				G
========	:====:	========		Groupie
PROGRAM	I GROUI	PIE		Groupie
VERSION	76-1	(NOVEMBER 1976	5)	Groupie
VERSION	79-1	(OCTOBER 1979)	CDC-7600 AND CRAY-1 VERSION.	Groupie
VERSION	80-1	(MAY 1980) IBN	1, CDC AND CRAY VERSION	Groupie
VERSION	1 81-1	(JANUARY 1981)	EXTENSION TO 3000 GROUPS	Groupie
		(MARCH 1981)		Groupie
			BUILT-IN 1/E WEIGHTING SPECTRUM	Groupie
			IMPROVED COMPUTER COMPATIBILITY	Groupie
VERSION	1 83-I	(JANUARY 1983)	*MAJOR RE-DESIGN. *ELIMINATED COMPUTER DEPENDENT CODING.	Groupie
			*NEW, MORE COMPATIBLE I/O UNIT NUMBERS.	Groupie Groupie
			*NEW MULTI-BAND LIBRARY BINARY FORMAT.	Groupie
VERSION	1 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	1 84-1	(APRIL 1984)	ADDED MORE BUILT IN MULTIGROUP ENERGY	Groupie
			STRUCTURES.	Groupie
VERSION	1 85-1	(APRIL 1985)	*UPDATED FOR ENDF/B-VI FORMATS.	Groupie
			*SPECIAL I/O ROUTINES TO GUARANTEE	Groupie
			ACCURACY OF ENERGY. *DOUBLE PRECISION TREATMENT OF ENERGY	Groupie 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	1 88-1	(JULY 1988)	*OPTIONINTERNALLY DEFINE ALL I/O	Groupie
			FILE NAMES (SEE, SUBROUTINES FILIO1	Groupie
			FILIO2 FOR DETAILS). *IMPROVED BASED ON USER COMMENTS.	Groupie Groupie
VERSION	T 89_1	(.TANIIARY 1989)	PSYCHOANALYZED BY PROGRAM FREUD TO	Groupie
VERBIO	. 05 1	(OIMOIMEI IDOD)	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	1 91-1	(JUNE 1991)	*INCREASED PAGE SIZE FROM 1002 TO 5010	Groupie
			POINTS **HDDATED DAGED ON HGED COMMENTS	Groupie
			*UPDATED BASED ON USER COMMENTS *ADDED FORTRAN SAVE OPTION	Groupie Groupie
			*COMPLETELY CONSISTENT ROUTINE TO READ	Groupie
			FLOATING POINT NUMBERS.	Groupie
VERSION	1 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. *ALLOW SELECTION OF ZA/MF/MT OR	Groupie 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
	- 00 -	/ 	RANGES.	Groupie
VERSION	1 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	(TANIJARY 1996)	*COMPLETE RE-WRITE	Groupie
721102011 70 2	(0111011111 1330)	*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	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
VIIIOIOIN JJ I	(Innien 1999)	POINT READ FOR MORE DIGITS	Groupie
		*UPDATED TEST FOR ENDF/B FORMAT	Groupie
		VERSION BASED ON RECENT FORMAT CHANGE	-
		*GENERAL IMPROVEMENTS BASED ON	Groupie
		USER FEEDBACK	Groupie
VERSION 99-2	(TIINE 1999)	*ASSUME ENDF/B-VI, NOT V, IF MISSING	Groupie
V21101011	(00112 2)))	MF=1, MT-451.	Groupie
VERS. 2000-1	(FEBRUARY 2000) *ADDED MF=10, ACTIVATION CROSS SECTION	-
121121 2000 2	(122101111 2000	PROCESSING.	Groupie
		*GENERAL IMPROVEMENTS BASED ON	Groupie
		USER FEEDBACK	Groupie
VERS 2002-1	(FEBRIJARY 2002	*)*ADDED TART 700 GROUP STRUCTURE	Groupie
	,	*ADDED VARIABLE SIGMAO INPUT OPTION	Groupie
	(MAY 2002)	*OPTIONAL INPUT PARAMETERS	Groupie
	(NOV. 2002)	*ADDED SAND-II EXTENDED DOWN TO	Groupie
	/	1.0E-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 -	-
		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. OF	Groupie
	,	2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2	

*INCREASED PAGE SIZE FROM 120,000 TO	Groupie
600,000 POINTS	Groupie
	Groupie
OWNED, MAINTAINED AND DISTRIBUTED BY	Groupie
THE WISTERS DATA COCKETON	Groupie
THE NUCLEAR DATA SECTION	Groupie
INTERNATIONAL ATOMIC ENERGY AGENCY P.O. BOX 100	Groupie Groupie
A-1400, VIENNA, AUSTRIA	Groupie
EUROPE	Groupie
201012	Groupie
ORIGINALLY WRITTEN BY	Groupie
	Groupie
DERMOTT E. CULLEN	Groupie
UNIVERSITY OF CALIFORNIA	Groupie
LAWRENCE LIVERMORE NATIONAL LABORATORY	Groupie
L-159	Groupie
P.O. BOX 808	Groupie
LIVERMORE, CA 94550	Groupie
U.S.A. TELEPHONE 925-423-7359	Groupie
E. MAIL CULLEN1@LLNL.GOV	Groupie Groupie
WEBSITE HTTP://WWW.LLNL.GOV/CULLEN1	Groupie
WEDDITE WITH A PROPERTY OF THE	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	_
READ ALL OF THESE COMMENTS BEFORE IMPLEMENTATION, PARTICULARLY	Groupie
THE COMMENTS CONCERNING MACHINE DEPENDENT CODING.	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	_
IT WOULD BE APPECIATED 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
THE PROGRAM TO REGIONER TO GALOW ATTE AND COMPLICATION OF	Groupie
THIS PROGRAM IS DESIGNED TO CALCULATE ANY COMBINATION OF THE FOLLOWING QUANTITIES FROM LINEARLY INTERPOLABLE TABULATED	Groupie Groupie
CROSS SECTIONS IN THE ENDF/B FORMAT	Groupie
CROSS SECTIONS IN THE EMDITE FORMAL	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 TERMINOLOGYENDF/B	Groupie
TAPEWILL BE USED. IN FACT THE ACTUAL MEDIUM MAY BE TAPE, CARDS,	Groupie
DISK OR ANY OTHER MEDIUM.	Groupie
ENDE /D EODMAT	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 CORRECTLY OUTPUT ON ALL CARDS. THE FORMAT OF SECTION MF=1, MT=451 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 THE CORRECTNESS OR INCORRECTNESS OF ALL OTHER SECTIONS.

ALL FILE 3 CROSS SECTIONS THAT ARE USED BY THIS PROGRAM MUST BE LINEARLY INTERPOLABLE IN ENERGY AND CROSS SECTION (ENDF/B INTERPOLATION LAW 2). FILE 3 BACKGROUND CROSS SECTIONS MAY BE MADE Groupie LINEARLY INTERPOLABLE USING PROGRAM LINEAR (UCRL-50400, VOL. 17, 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 LINEARLY INTERPOLABLE THIS PROGRAM WILL TERMINATE EXECUTION.

CONTENTS OF OUTPUT

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.

DOCUMENTATION

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.

****************** PROGRAM GROUPIE (2007-1) **********

UNSHIELDED GROUP AVERAGES USING 69 GROUPS (WIMS) MAXWELLIAN, 1/E AND FISSION WEIGHTING SPECTRUM

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

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 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.

REACTION INDEX

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

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 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.

SECTION SIZE

Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie

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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 SECTION MAY BE REPRESENTED BY 200,000 DATA POINTS. Groupie Groupie Groupie SELECTION OF DATA 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 Groupie INTERPOLABLE. 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) 650 GROUP (TART STRUCTURE) Groupie (12) 700 GROUP (TART STRUCTURE) Groupie (13) 665 GROUP (SAND-II STRUCTURE, 1.0e-5 eV, UP TO 18 MEV) Groupie (14) 685 GROUP (SAND-II STRUCTURE, 1.0e-5 eV, UP TO 20 MEV) Groupie Groupie GROUP AVERAGES Groupie Groupie THIS PROGRAM DEFINES GROUP AVERAGED CROSS SECTIONS AS... Groupie Groupie (INTEGRAL E1 TO E2) (SIGMA(E)*S(E)*WT(E)*DE) Groupie Groupie (INTEGRAL E1 TO E2) (S(E)*WT(E)*DE) Groupie

AVERAGE = GROUP AVERAGED CROSS SECTION

Groupie Groupie

Groupie

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E1, E2 = ENERGY LIMITS OF THE GROUP
                                                                  Groupie
SIGMA(E) = ENERGY DEPENDENT CROSS SECTION FOR ANY GIVEN REACTION
                                                                  Groupie
       = ENERGY DEPENDENT WEIGHTING SPECTRUM
                                                                  Groupie
       = 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
                                                                  Groupie
(3) MAXWELLIAN = E*EXP(-E/KT)/KT
                                               (0.0 \text{ TO } 4*\text{KT})
        = C1/E
                                               (4*KT TO 67 KEV)
                                                                  Groupie
    1/E
              = C2*EXP(-E/WA)*SINH(SQRT(E*WB)) (ABOVE 67 KEV)
                                                                  Groupie
                                                                  Groupie
          = 0.253 EV (293 KELVIN)
   KТ
                                                                  Groupie
          = 9.65E+5
   WA
                                                                  Groupie
           = 2.29E-6
   WB
                                                                  Groupie
   C1, C2 = 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
                                                                  Groupie
S(E)
         - ENERGY DEPENDENT WEIGHTING SPECTRUM (DEFINED BY
           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
          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
```

Groupie Groupie

THE PROGRAM WILL USE ONE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E) AND 25 DIFFERENT BONDERENKO TYPE SELF-SHIELDING FACTORS (25 SIGMAO Groupie AND N COMBINATIONS) TO DEFINE 25 DIFFERENT AVERAGE CROSS SECTIONS, Groupie FOR EACH REACTION, WITHIN EACH GROUP.

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THE 25 WEIGHTING FUNCTIONS USED ARE....

- UNSHIELDED CROSS SECTIONS (N=0) (1)
- (2-22) PARTIALLY SHIELDED CROSS SECTIONS (N=1 , VARIOUS SIGMA0) THE VALUES OF SIGMAO USED WILL BE EITHER,
 - (A) THE VALUES OF SIGMAO THAT ARE USED VARY FROM 1024 TIMES THE UNSHIELDED TOTAL CROSS SECTIONS IN STEPS OF 1/2 Groupie DOWN TO 1/1024 TIMES THE UNSHIELDED TOTAL CROSS SECTION (A RANGE OF OVER 1 MILLION, CENTERED ON THE UNSHIELDED TOTAL CROSS SECTION WITHIN EACH GROUP).
 - (B) THE SAME CONSTANT VALUES OF SIGMAO IN EACH GROUP. THE Groupie VALUES OF SIGMAO USED INCLUDE 40000, 20000, 10000, 7000, 4000, 2000, 1000, 700, 400, 200, 100, 70, 40, 20, 10, 7, 4, 2, 1, 0.7, 0.4 (A RANGE OF 100,000 SPANNING MORE THAN THE RANGE OF SIGMAO VALUES THAT MAY BE ENCOUNTERED IN ACTUAL APPLICATIONS)
- (23) TOTALLY SHIELDED FLUX WEIGHTED CROSS SECTION (N=1, SIGMA0=0)
- TOTALLY SHIELDED CURRENT WEIGHTED CROSS SECTION (N=2, SIGMA0=0)
- (25) TOTALLY SHIELDED COSINE SQUARED WEIGHTED CROSS SECTION (N=3, SIGMA0=0)

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

CALCULATION OF RESONANCE INTEGRALS

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

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

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

THE RESONANCE INTEGRAL IS DEFINED AS,

= (INTEGRAL E1 TO E2) (SIGMA(E)*S(E)*WT(E)*DE) RΙ

WHERE NORMALLY,

= 1/ES(E)

WT(E) = 1 - NO SELF-SHIELDING

FROM THE ABOVE DEFINITION OF GROUP AVERAGED CROSS SECTIONS THE RESONANCE INTEGRAL IS,

RΙ = AVERAGE * (INTEGRAL E1 TO E2) (S(E)*WT(E)*DE) Groupie Groupie Groupie

Groupie Groupie

Groupie Groupie

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FOR A 1/E SPECTRUM AND NO SELF-SHIELDING THIS REDUCES TO, = AVERAGE* LOG(E2/E1)

IN ANY OTHER SITUATION, INCLUDING ABSORPTION AND/OR ENERGY DEPENDENT CROSS SECTIONS, THE SPECTRUM WILL NOT BE 1/E -ABSORPTION WILL TEND TO DECREASE THE SPECTRUM PROGRESSIVELY MORE AT LOWER ENERGIES - ENERGY DEPENDENCE OF THE CROSS SECTION WILL LEAD TO SELF-SHIELDING.

HERE WE WILL NOT ATTEMPT TO PERFORM A DETAILED SPECTRUM CALCULATION TO ACCOUNT FOR ABSORPTION.

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

= AVERAGE* LOG(E2/E1) RΙ

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

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

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 HAVE SELECTED A 1/E SPECTRUM - THERE IS NO CHECK ON THIS - THE PROGRAM MERELY MULTIPLIES THE GROUP AVERAGED CROSS SECTIONS BY,

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

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 Groupie

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

RESOLVED RESONANCE REGION

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

UNRESOLVED RESONANCE REGION

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

ACCURACY OF RESULTS

ALL INTEGRALS ARE PERFORMED ANALYTICALLY. THEREFORE NO ERROR IS INTRODUCED DUE TO THE USE OF TRAPAZOIDAL OR OTHER INTEGRATION SCHEME. THE TOTAL ERROR THAT CAN BE ASSIGNED TO THE RESULTING AVERAGES IS JUST THAT DUE TO THE ERROR IN THE CROSS SECTIONS AND ENERGY DEPENDENT WEIGHTING SPECTRUM. GENERALLY SINCE THE THE ENERGY DEPENDENT WEIGHTING SPECTRUM APPEARS IN BOTH THE NUMERATOR AND THE DENOMINATOR THE AVERAGES RAPIDLY BECOME INSENSITIVE TO THE WEIGHTING SPECTRUM AS MORE GROUPS ARE USED. SINCE THE WEIGHTING SPECTRUM IS LOADED IN THE PAGING SYSTEM THE USER CAN DESCRIBE THE SPECTRUM TO ANY REQUIRED ACCURACY USING ANY NUMBER OF ENERGY VS. SPECTRUM PAIRS.

MULTI-BAND PARAMETERS

MULTI-BAND PARAMETERS ARE CALCULATED FOR THE TOTAL, ELASTIC, CAPTURE AND FISSION REACTIONS. WITH THE NUMBER OF GROUPS THAT ARE NORMALLY USED (SEE BUILT IN GROUP STRUCTURES) ALL OTHER REACTIONS RESULT IN A NEGLIGABLE AMOUNT OF SELF-SHIELDING. AS SUCH THEIR EQUIVALENT BAND CROSS SECTION WILL MERELY BE THEIR UNSHIELDED VALUE WITHIN EACH BAND.

FOR ANY GIVEN EVALUATION, WITHIN ANY GIVEN GROUP THIS PROGRAM WILL GENERATE THE MINIMUM NUMBER OF BANDS REQUIRED WITHIN THAT GROUP. AS OUTPUT TO THE COMPUTER READABLE DISK FILE THE BAND PARAMETERS FOR EACH EVALUATION WILL BE FORMATTED TO HAVE THE SAME NUMBER OF BANDS IN ALL GROUPS (WITH ZERO WEIGHT FOR SOME BANDS WITHIN ANY GROUP). THE USER MAY DECIDE TO HAVE OUTPUT

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EITHER WITH THE MINIMUM NUMBER OF BANDS REQUIRED FOR EACH EVALUATION (E.G. 2 BANDS FOR HYDROGEN AND 4 BANDS FOR U-233) OR THE SAME NUMBER OF BANDS FOR ALL EVALUATIONS (E.G. 4 BANDS FOR BOTH HYDROGEN AND U-233).

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

ENDF/B FORMATTED UNSHIELDED AVERAGES

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

MIXTURES OF MATERIALS AND RESONANCE OVERLAP

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

MULTI-BAND PARAMETER OUTPUT FORMAT

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

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

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

THE FORMAT OF THE CHARACTER FILE IS,

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	12-22	I11	NUMBER GROUPS	Groupie
	23-33	I11	NUMBER OF BANDS	Groupie

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UNIT 31 32 33 34 311 SCRATT UNIT 8 9 12 13	DESCRIPTION MULTI-BAND PARAMETE (BCD - 80 CHARACTER SELF-SHIELDED CROSS (BCD - 120 CHARACTE MULTI-BAND PARAMETE (BCD - 120 CHARACTE UNSHIELDED CROSS SE (BCD - 120 CHARACTE OUTPUT REPORT (BCD MULTI-GROUP ENDF/B (BCD - 80 CHARACTER CH FILES FILENAME DESCRIPTI ENERGY DEPENDENT WE (BINARY - 40080 WOR TOTAL CROSS SECTION (BINARY - 40080 WOR CAPTURE CROSS SECTI (BINARY - 40080 WOR	S/RECORD) SECTION LISTING - OPTIONAL RS/RECORD) R LISTING - OPTIONAL RS/RECORD) CTION LISTING - OPTION RS/RECORD) - 80 CHARACTERS/RECORD) DATA - OPTIONAL S/RECORD) ON IGHTING SPECTRUM DS/BLOCK) ON - ONLY FOR SELF-SHIELDING CALCULATION	Groupie

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UNIT	FILE N	 лмг		Groupie Groupie
				Groupie
2	GROUPI			Groupie
3	GROUPI			Groupie
8	(SCRAT			Groupie
9	(SCRAT			Groupie
10	ENDFB.			Groupie
11	ENDIB.			Groupie
12	(SCRAT			Groupie
13	(SCRAT			Groupie
14	(SCRAT			Groupie
31	MULTBA			Groupie
32	SHIELD			Groupie
33	MULTBA			Groupie
34	UNSHIE			Groupie
54	ONDITE	пр.прі		Groupie
т/О	UNITS U	CED.		Groupie
				Groupie
			10 WILL ALWAYS BE USED.	Groupie
			AND 11 ARE OPTIONALLY USED DEPENDING ON THE	Groupie
	UT REQU		AND II ARE OFFICINALLY OSED DEFENDING ON THE	Groupie
			WILL ONLY BE USED IF SELF-SHIELDED OR	Groupie
	-		REQUESTED.	Groupie
МОПТ	IBAND O	01701 15	KEQUESTED.	Groupie
TNIDIT	CARDS			Groupie
				Groupie
	COLS.	FORMAT	DESCRIPTION	Groupie
				Groupie
1	1-11	I11	SELECTION CRITERIA (0=MAT, 1=ZA)	Groupie
1	12-22	I11	NUMBER OF GROUPS.	Groupie
	12 22	111	=.GT.0 - ARBITRARY GROUP BOUNDARIES ARE READ	Groupie
			FROM INPUT FILE (N GROUPS REQUIRE	Groupie
			N+1 GROUP BOUNDARIES). CURRENT	Groupie
			PROGRAM MAXIMUM IS 1000 GROUPS.	Groupie
			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
			- 5 ORIVE 171 GROOLD	GIGUPIC
			= -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 650 GROUPS	Groupie
			=-12 - TART 700 GROUPS	Groupie
			=-13 - SAND-II 665 GROUPS TO 18 MEV	Groupie
			=-14 - SAND-II 685 GROUPS TO 20 MEV	Groupie
1	23-33	I11	MULTI-BAND SELECTOR	Groupie
_	20 00		= 0 - NO MULTI-BAND CALCULATIONS	Groupie
			= 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+SIGMAO)) WHERE	Groupie
			SIGMAO = AV(TOT) IN EACH GROUP	Groupie
			= 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
				CICAPIC

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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
                                                                      Groupie
                        BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR
                        ALL ISOTOPES.
                                                                      Groupie
        34-44 I11
                        NUMBER OF POINTS USED TO DESCRIBE ENERGY
                                                                      Groupie
                        DEPENDENT WEIGHTING SPECTRUM S(E).
                                                                      Groupie
                               - MAXWELLIAN - UP TO 0.1 EV
                                                                      Groupie
                                           - 0.1 EV TO 67 KEV
- ABOVE 67 KEV
                                  1/E
                                                                      Groupie
                                  FISSION
                                                                      Groupie
05/01/20-----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
        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
                                                                     Groupie
        56-66
                  I11
                        SIGMA-0 DEFINITION SELECTOR.
                                                                      Groupie
                         < 0 - 21 VALUES OF SIGMAO 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).
                                                                      Groupie
                        IF SIGMA-0 DEFINITION SELECTOR < 0, THE NEXT Groupie
    2-4 1-66 6D11.4
                         4 LINES OF INPUT ARE THE 22 VALUES OF SIGMAO, Groupie
                         6 PER LINE.
                                                                      Groupie
                        ENDF/B INPUT DATA FILENAME
                                                                      Groupie
      2
          1-60
                  A60
                        (STANDARD OPTION = ENDFB.IN)
                                                                      Groupie
                        ENDF/B OUTPUT DATA FILENAME
                                                                      Groupie
          1-60
                  A60
                         (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
                        SELF-SHIELDED CROSS SECTION LISTING
                                                                      Groupie
                         = 1 - CROSS SECTIONS
                                                                      Groupie
                         = 2 - RESONANCE INTEGRALS
                                                                      Groupie
          12-22
                        MULTI-BAND PARAMETER LISTING
                  T11
                                                                      Groupie
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4 4	23-33 34-44	I11 I11	MULTI-BAND PARAMETERS COMPUTER READABLE UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT = 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1)	Groupie Groupie Groupie
4	45-55	I11	= 2 - LINEAR-LINEAR (INTERPOLATION LAW 2) UNSHIELDED CROSS SECTIONS LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS	Groupie Groupie Groupie Groupie
05/01/20 - 4	ADDED 56-66		IF THE STANDARD BUILT-IN SPECTRA IS USED, INPUT LINE 1, COLUMNS 34-44 = 2, THIS FIELD CAN BE USED TO OPTIONALLY CHANGE TEMPERATURE OF THE MAXWELLIAN. INPUT IS IN EV (0.0253 EV = ROOM TEMPERATURE) = 0 - USE DEFAULT 0.0253 EV, ROOM TEMPERATURE > 0 - USE THIS AS THE TEMPERATURE RESTRICTION - TEMPERATURE CANNOT EXCEED 1000 EV.	_
5	1-80	18A4	LIBRARY IDENTIFICATION. ANY TEXT THAT THE USER WISHES TO IDENTIFY THE MULTI-BAND PARAMETERS. THIS LIBRARY IDENTIFICATION IS WRITTEN INTO THE COMPUTER READABLE MULTI-BAND DATA FILE.	Groupie Groupie Groupie
6-N	1- 6	16	LOWER MAT OR ZA LIMIT	Groupie
	7- 8	I2	LOWER MF LIMIT	Groupie
	9-11	I3	LOWER MT LIMIT	Groupie
	12-17 18-19	I11 I2	UPPER MAT OR ZA LIMIT UPPER MF LIMIT	Groupie Groupie
	20-22	12 13	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	_
			MF OR MT LIMIT IS ZERO IT WILL BE SET EQUAL 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	6D11.4	ENERGY GROUP BOUNDARIES. ONLY REQUIRED IF	Groupie
			THE NUMBER OF GROUPS INDICATED ON THE FIRST	Groupie
			INPUT CARD IS POSITIVE. ALL ENERGIES MUST BE IN ASCENDING ENERGY IN EV. THE PRESENT	Groupie Groupie
			LIMITS ARE 1 TO 1000 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
			WILL BE READ FROM THE INPUT FILE.	Groupie
VARY	1-66	6D11 /	ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY	Groupie Groupie
VARI	1-00	0011.4	REQUIRED IF THE NUMBER OF POINTS INDICATED	Groupie
			ON FIRST CARD IS MORE THAN ONE. DATA IS	Groupie
			GIVEN IN (ENERGY, WEIGHT) PAIRS, UP TO 3	Groupie
			PAIRS PER CARD, USING ANY NUMBER OF CARDS	Groupie
			REQUIRED. ENERGIES MUST BE IN ASCENDING	Groupie
			ORDER IN EV. THE SPECTRUM VALUES MUST BE	Groupie
			NON-NEGATIVE. THE ENERGY RANGE OF SPECTRUM MUST AT LEAST SPAN THE ENERGY RANGE OF THE	Groupie Groupie
			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

т.							Groupi
	EXAMPLE INPU	וי אור דו					Groupi
							Groupi
			ND PROCESS A	LL DATA (AL	I MAT BETWEE	EN 1 AND	Groupi
				UCTURE, GENI			Groupi
				TO 0.1 PER-			Groupi
				OUT ALL LIST			Groupi
			ORMAT GROUP				Groupi
		•					Groupi
F	EXPLICITLY S	SPECIFY THE	E STANDARD F	'ILENAMES.			Groupi
							Groupi
T	THE FOLLOWIN	IG 7 INPUT	LINES ARE R	EQUIRED.			Groupi
							Groupi
	0	0	-2	0 1.000	000-03	0	Groupi
ENDFE	B.IN						Groupi
ENDFF	3.OUT						Groupi
	1	1	1	1	1		Groupi
TART			BRARY TO 0.1	PER-CENT A	CCURACY		Groupi
1	l 1 1 9999	0 0					Groupi
		(BLA	ANK CARD TER	MINATES REQU	UEST LIST)		Groupi
							Groupi
	EXAMPLE INPU						Groupi
							Groupi
				Y THE ENDF/			Groupi
	•		•	(U-238 AT :	300 KELVIN)	AND	Groupi
V	VRITTEN TO \	ENDFB6\GR	OUPIE\K300\Z	A092238			Groupi
							Groupi
1	THE FOLLOWIN	NG 7 INPUT	LINES ARE R	EQUIRED.			Groupi
				~			Groupi Groupi
	0	0	-2	EQUIRED.	000-03	0	Groupi Groupi Groupi
\ENDF	0 FB6\SIGMA1\K	0 (300\ZA092)	-2 238	~	000-03	0	Groupi Groupi Groupi Groupi
\ENDF	0 FB6\SIGMA1\K FB6\GROUPIE\	0 300\zA092 K300\zA092	-2 238 2238	0 1.000		0	Groupi Groupi Groupi Groupi Groupi
\ENDF \ENDF	0 FB6\SIGMA1\K FB6\GROUPIE\ 1	0 3300\ZA0922 K300\ZA092	-2 238 2238	0 1.000	1	0	Groupi Groupi Groupi Groupi Groupi Groupi
\ENDF \ENDF TART	0 FB6\SIGMA1\K FB6\GROUPIE\ 1 175 GROUP,	0 3300\ZA0922 K300\ZA092 1 2 BAND LIF	-2 238 2238	0 1.000	1	0	Groupi Groupi Groupi Groupi Groupi Groupi Groupi
\ENDF \ENDF TART	0 FB6\SIGMA1\K FB6\GROUPIE\ 1	0 3300\za0922 K300\za092 1 2 BAND LIE	-2 238 2238 1 BRARY TO 0.1	0 1.000	1 CCURACY	0	Groupi Groupi Groupi Groupi Groupi Groupi Groupi Groupi
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\ENDF \ENDF TART 1	0 FB6\SIGMA1\K FB6\GROUPIE\ 1 175 GROUP,	0 3300\ZA0922 \K300\ZA092 1 2 BAND LIH 0 0 0 (BL2	-2 238 2238 1 BRARY TO 0.1	0 1.000	1 CCURACY	0	Groupi
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\ENDF \ENDF TART 1 E	0 FB6\SIGMA1\K FB6\GROUPIE\ 1 175 GROUP, 1 1 9999 EXAMPLE INPU	0 (3300\ZA092) (K300\ZA092) 1 2 BAND LIH 0 0 0 (BLA UT NO. 3 DATA. USE	-2 238 2238 1 BRARY TO 0.1 ANK CARD TER	0 1.000 1 PER-CENT AGE MINATES REQUE ING IN ORDER	1 CCURACY UEST LIST) TO CALCULAT	TE.	Groupi
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