				Groupie
PROGRAM	GROUI	PIE		Groupie
		==		Groupie
VERSION	76-1	(NOVEMBER 1976	)	Groupie
			CDC-7600 AND CRAY-1 VERSION.	Groupie
			, CDC AND CRAY VERSION	Groupie
			EXTENSION TO 3000 GROUPS	Groupie
		(MARCH 1981) I		Groupie
			BUILT-IN 1/E WEIGHTING SPECTRUM	Groupie
			IMPROVED COMPUTER COMPATIBILITY	Groupie
			*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
ZERSION	86-2	(JUNE 1986)	*BUILT-IN MAXWELLIAN, 1/E AND FISSION	Groupie
			WEIGHTING SPECTRUM.	Groupie
VERSION	88-1	(JULY 1988)	*OPTIONINTERNALLY 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
			*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
/ERSION	94-1	(JANUARY 1994)	*VARIABLE ENDF/B DATA FILENAMES	Groupie
			TO ALLOW ACCESS TO FILE STRUCTURES	Groupie

		HAS BEEN CHANGED)	Crownia
		*CLOSE ALL FILES BEFORE TERMINATING	Groupie 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 INDEPENDENCE	Groupie
		*ALL DOUBLE PRECISION	Groupie
		*ON SCREEN OUTPUT *UNIFORM TREATMENT OF ENDF/B I/O	Groupie 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 (PREVIOUSLY 100)	Groupie 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 *UPDATED TEST FOR ENDF/B FORMAT	Groupie Groupie
		VERSION BASED ON RECENT FORMAT CHANGE	-
		*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	
		PROCESSING. *GENERAL IMPROVEMENTS BASED ON	Groupie Groupie
		USER FEEDBACK	Groupie
VERS. 2002-1	(FEBRUARY 2002	2) *ADDED TART 700 GROUP STRUCTURE	Groupie
		*ADDED VARIABLE SIGMA0 INPUT OPTION	Groupie
	(MAY 2002)	<b>*OPTIONAL INPUT PARAMETERS</b>	Groupie
	(NOV. 2002)	*ADDED SAND-II EXTENDED DOWN TO	Groupie
	(	1.0D-5 EV.	Groupie
	(JUNE 2003)	*CORRECTED SAND-II 620 AND 640 GROUP ENERGY BOUNDARIES DEFINITIONS.	Groupie 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. *DO NOT SKIP LOW TOTAL ENERGY RANGES	Groupie 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
	( TAN 0007)	BUILT-IN STANDARD SPECTRUM.	Groupie
VERS. 2007-1	(JAN. 2007)	*CHECKED AGAINST ALL ENDF/B-VII.	Groupie
		*INCREASED PAGE SIZE FROM 120 000 TO	Ground
		<pre>*INCREASED PAGE SIZE FROM 120,000 TO 600,000 POINTS</pre>	Groupie Groupie
VERS. 2008-1	(JAN. 2008)	*INCREASED PAGE SIZE FROM 120,000 TO 600,000 POINTS *72 CHARACTER FILE NAMES.	Groupie Groupie Groupie
VERS. 2008-1	(JAN. 2008)	600,000 POINTS	Groupie
VERS. 2008-1 VERS. 2010-1		600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000	Groupie Groupie Groupie Groupie
		600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS.	Groupie Groupie Groupie Groupie Groupie
		600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS. *ADDED OUTPUT TO PLOT/COMPARE SHIELDED	Groupie Groupie Groupie Groupie Groupie
VERS. 2010-1	(Apr. 2010)	600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS. *ADDED OUTPUT TO PLOT/COMPARE SHIELDED AND UNSHIELDED CROSS SECTIONS.	Groupie Groupie Groupie Groupie Groupie Groupie
	(Apr. 2010)	600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS. *ADDED OUTPUT TO PLOT/COMPARE SHIELDED AND UNSHIELDED CROSS SECTIONS. *Corrected TART 700 groups to extend up	Groupie Groupie Groupie Groupie Groupie Groupie Groupie
VERS. 2010-1	(Apr. 2010)	600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS. *ADDED OUTPUT TO PLOT/COMPARE SHIELDED AND UNSHIELDED CROSS SECTIONS.	Groupie Groupie Groupie Groupie Groupie Groupie
VERS. 2010-1	(Apr. 2010) (June 2011)	600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS. *ADDED OUTPUT TO PLOT/COMPARE SHIELDED AND UNSHIELDED CROSS SECTIONS. *Corrected TART 700 groups to extend up to 1 GeV (1,000 MeV) - previously it	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
VERS. 2010-1 VERS. 2011-1	(Apr. 2010) (June 2011)	600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS. *ADDED OUTPUT TO PLOT/COMPARE SHIELDED AND UNSHIELDED CROSS SECTIONS. *Corrected TART 700 groups to extend up to 1 GeV (1,000 MeV) - previously it was ERRONEOUSLY cutoff at 20 MeV. *Corrected TART 616 groups lowest energy from 1.0D-4 eV to 1.0D-5 eV.	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
VERS. 2010-1 VERS. 2011-1	(Apr. 2010) (June 2011)	600,000 POINTS *72 CHARACTER FILE NAMES. *GENERAL UPDATES *INCREASED WEIGHTING SPECTRUM TO 30,000 FROM 3,000 ENERGY POINTS. *ADDED OUTPUT TO PLOT/COMPARE SHIELDED AND UNSHIELDED CROSS SECTIONS. *Corrected TART 700 groups to extend up to 1 GeV (1,000 MeV) - previously it was ERRONEOUSLY cutoff at 20 MeV. *Corrected TART 616 groups lowest	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie 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	
				<pre>is not explicitly defined - extension = 0 is standard, constant is NOT, so</pre>	Groupie Groupie
				constant extension is NOT RECOMMENDED.	_
VERS.	2012-1	(Aug.	2012)	*Added CODENAME	Groupie
		(y ·	,	*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 weigthing spectrum	
				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. *Extended SAND-II to 60, 150, 200 MeV.	Groupie Groupie
VERS	2015-2	(Mar	2015)	*Deleted 1P from formats reading input	Groupie
	2	·		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 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
				explcitly 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
				*METHODB was incorrecctly named	Groupie
				METHOD in one routine = corrected.	Groupie
				*Default multi-band is method #2 =	Groupie
				conserve <x>, &lt;1/(x+<x>&gt;, &lt;1/x&gt;.</x></x>	Groupie
				*Definition of built-in group structure	-
				using SUBROUTINE GROPE is identical for GROUPIE and VIRGIN.	Groupie 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
	2018-1			*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

from Resolved (supersedes URRDO and Groupie URRFIT codes). Groupie \*Added entire self-shielding array to Groupie memory - previously only one group Groupie results were in memory - saving ALL Groupie greatly simplifies the logic. Groupie \*Additional Interpolation Law Tests Groupie \*Check maximum Tabulated Energy of MTs Groupie Groupie to insure they ALL end at the same energy. Groupie \*Multi-band = 1 no longer allowed. Groupie The only allowed values are, Groupie 0 = no multi-band calculations, or, Groupie 2 = Conserve 1/[total + <total>] Groupie \*Unresolved Resonance Region Groupie Self-Shielding Requires all of these, Groupie 1) Unresolved data with ENDF input Groupie 2) 616 TART Groups (input -11) Groupie 3) Define Sigma0 standard (input = 0) Groupie \*Unresolved Resonance Region Groupie Self-Shielding Always Outputs, Groupie 1) LSSF = 0 = Output cross sections Groupie 2) INTUNR = 2 = Interpolation law Groupie \*Added ZAzzzaaa to filenames. Groupie Groupie 2015-2 Acknowledgment 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 ------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 RedCullen1.net/HOMEPAGE.NEW Website 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 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 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 Groupie CONTENTS OF OUTPUT 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 Groupie 69 GROUPS (WIMS) UNSHIELDED GROUP AVERAGES USING 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 20,000. 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 (19)1102 GROUP (UKAEA STRUCTURE, 1.0D-5 eV, UP TO 1 GeV) 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 AVERAGE = -----Groupie Groupie (INTEGRAL E1 TO E2) (S(E)\*WT(E)\*DE) WHERE . . . Groupie Groupie AVERAGE = GROUP AVERAGED CROSS SECTION Groupie = ENERGY LIMITS OF THE GROUP E1, E2 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 =  $E \times EXP(-E/KT)/KT$ (0.0 TO 4\*KT) Groupie 1/E = C1/E(4\*KT TO 67 KEV) Groupie FISSION = C2\*EXP(-E/WA)\*SINH(SQRT(E\*WB)) (ABOVE 67 KEV) Groupie Groupie = 0.253 EV (293 KELVIN) кт Groupie WA = 9.65D+5 Groupie WB = 2.29D-6Groupie 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 20,000 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 20,000 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) \* NGroupie Groupie WHERE... Groupie Groupie S(E) - ENERGY DEPENDENT WEIGHTING SPECTRUM (DEFINED BY Groupie TABULATED VALUES AND LINEAR INTERPOLATION BETWEEN Groupie

	Crownia
TABULATED VALUES). TOTAL(E) - ENERGY DEPENDENT TOTAL CROSS SECTION FOR ONE MATERIAL	Groupie 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 THAT GROUP OR POWERS OF 10 - INPUT OPTION).	Groupie 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 SIGMAO	-
AND N COMBINATIONS) TO DEFINE 25 DIFFERENT AVERAGE CROSS SECTIONS,	
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 SIGMAO THAT ARE USED VARY FROM 1024	Groupie
TIMES THE UNSHIELDED TOTAL CROSS SECTIONS IN STEPS OF 1/2 DOWN TO 1/1024 TIMES THE UNSHIELDED TOTAL CROSS SECTION	Groupie
(A RANGE OF OVER 1 MILLION, CENTERED ON THE UNSHIELDED	Groupie Groupie
TOTAL CROSS SECTION WITHIN EACH GROUP).	Groupie
(B) THE SAME CONSTANT VALUES OF SIGMA0 IN EACH GROUP. THE	Groupie
VALUES OF SIGMAO 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 SIGMAO VALUES THAT MAY BE ENCOUNTERED IN	Groupie
ACTUAL APPLICATIONS) (23) - TOTALLY SHIELDED FLUX WEIGHTED CROSS SECTION	Groupie 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
FOR ALL OTHER REACTIONS (EXCEPT TOTAL, ELASTIC, CAPTURE AND	Groupie
FISSION) THE PROGRAM WILL USE THE ENERGY DEPENDENT WEIGHTING	Groupie 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
IN A PURE ELASTIC ISOTROPICALLY SCATTERING MATERIAL WITH A	Groupie 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 THERE IS ONLY SCATTERING.	Groupie Groupie
THERE IS ONLY SCATTERING.	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 = (INTEGRAL E1 TO E2) (SIGMA(E) $\times$ S(E) $\times$ WT(E) $\times$ DE)	Groupie
	Groupie
WHERE NORMALLY,	Groupie
S(E) = 1/E	
WT(E) = 1 - NO SELF-SHIELDING	Groupie
WT(E) = 1 - NO SELF-SHIELDING	Groupie
	Groupie Groupie
FROM THE ABOVE DEFINITION OF GROUP AVERAGED CROSS SECTIONS THE	Groupie Groupie Groupie
	Groupie Groupie Groupie Groupie
FROM THE ABOVE DEFINITION OF GROUP AVERAGED CROSS SECTIONS THE	Groupie Groupie Groupie
FROM THE ABOVE DEFINITION OF GROUP AVERAGED CROSS SECTIONS THE RESONANCE INTEGRAL IS,	Groupie Groupie Groupie Groupie Groupie

FOR A 1/E SPECTRUM AND NO SELF-SHIELDING THIS REDUCES TO,	Groupie
	Groupie
RI = AVERAGE * 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 Groupie
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,	Groupie
AND THEN DEFINING THE RESONANCE INTEGRAL AS,	Groupie Groupie
RI = AVERAGE* LOG (E2/E1)	Groupie
KI - AVERAGE DOG(EZ/EI)	Groupie
IN ORDER TO CALCULATE RESONANCE INTEGRALS YOU MUST FOLLOW THESE	Groupie
STEPS,	Groupie
51215,	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	-
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	Groupie
STEPS OUTLINED ABOVE IF YOU WISH TO OBTAIN MEANINGFUL	Groupie
RESULTS.	Groupie
	Groupie
NOTE - OUTPUT IN THE ENDF/B FORMAT IS ALWAYS GROUP AVERAGED CROSS	Groupie
SECTIONS, REGARDLESS OF WHETHER YOU ASK FOR AVERAGED CROSS	Groupie
SECTIONS OR RESONANCE INTEGRALS - THIS IS BECAUSE DATA IN	Groupie
THE ENDF/B FORMAT IS EXPLICITLY DEFINED TO BE CROSS	Groupie
SECTIONS.	Groupie
	Groupie
RESONANCE INTEGRAL OUTPUT CAN ONLY BE OBTAINED IN THE	Groupie
LISTING FORMATS.	Groupie
	Groupie
MINIMUM TOTAL CROSS SECTION TREATMENT	Groupie
	Groupie
SINCE THE BONDARENKO SELF-SHIELDING DEPENDS ON 1/TOTAL CROSS	
SECTION, THE ALGORITHM WILL BECOME NUMERICALLY UNSTABLE IF THE	Groupie
TOTAL CROSS SECTION IS NEGATIVE (AS OCCURS IN MANY ENDF/B	Groupie
EVALUATIONS). IF THE TOTAL IS LESS THAN SOME MINIMUM ALLOWABLE	Groupie Groupie
	Groupie Groupie Groupie
VALUE (DEFINE BY OKMIN, PRESENTLY 1 MILLI-BARN) AN ERROR MESSAGE	Groupie Groupie Groupie Groupie
VALUE (DEFINE BY OKMIN, PRESENTLY 1 MILLI-BARN) AN ERROR MESSAGE WILL BE PRINTED AND FOR THE SELF-SHIELDING CALCULATION ALL ENERGY	Groupie Groupie Groupie Groupie Groupie
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	Groupie Groupie Groupie Groupie Groupie
VALUE (DEFINE BY OKMIN, PRESENTLY 1 MILLI-BARN) AN ERROR MESSAGE WILL BE PRINTED AND FOR THE SELF-SHIELDING CALCULATION ALL ENERGY	Groupie Groupie Groupie Groupie Groupie Groupie
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.	Groupie Groupie Groupie Groupie Groupie Groupie Groupie
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	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
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	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
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	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
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 Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
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	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie

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

## 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 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

Groupie Groupie

Groupie Groupie

Groupie

Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie

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 Groupie THE FORMAT OF THE CHARACTER FILE IS, Groupie Groupie RECORD COLUMNS FORMAT DESCRIPTION Groupie 1-72 18A4 LIBRARY DESCRIPTION (AS READ) Groupie 1 MATERIAL ZA 2 1-11 I11 Groupie 12-22 I11 NUMBER GROUPS Groupie NUMBER OF BANDS 23-33 I11 Groupie 34-44 E11.4 TEMPERATURE (KELVIN) Groupie 45-55 1X,10A1 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 1-11 BLANK 4 \_\_\_\_ Groupie 12-22 E11.4 TOTAL (SECOND BAND) Groupie 23-33 E11.4 ELASTIC Groupie 34-44 E11.4 CAPTURE Groupie 35-55 FISSION E11.4 Groupie Groupie LINES 3 AND 4 ARE REPEATED FOR EACH GROUP. THE LAST LINE FOR EACH Groupie MATERIAL (ZA) IS, Groupie Groupie N 1-11 E11.4 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 \_\_\_\_\_ OUTPUT REPORT (BCD - 80 CHARACTERS/RECORD) 3 Groupie 11 MULTI-GROUP ENDF/B DATA - OPTIONAL Groupie (BCD - 80 CHARACTERS/RECORD) Groupie 16 PLOTTAB FORMATTED SELF-SHIELDING RESULTS Groupie (BCD - 80 CHARACTERS/RECORD) Groupie MULTI-BAND PARAMETERS CHARACTER FILE - OPTIONAL 31 Groupie (BCD - 80 CHARACTERS/RECORD) Groupie UNRESOLVED FSELF-SHIELDED PSEUDO ENDF FORMAT - OPTIONAL 32 Groupie (BCD - 120 CHARACTERS/RECORD) Groupie SELF-SHIELDED CROSS SECTION LISTING - OPTIONAL 33 Groupie (BCD - 120 CHARACTERS/RECORD) Groupie MULTI-BAND PARAMETER LISTING - OPTIONAL 34 Groupie (BCD - 120 CHARACTERS/RECORD) Groupie 35 UNSHIELDED CROSS SECTION LISTING - OPTION Groupie (BCD - 120 CHARACTERS/RECORD) Groupie Groupie SCRATCH FILES Groupie \_\_\_\_\_ Groupie UNIT FILENAME DESCRIPTION Groupie \_\_\_\_ -----Groupie ENERGY DEPENDENT WEIGHTING SPECTRUM 8 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 FISSION CROSS SECTION - ONLY FOR SELF-SHIELDING CALCULATION 14 Groupie (BINARY - 40080 WORDS/BLOCK) Groupie Groupie OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINES FILIO1 AND FILIO2) Groupie Groupie UNIT FILE NAME Groupie \_\_\_\_ \_\_\_\_\_ Groupie GROUPIE.INP 2 Groupie 3 GROUPIE.LST Groupie \_\_\_\_\_ Groupie 8 (SCRATCH) Groupie 9 (SCRATCH) Groupie 10 ENDFB.IN Groupie ENDFB.OUT 11 Groupie 12 (SCRATCH) Groupie 13 (SCRATCH) Groupie 14 (SCRATCH) Groupie -----2019/6/23 - New Filenames (added ZAzzzaaa at Beginning)------ Groupie ----- (OLD) ------ (NEW) ------ Groupie 16 PLOTTAB.CUR ZAzzzaaa.PLOT.CUR Groupie 31 MULTBAND. TAB ZAzzzaaa.MULTBAND.TAB Groupie 32 ZAzzzaaa.URR.ENDF Groupie ZAzzzaaa.SHIELD.LST 33 SHIELD.LST Groupie 34 MULTBAND.LST ZAzzzaaa.MULTBAND.LST Groupie ZAzzzaaa.UNSHIELD.LST 35 UNSHIELD.LST Groupie Groupie I/O UNITS USED Groupie Groupie UNITS 2, 3 8, 9 AND 10 WILL ALWAYS BE USED. Groupie UNITS 31 THROUGH 35, 11 AND 16 ARE OPTIONALLY USED DEPENDING Groupie ON THE OUTPUT REQUESTED. Groupie UNITS 12, 13 AND 14 WILL ONLY BE USED IF SELF-SHIELDED OR Groupie MULTIBAND OUTPUT IS REQUESTED. Groupie

		Crownia
INPUT CARDS		Groupie Groupie
		Groupie
CARD COLS. FORMAT	DESCRIPTION	Groupie
		Groupie
1 1-11 I11	SELECTION CRITERIA (0=MAT, 1=ZA)	Groupie
1 12-22 I11		Groupie
	-2019/6/23 -11 (TART 616 groups) required for	Groupie
	Unresolved Resonance Region Self-Shielding	Groupie
	calculation.	Groupie
	=.GT.0 - ARBITRARY GROUP BOUNDARIES ARE READ	Groupie
	FROM INPUT FILE (N GROUPS REQUIRE N+1 GROUP BOUNDARIES). CURRENT	Groupie Groupie
	PROGRAM MAXIMUM IS 20,000 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
	= -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 $= -10 - ABBN 28 GROUPS$	Groupie Groupie
Current TART Standard		Groupie
(-11 is required for		Groupie
unresolved resonance		Groupie
region self-shieldin		Groupie
calculations)	=-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
	=-19 - UKAEA 1102 GROUPS TO 1 GeV	Groupie
1 23-33 I11	MULTI-BAND SELECTOR	Groupie
	-2019/6/23 - ONLY 0 or 2 allowed = the = 1	Groupie
	option has proven to give very poor results,	Groupie
	and therefore is no longer allowed. = 0 - NO MULTI-BAND CALCULATIONS	Groupie
No longer allowed	= 0 - NO MOLTI-BAND CALCULATIONS = 1 - 2 BAND. CONSERVE AV(TOT), $AV(1/TOT)$	Groupie Groupie
no ionger allowed	= 1 = 2 BAND. CONSERVE AV(101), AV(1/101) 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 EACH ISOTOPE INDEPENDENTLY. IF THE SELECTOR	Groupie
	EACH ISOTORE INDERENDENTLY IT THE SELECTOR	
		Groupie
	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF	Groupie
	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR	Groupie Groupie
1 34-44 т11	IS NEGATIVE $(-1 \text{ TO } -5)$ THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES.	Groupie Groupie Groupie
1 34-44 I11	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY	Groupie Groupie Groupie Groupie
1 34-44 I11	IS NEGATIVE $(-1 \text{ TO } -5)$ THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES.	Groupie Groupie Groupie Groupie Groupie
1 34-44 I11	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E).	Groupie Groupie Groupie Groupie
1 34-44 I11	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant)	Groupie Groupie Groupie Groupie Groupie
1 34-44 I11	IS NEGATIVE $(-1 \text{ TO } -5)$ THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies	Groupie Groupie Groupie Groupie Groupie Groupie
1 34-44 I11	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV	Groupie Groupie Groupie Groupie Groupie Groupie Groupie
1 34-44 I11	IS NEGATIVE $(-1 \text{ TO } -5)$ THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV 1/E - 0.1 EV TO 67 KEV	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
-	IS NEGATIVE $(-1 \text{ TO } -5)$ THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV 1/E - 0.1 EV TO 67 KEV FISSION - ABOVE 67 KEV	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
-	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV 1/E - 0.1 EV TO 67 KEV FISSION - ABOVE 67 KEV = > 1 - Read input table -ADDED OPTION TO ALLOW TEMPERATURE OF THE MAXWELLIAN TO BE CHANGED - SEE INPUT LINE 4,	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
-	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV 1/E - 0.1 EV TO 67 KEV FISSION - ABOVE 67 KEV = > 1 - Read input table -ADDED OPTION TO ALLOW TEMPERATURE OF THE MAXWELLIAN TO BE CHANGED - SEE INPUT LINE 4, COLUMNS 55 - 66.	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
-	IS NEGATIVE $(-1 \text{ TO } -5)$ THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV 1/E - 0.1 EV TO 67 KEV FISSION - ABOVE 67 KEV = > 1 - Read input table -ADDED OPTION TO ALLOW TEMPERATURE OF THE MAXWELLIAN TO BE CHANGED - SEE INPUT LINE 4, COLUMNS 55 - 66. = -1 - 1/E	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
-	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV 1/E - 0.1 EV TO 67 KEV FISSION - ABOVE 67 KEV = > 1 - Read input table -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	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
-	IS NEGATIVE $(-1 \text{ TO } -5)$ THE SAME NUMBER OF BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR ALL ISOTOPES. NUMBER OF POINTS USED TO DESCRIBE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E). = 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 EV 1/E - 0.1 EV TO 67 KEV FISSION - ABOVE 67 KEV = > 1 - Read input table -ADDED OPTION TO ALLOW TEMPERATURE OF THE MAXWELLIAN TO BE CHANGED - SEE INPUT LINE 4, COLUMNS 55 - 66. = -1 - 1/E	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie

				NO IINTE EO EUE NUMPER OF DOINES	<b>0</b>
				NO LIMIT TO THE NUMBER OF POINTS USED TO DESCRIBE WEIGHTING.	Groupie 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. = .LT. 0.0001 - USE STANDARD 0.001	Groupie
				(0.1 PER-CENT)	Groupie 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 alllowed = mulriples 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 INPUT VALUES MUST ALL BE,	Groupie 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 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
	1	67-70	I4	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. $(abardard ENDE/D)$	Groupie
				= 0 = cross section = 0 (standard ENDF/B) = 1 = cross section = constant (equal to	Groupie
					Groupie Groupie
:	2-4	1-66 6	E11.4	value at highest tabulated energy). SIGMA-0 Definition	Groupie Groupie
	2-4			value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0	Groupie Groupie Groupie
	2-4			value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region	Groupie Groupie Groupie Groupie
:	2-4			value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation	Groupie Groupie Groupie Groupie Groupie
:	2-4			<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT</pre>	Groupie Groupie Groupie Groupie Groupie Groupie
	2-4			<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0,</pre>	Groupie Groupie Groupie Groupie Groupie Groupie
	2-4			<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT</pre>	Groupie Groupie Groupie Groupie Groupie Groupie
:				<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE.</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie
				<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2	1-72	A72	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPLE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN)</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3	1-72 1-72	A72 A72	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT)</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE F	1-72 1-72 DURTH IN	A72 A72 IPUT CARI	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT)</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE F EACH (	1-72 1-72 DURTH IN DUTPUT D	A72 A72 IPUT CARI DEVICE M2	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE F EACH ( THERE)	1-72 1-72 DURTH IN DUTPUT D FORE EAC	A72 A72 IPUT CARI EVICE MA CH OF THE	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT)</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE F EACH ( THERE)	1-72 1-72 DURTH IN DUTPUT D FORE EAC	A72 A72 IPUT CARI EVICE MA CH OF THE	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPLE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE F EACH ( THERE)	1-72 1-72 DURTH IN DUTPUT D FORE EAC	A72 A72 IPUT CARI EVICE MA CH OF THE	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPLE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE FC EACH ( THERE) ZERO (	1-72 1-72 DURTH IN DUTPUT D FORE EAC FO INDIC	A72 A72 IPUT CARI IEVICE M2 IH OF THE IATE NO (	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE F EACH ( THERE ZERO ( 2 ZERO (	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11	A72 A72 PUT CARL EVICE MA CH OF THE CATE NO C I11	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE FO EACH ( THERE) ZERO ? 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22	A72 A72 IPUT CARL EVICE MA EVICE MA CH OF THE CATE NO C II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETER LISTING</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE FO EACH O THERE ZERO 9 4 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22 23-33	A72 A72 PUT CARI EVICE MA H OF THE CATE NO C II1 II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETERS COMPUTER READABLE</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE FO EACH ( THERE) ZERO ? 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22	A72 A72 IPUT CARL EVICE MA EVICE MA CH OF THE CATE NO C II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPLE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETERS LISTING MULTI-BAND PARAMETERS COMPUTER READABLE UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT</pre>	Groupie Groupie
	2 3 THE FO EACH O THERE ZERO 9 4 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22 23-33	A72 A72 PUT CARI EVICE MA H OF THE CATE NO C II1 II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETERS COMPUTER READABLE</pre>	Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie Groupie
	2 3 THE FO EACH O THERE ZERO 9 4 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22 23-33	A72 A72 PUT CARI EVICE MA H OF THE CATE NO C II1 II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETERS COMPUTER READABLE UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT = 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1)</pre>	Groupie Groupie
	2 3 THE F( THERE ZERO 2 4 4 4 4 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22 23-33 34-44	A72 A72 IPUT CARI EVICE M2 IH OF THE IATE NO C II1 II1 II1 II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETERS COMPUTER READABLE UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT = 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1) = 2 - LINEAR-LINEAR (INTERPOLATION LAW 2) UNSHIELDED CROSS SECTIONS = 1 - CROSS SECTIONS = 1 - CROSS SECTIONS</pre>	Groupie Groupie
	2 3 FHE F EACH O THERE ZERO 7 4 4 4 4 4 4 4 4 4 4 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22 23-33 34-44 45-55	A72 A72 PPUT CARL EVICE MA CH OF THE CATE NO C II11 II11 II11 II11 II11 II11	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETERS COMPUTER READABLE UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT = 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1) = 2 - LINEAR-LINEAR (INTERPOLATION LAW 2) UNSHIELDED CROSS SECTIONS = 2 - RESONANCE INTEGRALS = 1 - CROSS SECTIONS</pre>	Groupie Groupie
	2 3 THE FC EACH 0 THERE ZERO 9 4 4 4 4 4 4 4 4 4 4 1/20	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22 23-33 34-44 45-55 - ADDED	A72 A72 PUT CARL EVICE MA CH OF THE CATE NO C II1 II1 II1 II1 II1 II1 II1 II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETERS COMPUTER READABLE UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT = 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1) = 2 - LINEAR-LINEAR (INTERPOLATION LAW 2) UNSHIELDED CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-DEARD FORMAT (INTERPOLATION LAW 2) UNSHIELDED CROSS SECTIONS LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-DEARD FORMAT (INTERPOLATION LAW 2) UNSHIELDED CROSS SECTIONS LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS W OPTION</pre>	Groupie Groupie
	2 3 FHE F EACH O THERE ZERO 7 4 4 4 4 4 4 4 4 4 4 4	1-72 1-72 DURTH IN DUTPUT D FORE EAC TO INDIC 1-11 12-22 23-33 34-44 45-55	A72 A72 PUT CARL EVICE MA CH OF THE CATE NO C II1 II1 II1 II1 II1 II1 II1 II1 II1	<pre>value at highest tabulated energy). SIGMA-0 Definition -2019/6/23 - Only the GROUPIE standard = 0 in allowed for Unresolved Resonance Region Self-Shielding calculation IF SIGMA-0 DEFINITION SELECTOR &lt; 0, THE NEXT 4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0, 6 PER LINE. ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME (STANDARD OPTION = ENDFB.OUT) O IS USED TO SELECT ALL DESIRED OUTPUT MODES. AY BE TURNED OFF (0) OR ON (1). THEREFORE E FOLLOWING INPUT PARAMETERS MAY BE EITHER DUTPUT OR NON-ZERO TO INDICATE OUTPUT. SELF-SHIELDED CROSS SECTION LISTING = 1 - CROSS SECTIONS = 2 - RESONANCE INTEGRALS MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETERS COMPUTER READABLE UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT = 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1) = 2 - LINEAR-LINEAR (INTERPOLATION LAW 2) UNSHIELDED CROSS SECTIONS = 2 - RESONANCE INTEGRALS = 1 - CROSS SECTIONS</pre>	Groupie Groupie

6-N 1-6 16 LOWER MAT OR ZA LIMIT Grou 7-8 12 LOWER MF LIMIT Grou 9-11 13 LOWER MT LIMIT Grou 12-17 111 UPPER MAT OR ZA LIMIT Grou 18-19 12 UPPER MT LIMIT Grou 20-22 I3 UPPER MT LIMIT Grou UP TO 100 RANGES MAY BE SPECIFIED, ONE RANGE Grou PER LINE. THE LIST OF RANGES IS TERMINATED Grou BY A BLANK CARD. IF THE UPPER MAT OR ZA Grou LIMIT IS LESS THAN THE LOWER LIMIT. IF THE UPPER Grou IS SET EQUAL TO THE LOWER LIMIT. IF THE UPPER Grou MF OR MT LIMIT IS ZERO IT WILL BE SET EQUAL Grou IF THE FIRST REQUEST LINE IS BLANK IT WILL Grou TERMINATE THE LIST OF REQUESTS AND CAUSE ALL Grou DATA TO BE RETRIEVED (SEE EXAMPLE INPUT). Grou Grou VARY 1-66 6E11.4 ENERGY GROUP BOUNDARIES. ONLY REQUIRED IF THE NUMBER OF GROUPS INDICATED ON THE FIRST Grou NH1 BOUNDARIES WILL BE READ FROM THE GROUPS GROU NH1 BOUNDARIES WILL BE READ FROM THE GROU NH1 BOUNDARIES WILL BE READ FROM THE GROU WILL BE READ FROM THE INPUT FILE. Grou VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY REQUIRED IF THE NUMBER OF POINTS INDICATED ON THE FIRST GROU VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY WILL BE READ FROM THE INPUT FILE. GROUPS VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY REQUIRED IF THE NUMBER OF POINTS INDICATED GROUPS VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY WILL BE READ FROM THE INPUT FILE. GROUPS WILL BE READ FROM THE INPUT FILE. GROUPS WILL BE READ FROM THE INPUT FILE. GROUPS WILL BE READ FROM THE INPUT FILE. GROUPS VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY WILL BE READ FROM THE INPUT FILE. GROUPS WILL BE READ FROM THE IN	upie upie upie upie upie upie upie upie
THE NUMBER OF GROUPS INDICATED ON THE FIRST Grou INPUT CARD IS POSITIVE. ALL ENERGIES MUST Grou BE IN ASCENDING ENERGY IN EV. THE PRESENT Grou LIMITS ARE 1 TO 20,000 GROUPS. FOR N GROUPS Grou N+1 BOUNDARIES WILL BE READ FROM THE Grou INPUT FILE, E.G. IF THE FIRST INPUT CARD Grou INDICATES 20 GROUPS, 21 ENERGY BOUNDARIES Grou WILL BE READ FROM THE INPUT FILE. Grou Grou VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY Grou	npie npie npie npie npie npie npie npie
VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY Grou	ipie ipie ipie ipie ipie ipie ipie
ON FIRST CARD IS MORE THAN ONE. DATA IS Grou GIVEN IN (ENERGY, WEIGHT) PAIRS, UP TO 3 Grou PAIRS PER CARD, USING ANY NUMBER OF CARDS Grou REQUIRED. ENERGIES MUST BE IN ASCENDING Grou ORDER IN EV. THE SPECTRUM VALUES MUST BE Grou NON-NEGATIVE. THE ENERGY RANGE OF SPECTRUM Grou MUST AT LEAST SPAN THE ENERGY RANGE OF THE Grou ENERGY GROUPS. SINCE SPECTRUM IS STORED IN Grou PAGING SYSTEM THERE IS NO LIMIT TO NUMBER Grou OF POINTS THAT CAN BE USED TO DESCRIBE THE Grou WEIGHTING SPECTRUM. Grou	npie npie npie npie npie npie npie npie
Grou EXAMPLE INPUT NO. 1 Grou	-
Grou REQUEST DATA BY MAT AND PROCESS ALL DATA (ALL MAT BETWEEN 1 AND 9999). USE THE TART 175 GROUP STRUCTURE, GENERATE 2 BAND PARAMETERS (THE FOR ALL ISOTOPES) TO 0.1 PER-CENT ACCURACY IN THE SELF-SHIELDING CURVE. OUTPUT ALL LISTING, COMPUTER READABLE AND ENDF/B FORMAT GROUP AVERAGES. Grou EXPLICITLY SPECIFY THE STANDARD FILENAMES. Grou	ipie ipie ipie ipie ipie ipie ipie
Grou THE FOLLOWING 7 INPUT LINES ARE REQUIRED. Grou	pie
Group         Group <th< td=""><td>ipie ipie ipie</td></th<>	ipie ipie ipie

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 THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ Groupie FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND Groupie WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238 Groupie Groupie THE FOLLOWING 7 INPUT LINES ARE REQUIRED. Groupie Groupie 0 1.00000-03 Groupie 0 0 -2 0 \ENDFB6\SIGMA1\K300\ZA092238 Groupie \ENDFB6\GROUPIE\K300\ZA092238 Groupie 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. 3 Groupie Groupie PROCESS ALL DATA. USE 1/E WEIGHTING IN ORDER TO CALCULATE Groupie UNSHIELDED ONE GROUP CROSS SECTIONS OVER THE ENERGY RANGE 0.5 EV Groupie TO 1 MEV (NOTE THAT THE RESULTS ARE SIMPLY PROPORTIONAL TO THE Groupie RESONANCE INTEGRAL FOR EACH REACTION). OUTPUT UNSHIELDED LISTING. Groupie Groupie LEAVE THE DEFINITION OF THE FILENAMES BLANK - THE PROGRAM WILL Groupie THEN USE STANDARD FILENAMES. Groupie Groupie THE FOLLOWING 7 INPUT CARDS ARE REQUIRED. Groupie Groupie ٥ ٥ 1 -1 0 Groupie (USE STANDARD FILENAME = ENDFB.IN) Groupie (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 EXAMPLE INPUT NO. 4 Groupie \_\_\_\_\_ Groupie THIS EXAMPLE USES A USER DEFINED GROUP STRUCTURE AND WEIGHTING Groupie FUNCTION - THESE ARE NOT REALISTIC IN TERMS OF ACTUAL ENERGIES Groupie AND WEIGHTS - THEY ARE ONLY INTENDED TO ILLUSTRATE THE ORDER OF Groupie THE INPUT PARAMETERS. Groupie Groupie 0 11 0 6 0 Groupie RECENT.OUT Groupie GROUPIE.OUT Groupie 1 1 1 1 1 Groupie Example with users defined groupus and spectrum weighting Groupie 1 1 1 999999999 Groupie (blabk line terminates request list) Groupie 1.00000-05 1.00000-04 1.00000-03 1.00000-02 1.00000-01 1.00000+00 grou Groupie 1.00000+01 1.00000+02 1.00000+03 1.00000+04 1.00000+05 1.00000+06 grou Groupie 1.00000-02 0.1 1.00000+00 0.01 1.00000-05 1.0 weig Groupie 1.00000+02 0.001 1.00000+04 0.0001 1.00000+06 0.000001 weig Groupie Groupie