUM BETWEEN TABULATED DATA POINTS.

THE PRESENT VERSION OF THE CODE HAS THREE BUILT-IN WEIGHTING SPECTRA,

- (1) CONSTANT
- (2) 1/E
- (3) MAXWELLIAN = $E * \exp(-E/KT) / KT$ (0.0 TO 4*KT)
- 1/E = $C1/E$ (4*KT TO 67 KeV)
- FISSION = $C2 * \exp(-E/WA) * \sinh(\text{SQRT}(E * WB))$ (67 KeV, 10 MeV)
- CONSTANT = Equal to Fission at 10 MeV (above 10 MeV)

RECORD	COLUMNS	FORMAT	DESCRIPTION	GROUPIE
1	1-72	18A4	LIBRARY DESCRIPTION (AS READ)	GROUPIE
2	1-11	I11	MATERIAL ZA	GROUPIE
	12-22	I11	NUMBER GROUPS	GROUPIE
	23-33	I11	NUMBER OF BANDS	GROUPIE
	34-44	E11.4	TEMPERATURE (KELVIN)	GROUPIE
	45-57	1X,12A1	HOLLERITH DESCRIPTION OF ZA	GROUPIE
3	1-11	E11.4	ENERGY (EV) - GROUP BOUNDARY.	GROUPIE
	12-22	E11.4	TOTAL (FIRST BAND)	GROUPIE
	23-33	E11.4	ELASTIC	GROUPIE
	34-44	E11.4	CAPTURE	GROUPIE
	35-55	E11.4	FISSION	GROUPIE
4	1-11	----	BLANK	GROUPIE
	12-22	E11.4	TOTAL (SECOND BAND)	GROUPIE
	23-33	E11.4	ELASTIC	GROUPIE
	34-44	E11.4	CAPTURE	GROUPIE
	35-55	E11.4	FISSION	GROUPIE

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

N	1-11	E11.4	ENERGY (EV) - UPPER ENERGY LIMIT OF LAST GROUP.	GROUPIE
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FOR EXAMPLE, A 175 GROUP, 2 BAND FILE, FOR EACH MATERIAL WILL CONTAIN 352 LINES = 1 HEADER LINE, 175 * 2 LINES OF PARAMETERS, AND 1 FINAL LINE WITH THE UPPER ENERGY LIMIT OF THE LAST GROUP.

INPUT FILES

UNIT	DESCRIPTION	GROUPIE
2	INPUT DATA (BCD - 80 CHARACTERS/RECORD)	GROUPIE
10	ORIGINAL ENDF/B DATA (BCD - 80 CHARACTERS/RECORD)	GROUPIE

OUTPUT FILES

UNIT	DESCRIPTION	GROUPIE
3	OUTPUT REPORT (BCD - 80 CHARACTERS/RECORD)	GROUPIE
11	MULTI-GROUP ENDF/B DATA - OPTIONAL (BCD - 80 CHARACTERS/RECORD)	GROUPIE
16	PLOTTAB FORMATTED SELF-SHIELDING RESULTS (BCD - 80 CHARACTERS/RECORD)	GROUPIE
31	MULTI-BAND PARAMETERS CHARACTER FILE - OPTIONAL (BCD - 80 CHARACTERS/RECORD)	GROUPIE
32	UNRESOLVED FSELF-SHIELDED PSEUDO ENDF FORMAT - OPTIONAL (BCD - 120 CHARACTERS/RECORD)	GROUPIE
33	SELF-SHIELDED CROSS SECTION LISTING - OPTIONAL (BCD - 120 CHARACTERS/RECORD)	GROUPIE
34	MULTI-BAND PARAMETER LISTING - OPTIONAL (BCD - 120 CHARACTERS/RECORD)	GROUPIE
35	UNSHIELDED CROSS SECTION LISTING - OPTION (BCD - 120 CHARACTERS/RECORD)	GROUPIE

SCRATCH FILES

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

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OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINES FILIO1 AND FILIO2)
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UNIT  FILE NAME
-----
      2  GROUPIE.INP
      3  GROUPIE.LST
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      8  (SCRATCH)
      9  (SCRATCH)
     10  ENDFB.IN
     11  ENDFB.OUT
     12  (SCRATCH)
     13  (SCRATCH)
     14  (SCRATCH)
-----2019/6/23 - New Filenames (added ZAzzaaa at Beginning)-----
----- (OLD)----- (NEW)-----
     16  PLOTTAB.CUR      ZAzzaaa.PLOT.CUR
     31  MULTBAND.TAB    ZAzzaaa.MULTBAND.TAB
     32  ZAzzaaa.URR.ENDF
     33  SHIELD.LST     ZAzzaaa.SHIELD.LST
     34  MULTBAND.LST   ZAzzaaa.MULTBAND.LST
     35  UNSHIELD.LST   ZAzzaaa.UNSHIELD.LST

I/O UNITS USED
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UNITS 2, 3 8, 9 AND 10 WILL ALWAYS BE USED.
UNITS 31 THROUGH 35, 11 AND 16 ARE OPTIONALLY USED DEPENDING
ON THE OUTPUT REQUESTED.
UNITS 12, 13 AND 14 WILL ONLY BE USED IF SELF-SHIELDED OR
MULTIBAND OUTPUT IS REQUESTED.

INPUT CARDS
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CARD  COLS.  FORMAT  DESCRIPTION
-----
      1   1-11   I11    SELECTION CRITERIA (0=MAT, 1=ZA)
      1   12-22  I11    NUMBER OF GROUPS.
-----2019/6/23 -11 (TART 616 groups) required for
Unresolved Resonance Region Self-Shielding
calculation.
=.GT.0 - ARBITRARY GROUP BOUNDARIES ARE READ
FROM INPUT FILE (N GROUPS REQUIRE
N+1 GROUP BOUNDARIES). CURRENT
PROGRAM MAXIMUM IS 20,000 GROUPS.
BUILT-IN OPTIONS INCLUDE....
= 0 - TART 175 GROUPS
= -1 - ORNL 50 GROUPS
= -2 - ORNL 126 GROUPS
= -3 - ORNL 171 GROUPS
= -4 - SAND-II 620 (665) GROUPS TO 18 MEV
= -5 - SAND-II 640 (685) GROUPS TO 20 MEV
= -6 - WIMS 69 GROUPS
= -7 - GAM-I 68 GROUPS
= -8 - GAM-II 99 GROUPS
= -9 - MUFT 54 GROUPS
=-10 - ABBN 28 GROUPS
Current TART Standard =-11 - TART 616 GROUPS TO 20 MEV
(-11 is required for =-12 - TART 700 GROUPS TO 1 GEV
unresolved resonance =-13 - SAND-II 665 GROUPS TO 18 MEV
region self-shielding=-14 - SAND-II 685 GROUPS TO 20 MEV
calculations) =-15 - TART 666 GROUPS TO 200 MEV
=-16 - SAND-II 725 GROUPS TO 60 MEV
=-17 - SAND-II 755 GROUPS TO 150 MEV
=-18 - SAND-II 765 GROUPS TO 200 MEV
=-19 - UKAEA 1102 GROUPS TO 1 GeV
      1   23-33   I11    MULTI-BAND SELECTOR
-----2019/6/23 - ONLY 0 or 2 allowed = the = 1
option has proven to give very poor results,
and therefore is no longer allowed.
= 0 - NO MULTI-BAND CALCULATIONS

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No longer allowed = 1 - 2 BAND. CONSERVE AV(TOT), AV(1/TOT)      GROUPIE
                   AND AV(1/TOT**2)                             GROUPIE
= 2 - 2 BAND. CONSERVE AV(TOT), AV(1/TOT)                     GROUPIE
                   AND AV(1/(TOT+SIGMA0)) WHERE                GROUPIE
                   SIGMA0 = AV(TOT) IN EACH GROUP              GROUPIE
No longer allowed = 3-5- MULTI-BAND FIT. CONSERVE AV(TOT) AND  GROUPIE
                   MINIMIZE FRACTIONAL ERROR FOR ENTIRE       GROUPIE
                   SELF-SHIELDING CURVE (SIGMA0 = 0 TO        GROUPIE
                   INFINITY)                                   GROUPIE
                   IF THE SELECTOR IS POSITIVE (1 TO 5) THE    GROUPIE
                   MINIMUM NUMBER OF BANDS WILL BE OUTPUT FOR GROUPIE
                   EACH ISOTOPE INDEPENDENTLY. IF THE SELECTOR GROUPIE
                   IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF   GROUPIE
                   BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR    GROUPIE
                   ALL ISOTOPES.                               GROUPIE
1   34-44   I11  NUMBER OF POINTS USED TO DESCRIBE ENERGY    GROUPIE
                   DEPENDENT WEIGHTING SPECTRUM S(E).         GROUPIE
= 0 or 1 - Flat (Constant)                                    GROUPIE
= -1      - 1/E at ALL energies                               GROUPIE
= -2      - MAXWELLIAN - UP TO 0.1 eV                        GROUPIE
                   1/E      - 0.1 eV TO 67 KeV              GROUPIE
                   FISSION  - 67 KeV to 10 MeV              GROUPIE
                   CONSTANT - Above 10 MeV                  GROUPIE
= > 1     - Read input table                                 GROUPIE
2005/01/20-----ADDED OPTION TO ALLOW TEMPERATURE OF THE    GROUPIE
                   MAXWELLIAN TO BE CHANGED - SEE INPUT LINE 4, GROUPIE
                   COLUMNS 55 - 66.                          GROUPIE
= -1      - 1/E                                             GROUPIE
= 0 OR 1- ENERGY INDEPENDENT (SO CALLED FLAT                GROUPIE
                   WEIGHTING SPECTRUM).                      GROUPIE
= .GT.1 - READ THIS MANY POINTS FROM INPUT                   GROUPIE
                   TO DESCRIBE WEIGHTING SPECTRUM.           GROUPIE
                   NO LIMIT TO THE NUMBER OF POINTS          GROUPIE
                   USED TO DESCRIBE WEIGHTING.               GROUPIE
1   45-55   E11.4 MULTI-BAND CONVERGENCE CRITERIA.           GROUPIE
-----2019/6/23 - No longer used now that code              GROUPIE
                   is restricted to no more than 2 bands.     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                  GROUPIE
1   56-66   I11  SIGMA-0 DEFINITION SELECTOR.                GROUPIE
-----2019/6/23 - For multi-band calculations                GROUPIE
                   only 0 is allowed = multiples of unshielded GROUPIE
                   total in each group = This is required for GROUPIE
                   the BEST self-shielding results.           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. WITHINGROUPIE
                   EACH DECADE VALUES OF 10, 7, 4, 2, 1    GROUPIE
                   BARNS WILL BE USED).                       GROUPIE
1   67-70   I14  High energy extension = definition of cross  GROUPIE
                   section above highest tabulated energy.   GROUPIE
-----2019/6/23 - Ignored - will always use ENDF           GROUPIE
                   Standard Definition = 0.                   GROUPIE
= 0 = cross section = 0 (standard ENDF/B)                   GROUPIE
= 1 = cross section = constant (equal to                     GROUPIE

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			value at highest tabulated energy).	GROUPIE
2-4	1-66	6E11.4	SIGMA-0 Definition	GROUPIE
-----			2019/6/23 - Only the GROUPIE standard = 0	GROUPIE
			in allowed for Unresolved Resonance Region	GROUPIE
			Self-Shielding calculation	GROUPIE
			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	E11.4	IF THE STANDARD BUILT-IN SPECTRA IS USED,	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
			RESTRICTION - TEMPERATURE CANNOT EXCEED	GROUPIE
			1000 EV.	GROUPIE
				GROUPIE
5	1-80	18A4	LIBRARY IDENTIFICATION. ANY TEXT THAT THE	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
				GROUPIE
6-N	1- 6	I6	LOWER MAT OR ZA LIMIT	GROUPIE
	7- 8	I2	LOWER MF LIMIT	GROUPIE
	9-11	I3	LOWER MT LIMIT	GROUPIE
	12-17	I11	UPPER MAT OR ZA LIMIT	GROUPIE
	18-19	I2	UPPER MF LIMIT	GROUPIE
	20-22	I3	UPPER MT LIMIT	GROUPIE
			UP TO 100 RANGES MAY BE SPECIFIED, ONE RANGE	GROUPIE
			PER LINE. THE LIST OF RANGES IS TERMINATED	GROUPIE
			BY A BLANK CARD. IF THE UPPER MAT OR ZA	GROUPIE
			LIMIT IS LESS THAN THE LOWER LIMIT THE UPPER	GROUPIE
			IS SET EQUAL TO THE LOWER LIMIT. IF THE UPPER	GROUPIE
			MF OR MT LIMIT IS ZERO IT WILL BE SET EQUAL	GROUPIE
			TO ITS MAXIMUM VALUE, 99 OR 999, RESPECTIVELY	GROUPIE
			IF THE FIRST REQUEST LINE IS BLANK IT WILL	GROUPIE
			TERMINATE THE LIST OF REQUESTS AND CAUSE ALL	GROUPIE
			DATA TO BE RETRIEVED (SEE EXAMPLE INPUT).	GROUPIE
				GROUPIE
VARY	1-66	6E11.4	ENERGY GROUP BOUNDARIES. ONLY REQUIRED IF	GROUPIE
			THE NUMBER OF GROUPS INDICATED ON THE FIRST	GROUPIE
			INPUT CARD IS POSITIVE. ALL ENERGIES MUST	GROUPIE
			BE IN ASCENDING ENERGY IN EV. THE PRESENT	GROUPIE
			LIMITS ARE 1 TO 20,000 GROUPS. FOR N GROUPS	GROUPIE
			N+1 BOUNDARIES WILL BE READ FROM THE	GROUPIE
			INPUT FILE, E.G. IF THE FIRST INPUT CARD	GROUPIE
			INDICATES 20 GROUPS, 21 ENERGY BOUNDARIES	GROUPIE

