				GROUP
PROGRAM				GROUP
	-====	==		GROUP
		(NOVEMBER 1976		GROUP
VERSION	79-1	(OCTOBER 1979)	CDC-7600 AND CRAY-1 VERSION.	GROUP
VERSION	80-1	(MAY 1980) IBM	I, CDC AND CRAY VERSION	GROUP
VERSION	81-1	(JANUARY 1981)	EXTENSION TO 3000 GROUPS	GROUP
VERSION	81-2	(MARCH 1981) I	MPROVED SPEED	GROUP
			BUILT-IN 1/E WEIGHTING SPECTRUM	GROUP
			IMPROVED COMPUTER COMPATIBILITY	GROUP
			*MAJOR RE-DESIGN.	GROUP
VERSION	03-T	(JANOARI 1985)	*ELIMINATED COMPUTER DEPENDENT CODING.	
			*NEW, MORE COMPATIBLE I/O UNIT NUMBERS	
		(*NEW MULTI-BAND LIBRARY BINARY FORMAT.	
VERSION	83-2	(OCTOBER 1983)	ADDED OPTION TO ALLOW SIGMA-0 TO BE	GROUP
			DEFINED EITHER AS MULTIPLES OF	GROUP
			UNSHIELDED TOTAL CROSS SECTION IN EACH	
			GROUP, OR POWERS OF 10 IN ALL GROUPS.	GROUP
VERSION	84-1	(APRIL 1984)	ADDED MORE BUILT IN MULTIGROUP ENERGY	GROUP
			STRUCTURES.	GROUP
VERSION	85-1	(APRIL 1985)	*UPDATED FOR ENDF/B-VI FORMATS.	GROUP
			*SPECIAL I/O ROUTINES TO GUARANTEE	GROUP
			ACCURACY OF ENERGY.	GROUP
			*DOUBLE PRECISION TREATMENT OF ENERGY	
			(REQUIRED FOR NARROW RESONANCES).	GROUP
			*MINIMUM TOTAL CROSS SECTION TREATMENT	
	0F 0	(1005)		
			*FORTRAN-77/H VERSION	GROUP
			*ENDF/B-VI FORMAT	GROUP
VERSION	86-2	(JUNE 1986)	*BUILT-IN MAXWELLIAN, 1/E AND FISSION	GROUP
			WEIGHTING SPECTRUM.	GROUP
VERSION	88-1	(JULY 1988)	*OPTIONINTERNALLY DEFINE ALL I/O	GROUP
			FILE NAMES (SEE, SUBROUTINES FILIO1	GROUP
			FILIO2 FOR DETAILS).	GROUP
			*IMPROVED BASED ON USER COMMENTS.	GROUP
VERSTON	89-1	(JANUARY 1989)	*PSYCHOANALYZED BY PROGRAM FREUD TO	GROUP
1210201	0, 1	(011101111 1909)	INSURE PROGRAM WILL NOT DO ANYTHING	GROUP
			CRAZY.	GROUP
			*UPDATED TO USE NEW PROGRAM CONVERT	GROUP
			KEYWORDS.	GROUP
			*ADDED LIVERMORE CIVIC COMPILER	GROUP
			CONVENTIONS.	GROUP
VERSION	91-1	(JUNE 1991)	*INCREASED PAGE SIZE FROM 1002 TO 5010	
			POINTS	GROUP
			*UPDATED BASED ON USER COMMENTS	GROUP
			*ADDED FORTRAN SAVE OPTION	GROUP
			*COMPLETELY CONSISTENT ROUTINE TO READ	GROUP
			FLOATING POINT NUMBERS.	GROUP
VERSION	92-1	(JANUARY 1992)	*ADDED RESONANCE INTEGRAL CALCULATION -	
			UNSHIELDED AND/OR SHIELDED - FOR	GROUP
			DETAILS SEE BELOW	GROUP
			*INCREASED NUMBER OF ENERGY POINTS	GROUP
			IN BUILT-IN SPECTRA - TO IMPROVE	GROUP
			ACCURACY.	GROUP
			*ALLOW SELECTION OF ZA/MF/MT OR	GROUP
			MAT/MF/MT RANGES - ALL DATA NOT	GROUP
			SELECTED IS SKIPPED ON INPUT AND	GROUP
			NOT WRITTEN AS OUTPUT.	GROUP
			*COMPLETELY CONSISTENT I/O ROUTINES -	GROUP
			TO MINIMIZE COMPUTER DEPENDENCE.	GROUP
			*NOTE, CHANGES IN INPUT PARAMETER	GROUP
			FORMAT - FOR ZA/MF/MT OR MAT/MF/MT	GROUP
			RANGES.	GROUP
VERSION	92-2	(JUNE 1992)	*MULTIBAND PARAMETERS OUTOUT AS	GROUP
A RUCE TON	92-Z	(00ME 1992)		
	02 1	(ADDIT 1000)	CHARACTER (RATHER THAN BINARY) FILE.	
VERSTON	93-1	(APRIL 1993)	*INCREASED PAGE SIZE FROM 5010 TO	GROUP
1010101			30000 POINTS	GROUP
12102101			*ELIMINATED COMPUTER DEPENDENCE.	GROUP
	94-1	(JANUARY 1994)	*VARIABLE ENDF/B DATA FILENAMES	GROUP
	94-1	(JANUARY 1994)	*VARIABLE ENDF/B DATA FILENAMES TO ALLOW ACCESS TO FILE STRUCTURES	GROUP GROUP

		HAS BEEN CHANGED)	GROUPIE
		*CLOSE ALL FILES BEFORE TERMINATING	GROUPIE
		(SEE, SUBROUTINE ENDIT)	GROUPIE
VERSION 95-1	(JANUARY 1994)	*CORRECTED MAXWELLIAN WEIGHTING	GROUPIE
		*CHANGING WEIGHTING SPECTRUM FROM	GROUPIE
		0.1 TO 0.001 % UNCERTAINTY	GROUPIE
VERSION 96-1	(JANUARY 1996)	*COMPLETE RE-WRITE	GROUPIE
		*IMPROVED COMPUTER 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	GROUPIE
		(PREVIOUSLY 100)	GROUPIE
		*CORRECTED USE OF MAXWELLIAN + 1/E + FISSION SPECTRUM	GROUPIE GROUPIE
		*ONLY 2 BAND VERSION DISTRIBUTED	GROUPIE
		(CONTACT AUTHOR FOR DETAILS)	GROUPIE
		*DEFINED SCRATCH FILE NAMES	GROUPIE
VERSION 99-1	(MARCH 1999)	*CORRECTED CHARACTER TO FLOATING	GROUPIE
		POINT READ FOR MORE DIGITS	GROUPIE
		*UPDATED TEST FOR ENDF/B FORMAT	GROUPIE
		VERSION BASED ON RECENT FORMAT CHANG	
		*GENERAL IMPROVEMENTS BASED ON	GROUPIE
VERSION 99-2	(TITNE 1000)	USER FEEDBACK *ASSUME ENDF/B-VI, NOT V, IF MISSING	GROUPIE
VERSION 33 Z	(UONE 1999)	MF=1, MT-451.	GROUPIE
VERS. 2000-1	(FEBRUARY 2000))*ADDED MF=10, ACTIVATION CROSS SECTIO	
		PROCESSING.	GROUPIE
		*GENERAL IMPROVEMENTS BASED ON	GROUPIE
		USER FEEDBACK	GROUPIE
VERS. 2002-1	(FEBRUARY 2002	2) *ADDED TART 700 GROUP STRUCTURE	GROUPIE
	(10312 0000)	*ADDED VARIABLE SIGMA0 INPUT OPTION	GROUPIE
	(MAY 2002) (NOV. 2002)	*OPTIONAL INPUT PARAMETERS *ADDED SAND-II EXTENDED DOWN TO	GROUPIE GROUPIE
	(1007. 2002)	1.0D-5 EV.	GROUPIE
	(JUNE 2003)	*CORRECTED SAND-II 620 AND 640 GROUP	
		ENERGY BOUNDARIES DEFINITIONS.	GROUPIE
VERS. 2004-1	(SEPT. 2004)	*INCREASED PAGE SIZE FROM 30000 TO	GROUPIE
		120000 POINTS	GROUPIE
		*ADDED "OTHER" AS ADDITIONAL REACTION	
		TO IMPROVE MULTI-BAND FITTING *ADDED ITERATION FOR "BEST" PARTIAL	GROUPIE 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	
	(TAN 0007)	BUILT-IN STANDARD SPECTRUM.	GROUPIE
VERS. 2007-1	(UAIN. 2007)	*CHECKED AGAINST ALL ENDF/B-VII. *INCREASED PAGE SIZE FROM 120,000 TO	GROUPIE GROUPIE
		600,000 POINTS	GROUPIE
VERS. 2008-1	(JAN. 2008)	*72 CHARACTER FILE NAMES.	GROUPIE
	•	*GENERAL UPDATES	GROUPIE
VERS. 2010-1	(Apr. 2010)	*INCREASED WEIGHTING SPECTRUM TO 30,00	OGROUPIE
		FROM 3,000 ENERGY POINTS.	GROUPIE
		*ADDED OUTPUT TO PLOT/COMPARE SHIELDED	
VEDC 2011 1	(Tupo 2011)	AND UNSHIELDED CROSS SECTIONS.	GROUPIE
VERS. 2011-1	(June 2011)	*Corrected TART 700 groups to extend up to 1 GeV (1,000 MeV) - previously it	-
		was ERRONEOUSLY cutoff at 20 MeV.	GROUPIE
VERS. 2011-2	(Nov. 2011)	*Corrected TART 616 groups lowest	GROUPIE
	-	energy from 1.0D-4 eV to 1.0D-5 eV.	GROUPIE
		*Added TART 666 to 200 MeV (for TENDL)	
		*Optional high energy cross section	GROUPIE

				extension above tabulated energy range	GROUPIE
				(either = 0 = standard, or constant)	GROUPIE
				WARNING - ENDF/B standard convention is that the cross section = 0 where it	GROUPIE
				is not explicitly defined - extension	
				= 0 is standard, constant is NOT, so	GROUPIE
				constant extension is NOT RECOMMENDED.	GROUPIE
VERS.	2012-1	(Aug.	2012)	*Added CODENAME	GROUPIE
				*32 and 64 bit Compatible	GROUPIE
VERS	2013-1	(Now	2013)	*Added ERROR stop. *Extended OUT9.	GROUPIE GROUPIE
VERO.	2015 1	(1000.	2013)	*Uses OUTG, not OUT10 for energies.	GROUPIE
VERS.	2015-1	(Jan.	2015)	*Corrected SPECTM - handle ALL included	
				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. *Extended OUTG	GROUPIE GROUPIE
				*Replaced ALL 3 way IF Statements.	GROUPIE
				*Generalized TART Group Strructures.	GROUPIE
				*Generalized SAND-II Group Structures.	GROUPIE
				*Extended SAND-II to 60, 150, 200 MeV.	GROUPIE
VERS.	2015-2	(Mar.	2015)	*Deleted 1P from formats reading input	
				parameters, causing incorrect scaling *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	
				listings simpler.	GROUPIE
				*Corrected High Energy Extension = Can effect highest energy group.	GROUPIE GROUPIE
VERS.	2016-1	(Julv	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 explcitly handle small shielding	GROUPIE GROUPIE
				limit - without this update the small	
				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
				<pre>*Default multi-band is method #2 = conserve <x>, <1/(x+<x>>, <1/x>.</x></x></pre>	GROUPIE GROUPIE
				*Definition of built-in group structure	
				using SUBROUTINE GROPE is identical	GROUPIE
				for GROUPIE and VIRGIN.	GROUPIE
				*All floating input parameters changed	
				to character input + IN9 conversion.	GROUPIE
				*Output report identfies MF now that this code does more than just MF=3.	GROUPIE 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
VEDO	2010 1	(201 01	comments) = for multi-band output	GROUPIE
	2018-1 2019-1			*Added on-line output for ALL ENDERROR *Major re-write to re-order output to	GROUPIE
		(Same	/	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 GROUPTE results were in memory - saving ALL GROUPIE greatly simplifies the logic. GROUPIE GROUPTE *Additional Interpolation Law Tests *Check maximum Tabulated Energy of MTs GROUPIE to insure they ALL end at the same GROUPIE energy. GROUPIE *Multi-band = 1 no longer allowed. GROUPIE The only allowed values are, GROUPIE 0 = no multi-band calculations, or, GROUPTE 2 = Conserve 1/[total + <total>] GROUPTE *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, GROUPTE 1) LSSF = 0 = Output cross sections GROUPIE 2) INTUNR = 2 = Interpolation law GROUPTE *Added ZAzzzaaa to filenames. GROUPIE VERS. 2020-1 (Aug. 2020) *Major re-write to update for new URR GROUPIE self-shielding, MF/MT=2/152 and 2/153.GROUPIE *Corrected BOTH ends of unresolved GROUPTE for MF/MT=2/152 and 2/153 output. GROUPIE *Unresolved extrapolation ONLY to GROUPIE groups completely inside the URR + GROUPTE per ends for MF/MT=2/152 & 153 output.GROUPIE *Small shielding < 0.1 % = accuracy GROUPIE of reconstructed data. GROUPIE *Forced no self-shielding at upper end GROUPIE GROUPTE of unresolved = match high energy tabulated. GROUPIE *Corrected PLOTTAB output if no URR GROUPTE fit - it was outputting EMPTY tables GROUPIE for original and fit moments, which GROUPIE in this case did not exist. GROUPIE *Only 2 band, Method#2 [sigt + <sigt>] GROUPIE alloed for multi-band calculation. GROUPIE *WARNING - if input Requested MF range GROUPIE prevents unresolved region calculationGROUPIE *Added Target Isomer Flag GROUPIE *Correct MULTBAND.LST output format. GROUPIE VERS. 2021-1 (Jan. 2021) *Updated for FORTRAN 2018 GROUPIE GROUPIE 2020-1 Acknowledgment GROUPTE GROUPTE I Thank Jean-Christophe Sublet (NDS, IAEA, Vienna, Austria) for GROUPTE reporting the ERROR in GROUPIE (2019-1) that led to the update in GROUPIE GROUPIE (2020-1) to correctly define the PLOTTAB output, whether GROUPIE or not Unresolved Resonance Region (URR) fit was performed. 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. GROUPTE GROUPIE I thank Jean-Christophe Sublet (UKAEA) for contributing MAC GROUPIE executables and Bojan Zefran (IJS, Slovenia) for contributing GROUPTE LINUX (32 or 63 bit) executables. And most of all I must thank GROUPTE 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 GROUPIE OWNED, MAINTAINED AND DISTRIBUTED BY 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 -----GROUPTE -------GROUPIE Dermott E. Cullen GROUPIE PRESENT CONTACT INFORMATION GROUPIE -------GROUPIE Dermott E. Cullen GROUPIE 1466 Hudson Way GROUPTE Livermore, CA 94550 GROUPIE U.S.A. GROUPIE Telephone 925-443-1911 GROUPTE RedCullen1@Comcast.net GROUPIE E. Mail RedCullen1.net/HOMEPAGE.NEW GROUPIE Website GROUPIE AUTHORS MESSAGE GROUPIE GROUPTE THE REPORT DESCRIBED ABOVE IS THE LATEST PUBLISHED DOCUMENTATION GROUPIE FOR THIS PROGRAM. HOWEVER, THE COMMENTS BELOW SHOULD BE CONSIDEREDGROUPIE THE LATEST DOCUMENTATION INCLUDING ALL RECENT IMPROVEMENTS. PLEASEGROUPIE READ ALL OF THESE COMMENTS BEFORE IMPLEMENTATION, PARTICULARLY GROUPIE THE COMMENTS CONCERNING MACHINE DEPENDENT CODING. GROUPTE GROUPTE AT THE PRESENT TIME WE ARE ATTEMPTING TO DEVELOP A SET OF COMPUTERGROUPIE INDEPENDENT PROGRAMS THAT CAN EASILY BE IMPLEMENTED ON ANY ONE GROUPIE OF A WIDE VARIETY OF COMPUTERS. IN ORDER TO ASSIST IN THIS PROJECTGROUPIE 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 GROUPTE 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 GROUPTE (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. GROUPTE GROUPTE GROUPTE ENDF/B FORMAT 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 GROUPTE ASSUMED THAT THE MAT, MF AND MT ON EACH CARD IS CORRECT. SEQUENCE GROUPLE NUMBERS (COLUMNS 76-80) ARE IGNORED ON INPUT, BUT WILL BE GROUPTE 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. GROUPTE 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 MADEGROUPIE LINEARLY INTERPOLABLE USING PROGRAM LINEAR (UCRL-50400, VOL. 17, GROUPIE PART A). THE RESONANCE CONTRIBUTION MAY BE ADDED TO THE BACKGROUNDGROUPIE CROSS SECTIONS USING PROGRAM RECENT (UCRL-50400, VOL. 17, PART B).GROUPIE

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	GROUPIE GROUPIE
	GROUPIE
IF ENDF/B FORMATTED OUTPUT IS REQUESTED ENTIRE EVALUATIONS ARE	GROUPIE
	GROUPIE
	GROUPIE
	GROUPIE GROUPIE
	GROUPIE
THE FACT THAT THIS PROGRAM HAS OPERATED ON THE DATA IS DOCUMENTED	
	GROUPIE
HOLLERITH SECTION TO DESCRIBE THE GROUP STRUCTURE AND WEIGHTING	GROUPIE
SPECTRUM, E.G.	GROUPIE
	GROUPIE
	GROUPIE
	GROUPIE
	GROUPIE
THE ORDER OF ALL SIMILAR COMMENTS (FROM LINEAR, RECENT AND SIGMA1)	
	GROUPIE
THE DATA.	GROUPIE
	GROUPIE
THESE COMMENT CARDS ARE ONLY ADDED TO EXISTING HOLLERITH SECTIONS,	
I.E., THIS PROGRAM WILL NOT CREATE A HOLLERITH SECTION. THE FORMAT	
·	GROUPIE
	GROUPIE
	GROUPIE
	GROUPIE
·	GROUPIE
AS SUCH IT IS IMPOSSIBLE FOR THE PROGRAM TO DETERMINE WHAT FORMAT	GROUPIE
SHOULD BE USED TO CREATE A HOLLERITH SECTION.	GROUPIE
	GROUPIE
THIS CONVENTION HAS BEEN ADOPTED BECAUSE MOST USERS DO NOT	GROUPIE
REQUIRE A CORRECT REACTION INDEX FOR THEIR APPLICATIONS AND IT WAS	
NOT CONSIDERED WORTHWHILE TO INCLUDE THE OVERHEAD OF CONSTRUCTING	
A CORRECT REACTION INDEX IN THIS PROGRAM. HOWEVER, IF YOU REQUIRE A REACTION INDEX FOR YOUR APPLICATIONS, AFTER RUNNING THIS PROGRAM	
	GROUPIE
	GROUPIE
	GROUPIE
	GROUPIE
SINCE THIS PROGRAM USES A LOGICAL PAGING SYSTEM THERE IS NO LIMIT	GROUPIE
, , ,	GROUPIE
	GROUPIE
	GROUPIE
THE PROGRAM SELECTS MATERIALS TO BE PROCESSED BASED EITHER ON	
	GROUPIE GROUPIE
MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR	GROUPIE GROUPIE GROUPIE
MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS	GROUPIE GROUPIE GROUPIE GROUPIE
MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS USED TO SELECT MATERIALS, AND WILL TERMINATE WHEN A MAT OR ZA	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS USED TO SELECT MATERIALS, AND WILL TERMINATE WHEN A MAT OR ZA IS FOUND THAT IS ABOVE THE RANGE OF ALL REQUESTS.	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
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MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS USED TO SELECT MATERIALS, AND WILL TERMINATE WHEN A MAT OR ZA IS FOUND THAT IS ABOVE THE RANGE OF ALL REQUESTS. ENERGY ORDER AND UNITS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS USED TO SELECT MATERIALS, AND WILL TERMINATE WHEN A MAT OR ZA IS FOUND THAT IS ABOVE THE RANGE OF ALL REQUESTS. ENERGY ORDER AND UNITS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
MAT (ENDF/B MAT NO.) OR ZA. THE PROGRAM ALLOWS UP TO 100 MAT OR ZA RANGES TO BE SPECIFIED. THE PROGRAM WILL ASSUME THAT THE ENDF/B TAPE IS IN EITHER MAT OR ZA ORDER, WHICHEVER CRITERIA IS USED TO SELECT MATERIALS, AND WILL TERMINATE WHEN A MAT OR ZA IS FOUND THAT IS ABOVE THE RANGE OF ALL REQUESTS. ENERGY ORDER AND UNITS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE 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 GROUPTE 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 GROUPTE THE USER MAY INPUT AN ARBITRARY GROUP STRUCTURE OR THE USER MAY GROUPTE USE USE ONE OF THE 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) GROUPTE (6) 69 GROUP (WIMS STRUCTURE) GROUPIE (7) 68 GROUP (GAM-I STRUCTURE) GROUPTE (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) GROUPTE (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) GROUPTE (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) (19)1102 GROUP (UKAEA STRUCTURE, 1.0D-5 eV, UP TO 1 GeV) GROUPIE GROUPIE GROUPIE GROUPTE GROUP AVERAGES 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 (INTEGRAL E1 TO E2) (S(E)*WT(E)*DE) GROUPIE WHERE . . . GROUPTE GROUPIE AVERAGE = GROUP AVERAGED CROSS SECTION GROUPIE E1, E2 = ENERGY LIMITS OF THE GROUP GROUPIE SIGMA(E) = ENERGY DEPENDENT CROSS SECTION FOR ANY GIVEN REACTION GROUPIE S(E) = ENERGY DEPENDENT WEIGHTING SPECTRUM GROUPTE WT(E) = ENERGY DEPENDENT SELF-SHIELDING FACTOR. GROUPTE GROUPTE 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 REPRESENTATIONGROUPIE OF THE SPECTRUM BETWEEN TABULATED DATA POINTS. GROUPIE GROUPIE THE PRESENT VERSION OF THE CODE HAS THREE BUILT-IN WEIGHTING GROUPIE SPECTRA. GROUPTE GROUPIE (1) CONSTANT GROUPIE GROUPIE (2) 1/E(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 = $C2 \times EXP(-E/WA) \times SINH(SQRT(E \times WB))$ (67 KeV, 10 MeV) GROUPIE CONSTANT = Equal to Fission at 10 MeV (above 10 MeV) GROUPIE

GROUPIE = 0.253 EV (293 KELVIN) GROUPIE KT WA = 9.65D+5GROUPIE = 2.29D-6GROUPIE WB C1, C2 = DEFINED TO MAKE SPECTRUM CONTINUOUS GROUPIE GROUPIE FISSION SPECTRUM CONSTANTS FROM GROUPTE 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. GROUPTE GROUPTE 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 GROUPTE FOR THE TOTAL, ELASTIC, CAPTURE AND FISSION THE PROGRAM USES A GROUPIE WEIGHTING FUNCTION THAT IS A PRODUCT OF THE ENERGY DEPENDENT GROUPTE WEIGHTING SPECTRUM TIMES A BONDERENKO TYPE SELF-SHIELDING FACTOR. GROUPIE GROUPIE WT(E) = S(E) / (TOTAL(E) + SIGMA0) **NGROUPIE GROUPTE WHERE . . . GROUPIE GROUPIE - ENERGY DEPENDENT WEIGHTING SPECTRUM (DEFINED BY GROUPTE S(E) 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). GROUPTE - CROSS SECTION TO REPRESENT THE EFFECT OF ALL OTHER STGMA0 GROUPTE MATERIALS AND LEAKAGE (DEFINED WITHIN EACH GROUP TO BE GROUPIE A MULTIPLE OF THE UNSHIELDED TOTAL CROSS SECTION WITHINGROUPIE THAT GROUP OR POWERS OF 10 - INPUT OPTION). GROUPIE N - A POSITIVE INTEGER (0, 1, 2 OR 3). GROUPIE GROUPTE THE PROGRAM WILL USE ONE ENERGY DEPENDENT WEIGHTING SPECTRUM S(E) GROUPIE AND 25 DIFFERENT BONDERENKO TYPE SELF-SHIELDING FACTORS (25 SIGMA0GROUPIE AND N COMBINATIONS) TO DEFINE 25 DIFFERENT AVERAGE CROSS SECTIONS, GROUPIE FOR EACH REACTION, WITHIN EACH GROUP. GROUPIE GROUPIE THE 25 WEIGHTING FUNCTIONS USED ARE.... GROUPIE - UNSHIELDED CROSS SECTIONS (N=0) GROUPIE (1) (2-22) - PARTIALLY SHIELDED CROSS SECTIONS (N=1 ,VARIOUS SIGMA0) GROUPIE THE VALUES OF SIGMAO USED WILL BE EITHER, GROUPTE (A) THE VALUES OF SIGMAO THAT ARE USED VARY FROM 1024 GROUPTE TIMES THE UNSHIELDED TOTAL CROSS SECTIONS IN STEPS OF 1/2 GROUPIE DOWN TO 1/1024 TIMES THE UNSHIELDED TOTAL CROSS SECTION GROUPIE (A RANGE OF OVER 1 MILLION, CENTERED ON THE UNSHIELDED GROUPIE TOTAL CROSS SECTION WITHIN EACH GROUP). GROUPIE (B) THE SAME CONSTANT VALUES OF SIGMAO 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 GROUPTE ACTUAL APPLICATIONS) GROUPIE (23) - TOTALLY SHIELDED FLUX WEIGHTED CROSS SECTION GROUPIE (N=1, SIGMA0=0)GROUPIE (24)- TOTALLY SHIELDED CURRENT WEIGHTED CROSS SECTION GROUPIE (N=2, SIGMA0=0) GROUPIE TOTALLY SHIELDED COSINE SQUARED WEIGHTED CROSS SECTION (25)GROUPIE 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 SPECTRUM S(E) TO DEFINE THE UNSHIELDED (BONDERENKO N=0) GROUPIE

AVERAGED CROSS SECTION WITHIN EACH GROUP.	GROUPIE
	GROUPIE
CALCULATION OF RESONANCE INTEGRALS	GROUPIE GROUPIE
IN A PURE ELASTIC ISOTROPICALLY SCATTERING MATERIAL WITH A	GROUPIE
CONSTANT CROSS SECTION THE SPECTRUM WILL BE 1/E AND THERE WILL	GROUPIE
BE NO SELF-SHIELDING.	GROUPIE
IN THIS CASE IF THE CROSS SECTION VARIES WITH ENERGY THE	GROUPIE GROUPIE
SPECTRUM WILL STILL BE 1/E AND THE SELF-SHIELDING FACTOR WILL	GROUPIE
BE EXACTLY 1/SIG-TOT(E) - WHERE SIG-TOT(E) = SIG-EL(E), SINCE	GROUPIE
THERE IS ONLY SCATTERING.	GROUPIE
	GROUPIE
IF WE HAVE AN INFINITELY DILUTE AMOUNT OF A MATERIAL UNIFORMLY MIXED WITH A PURE ELASTIC ISOTROPICALLY SCATTERING MATERIAL WITH	GROUPIE GROUPIE
A CONSTANT CROSS SECTION THE STANDARD DEFINITION OF THE RESONANCE	
INTEGRAL CAN BE USED TO DEFINE REACTION RATES FOR EACH REACTION.	GROUPIE
	GROUPIE
THE RESONANCE INTEGRAL IS DEFINED AS,	GROUPIE
RI = (INTEGRAL E1 TO E2) (SIGMA(E) \times S(E) \times T(E) \rightarrow DE)	GROUPIE GROUPIE
$\mathbf{x} = (\mathbf{x} + \mathbf{b} +$	GROUPIE
WHERE NORMALLY,	GROUPIE
S(E) = 1/E	GROUPIE
WT(E) = 1 - NO SELF-SHIELDING	GROUPIE
FROM THE ABOVE DEFINITION OF GROUP AVERAGED CROSS SECTIONS THE	GROUPIE GROUPIE
RESONANCE INTEGRAL IS,	GROUPIE
	GROUPIE
RI = AVERAGE * (INTEGRAL E1 TO E2) $(S(E) *WT(E) *DE)$	GROUPIE
FOR A 1/E SPECTRUM AND NO SELF-SHIELDING THIS REDUCES TO,	GROUPIE GROUPIE
FOR A 1/E SPECIRUM AND NO SELF-SHIELDING THIS REDUCES 10,	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 - ABSORPTION WILL TEND TO DECREASE THE SPECTRUM PROGRESSIVELY	GROUPIE 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 CALCULATION TO ACCOUNT FOR ABSORPTION.	GROUPIE GROUPIE
CALCULATION TO ACCOUNT FOR ADSORPTION.	GROUPIE
HOWEVER, WE WILL EXTEND THE DEFINITION OF THE RESONANCE INTEGRAL	GROUPIE
TO ACCOUNT FOR SELF-SHIELDING EFFECTS BY ALLOWING FOR INCLUSION	GROUPIE
OF SELF-SHIELDING EFFECTS IN THE DEFINITION OF GROUP AVERAGES	GROUPIE
AND THEN DEFINING THE RESONANCE INTEGRAL AS,	GROUPIE GROUPIE
RI = AVERAGE * LOG (E2/E1)	GROUPIE
	GROUPIE
IN ORDER TO CALCULATE RESONANCE INTEGRALS YOU MUST FOLLOW THESE	GROUPIE
STEPS,	GROUPIE GROUPIE
1) SELECT A 1/E SPECTRUM - ON FIRST LINE OF INPUT PARAMETERS.	GROUPIE
2) SELECT THE ENERGY BOUNDARIES - NORMALLY ONLY 1 GROUP FROM	GROUPIE
0.5 EV UP TO 20 MEV - HOWEVER, YOU ARE FREE TO SELECT ANY	GROUPIE
ENERGY RANGE THAT YOU WISH - YOU MAY EVEN SELECT MORE THAN	GROUPIE
1 GROUP MERELY BY SPECIFYING MORE THAN 1 GROUP AS INPUT - THIS CAN BE USED TO DEFINE THE CONTRIBUTIONS TO THE RESONANCE	GROUPIE 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
WHEN THIS OPTION IS USED THE PROGRAM WILL CALCULATE GROUP AVERAGE	GROUPIE
CROSS SECTIONS - AS DEFINED ABOVE - PRIOR TO OUTPUT THE RESULTS	GROUPIE
WILL MERELY BE MULTIPLIED BY THE WIDTH OF THE GROUP ASSUMING YOU	GROUPIE
HAVE SELECTED A 1/E SPECTRUM - THERE IS NO CHECK ON THIS - THE	GROUPIE
PROGRAM MERELY MULTIPLIES THE GROUP AVERAGED CROSS SECTIONS BY,	GROUPIE
LOG(E2/E1) - WHERE E2 AND E1 ARE THE GROUP ENERGY BOUNDARIES.	GROUPIE GROUPIE
200(22, 22, MILICE DE FRE DI FRE THE GROOF ENERGI DOUNDARIES.	5100715

		GROUPIE
WARNING	STEPS OUTLINED ABOVE IF YOU WISH TO OBTAIN MEANINGFUL	GROUPIE GROUPIE GROUPIE
		GROUPIE
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	
		GROUPIE
	SECTIONS.	GROUPIE
		GROUPIE
	RESONANCE INTEGRAL OUTPUT CAN ONLY BE OBTAINED IN THE	GROUPIE
	LISTING FORMATS.	GROUPIE
		GROUPIE
	,	GROUPIE
	· · · · · · · · · · · · · · · · · · ·	GROUPIE
	· · · · · · · · · · · · · · · · · · ·	
	E PRINTED AND FOR THE SELF-SHIELDING CALCULATION ALL ENERGY	
	ALS IN WHICH THE TOTAL IS LESS THAN THE MINIMUM WILL BE	GROUPIE
IGNOREI		GROUPIE
		GROUPIE
NOTE, F	FOR THE UNSHIELDED CALCULATIONS ALL CROSS SECTIONS WILL BE	GROUPIE
CONSIDE	ERED WHETHER THEY ARE POSITIVE OR NEGATIVE. THEREFORE IF	GROUPIE
		GROUPIE
	ED RESULTS ARE UNRELIABLE AND SHOULD NOT BE USED. THEREFORE	
	S CASE THE PROGRAM WILL NOT ATTEMPT TO MODIFY THE UNSHIELDED S TO ELIMINATE THE EFFECT OF NEGATIVE CROSS SECTIONS, SINCE	
ACTUAL		GROUPIE
		GROUPIE
RESOLVE	ED RESONANCE REGION	GROUPIE
		GROUPIE
IN THE	RESOLVED RESONANCE REGION (ACTUALLY EVERYWHERE BUT IN THE	GROUPIE
	LVED RESONANCE REGION) THE CROSS SECTIONS OUTPUT BY LINEAR-	
	LCULATIONS BY THIS PROGRAM WILL YIELD ACTUAL SHIELDED AND	GROUPIE
UNSHIEL	LDED CROSS SECTIONS.	GROUPIE
INRESOI	LVED RESONANCE REGION	GROUPIE
	LVED RESONANCE REGION	GROUPIE
IN THE	UNRESOLVED RESONANCE REGION PROGRAM RECENT USES THE	GROUPIE
		GROUPIE
		GROUPIE
INFINIS	FELY DILUTE DATA AS IF IT WERE ENERGY DEPENDENT DATA AND	GROUPIE
		GROUPIE GROUPIE
GROUP A UNSHIEI	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT	GROUPIE
GROUP A UNSHIEI	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT	GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS.	GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME.	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME. AVERAGE	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME AVERAGE AND ENE	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME AVERAGE AND ENE THE ENE	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME. AVERAGE AND ENE THE ENE NUMERAT	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME. AVERAGE AND ENE THE ENE NUMERAT INSENSI	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS FEGRALS ARE PERFORMED ANALYTICALLY. THEREFORE NO ERROR IS JCED DUE TO THE USE OF TRAPAZOIDAL OR OTHER INTEGRATION . THE TOTAL ERROR THAT CAN BE ASSIGNED TO THE RESULTING ES IS JUST THAT DUE TO THE ERROR IN THE CROSS SECTIONS ERGY DEPENDENT WEIGHTING SPECTRUM. GENERALLY SINCE THE ERGY DEPENDENT WEIGHTING SPECTRUM APPEARS IN BOTH THE FOR AND THE DENOMINATOR THE AVERAGES RAPIDLY BECOME ITIVE TO THE WEIGHTING SPECTRUM AS MORE GROUPS ARE USED.	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME A AND ENE THE ENE NUMERAT INSENSI SINCE T	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME A AND ENE THE ENE NUMERAT INSENSI SINCE T USER CA	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
GROUP A UNSHIEI IT WILI ACCURAC ALL INT INTRODU SCHEME A AVERAGE THE ENE NUMERAT INSENSI SINCE T USER CA	AVERAGE IT. AS SUCH THIS PROGRAM WILL PRODUCE THE CORRECT LDED CROSS SECTION IN THE UNRESOLVED RESONANCE REGION, BUT L NOT PRODUCE THE CORRECT SELF-SHIELDING EFFECTS. CY OF RESULTS 	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE

	GROUPIE
MULTI-BAND PARAMETERS ARE CALCULATED FOR THE TOTAL, ELASTIC,	GROUPIE
CAPTURE AND FISSION REACTIONS. WITH THE NUMBER OF GROUPS THAT	GROUPIE
ARE NORMALLY USED (SEE BUILT IN GROUP STRUCTURES) ALL OTHER	GROUPIE
REACTIONS RESULT IN A NEGLIGABLE AMOUNT OF SELF-SHIELDING. AS SUCH THEIR EQUIVALENT BAND CROSS SECTION WILL MERELY BE THEIR	GROUPIE
UNSHIELDED VALUE WITHIN EACH BAND.	GROUPIE
UNSHIELDED VALUE WITHIN EACH DAND.	GROUPIE
FOR ANY GIVEN EVALUATION, WITHIN ANY GIVEN GROUP THIS PROGRAM	GROUPIE
WILL GENERATE THE MINIMUM NUMBER OF BANDS REQUIRED WITHIN THAT	GROUPIE
GROUP. AS OUTPUT TO THE COMPUTER READABLE DISK FILE THE BAND	GROUPIE
PARAMETERS FOR EACH EVALUATION WILL BE FORMATTED TO HAVE THE	GROUPIE
SAME NUMBER OF BANDS IN ALL GROUPS (WITH ZERO WEIGHT FOR SOME	GROUPIE
BANDS WITHIN ANY GROUP). THE USER MAY DECIDE TO HAVE OUTPUT	GROUPIE
EITHER WITH THE MINIMUM NUMBER OF BANDS REQUIRED FOR EACH	GROUPIE
EVALUATION (E.G. 2 BANDS FOR HYDROGEN AND 4 BANDS FOR U-233) OR	GROUPIE
THE SAME NUMBER OF BANDS FOR ALL EVALUATIONS (E.G. 4 BANDS FOR	GROUPIE
BOTH HYDROGEN AND U-233).	GROUPIE
	GROUPIE
FOR 2 OR FEWER BANDS THE PROGRAM USES AN ANALYTIC EXPRESSION	GROUPIE
TO DEFINE ALL MULTI-BAND PARAMETERS. FOR MORE THAN 2 BANDS THE	GROUPIE
PROGRAM PERFORMS A NON-LINEAR FIT TO SELECT THE MULTI-BAND PARAMETERS THAT MINIMIZE THE MAXIMUM FRACTIONAL ERROR AT ANY	GROUPIE
POINT ALONG THE ENTIRE SELF-SHIELDING CURVE. THE NUMBER OF BANDS	GROUPIE
REQUIRED WITHIN ANY GIVEN GROUP IS DEFINED BY INSURING THAT THE	GROUPIE
MULTI-BAND PARAMETERS CAN BE USED TO ACCURATELY DEFINE SELF-	GROUPIE
SHIELDED CROSS SECTIONS ALONG THE ENTIRE SELF-SHIELDING CURVE	GROUPIE
FROM SIGMA0 = 0 TO INFINITY. THE USER MAY DEFINE THE ACCURACY	GROUPIE
REQUIRED.	GROUPIE
	GROUPIE
ENDF/B FORMATTED UNSHIELDED AVERAGES	GROUPIE
	GROUPIE
UNSHIELDED MULTI-GROUP AVERAGED CROSS SECTIONS FOR ALL REACTIONS	GROUPIE
MAY BE OBTAINED IN THE ENDF/B FORTRAN IN EITHER HISTOGRAM	GROUPIE
(INTERPOLATION LAW 1) OR LINEARLY INTERPOLABLE (INTERPOLATION	GROUPIE
LAW 2) FORM. SEE INPUT BELOW FOR DETAILS.	GROUPIE
	GROUPIE
MIXTURES OF MATERIALS AND RESONANCE OVERLAP	GROUPIE
	GROUPIE
THE SELF-SHIELDED CROSS SECTIONS FOR THE INDIVIDUAL CONSTITUENTS 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	
FOR EACH CONSTITUENT AND TO RUN THIS PROGRAM. THIS CAN BE DONE BY	
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
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	GROUPIE
CAN IN TURN BE USED TO CALCULATE FULLY CORRELATED MULT-BAND	GROUPIE
MULTI-BAND PARAMETER OUTPUT FORMAT	GROUPIE 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
THE BINARY FORMAT USED IN EARLIER VERSIONS OF THIS CODE IS NO	GROUPIE
THE BINARY FORMAT USED IN EARLIER VERSIONS OF THIS CODE IS NO LONGER USED.	GROUPIE GROUPIE
	GROUPIE GROUPIE GROUPIE
	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
LONGER USED.	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
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	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
LONGER USED. CONTACT THE AUTHOR IF YOU WOULD LIKE TO RECEIVE A SIMPLE PROGRAM TO READ THE CHARACTER FORMATTED MULTI-BAND PARAMETER FILE AND	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
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	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE

RECORI	D COLUMNS FORMAT DESCH	IPTION	GROUPIE
1		RY DESCRIPTION (AS READ)	GROUPIE
2	1-11 I11 MATER		GROUPIE
		R GROUPS	GROUPIE
	23-33 I11 NUMBE	R OF BANDS	GROUPIE
	34-44 E11.4 TEMPE	RATURE (KELVIN)	GROUPIE
		RITH DESCRIPTION OF ZA	GROUPIE
3		Y (EV) - GROUP BOUNDARY.	GROUPIE
	12-22 E11.4 TOTAI		GROUPIE
	23-33 E11.4 ELAST 34-44 E11.4 CAPTU		GROUPIE GROUPIE
	35-55 E11.4 FISS		GROUPIE
4	1-11 BLANK		GROUPIE
	12-22 E11.4 TOTAI		GROUPIE
	23-33 E11.4 ELASI	'IC	GROUPIE
	34-44 E11.4 CAPTU	RE	GROUPIE
	35-55 E11.4 FISSI	ON	GROUPIE
			GROUPIE
		CH GROUP. THE LAST LINE FOR EACH	
MATER	IAL (ZA) IS,		GROUPIE
N	1-11 E11.4 ENERG	Y (EV) - UPPER ENERGY LIMIT OF	GROUPIE
М	I II EII.4 ENERG	LAST GROUP.	GROUPIE
			GROUPIE
FOR EX	XAMPLE, A 175 GROUP, 2 BAND	FILE, FOR EACH MATERIAL WILL	GROUPIE
CONTAI	IN 352 LINES = 1 HEADER LINE	, 175 * 2 LINES OF PARAMETERS,	GROUPIE
	AND 1 FINAL I	INE WITH THE UPPER ENERGY LIMIT	GROUPIE
	OF THE LAST O	ROUP.	GROUPIE
			GROUPIE
	FILES		GROUPIE
	DESCRIPTION		GROUPIE GROUPIE
			GROUPIE
2	INPUT DATA (BCD - 80 CHARAC	TERS/RECORD)	GROUPIE
	ORIGINAL ENDF/B DATA (BCD -		GROUPIE
			GROUPIE
OUTPUT	T FILES		GROUPIE
			GROUPIE
	DESCRIPTION		GROUPIE
			GROUPIE
3 11	OUTPUT REPORT (BCD - 80 CHA MULTI-GROUP ENDF/B DATA - 0		GROUPIE
	(BCD - 80 CHARACTERS/RECORD		GROUPIE
16	PLOTTAB FORMATTED SELF-SHIP	•	GROUPIE
	(BCD - 80 CHARACTERS/RECORD))	GROUPIE
31	MULTI-BAND PARAMETERS CHARA	CTER FILE - OPTIONAL	GROUPIE
	(BCD - 80 CHARACTERS/RECORD		GROUPIE
32		SEUDO ENDF FORMAT - OPTIONAL	GROUPIE
22	(BCD - 120 CHARACTERS/RECOR		GROUPIE
33	SELF-SHIELDED CROSS SECTION (BCD - 120 CHARACTERS/RECOR		GROUPIE GROUPIE
34	MULTI-BAND PARAMETER LISTIN	-	GROUPIE
51	(BCD - 120 CHARACTERS/RECOF		GROUPIE
35	UNSHIELDED CROSS SECTION LI	-	GROUPIE
	(BCD - 120 CHARACTERS/RECOR	D)	GROUPIE
			GROUPIE
	CH FILES		GROUPIE
			GROUPIE
UNIT	FILENAME DESCRIPTION		GROUPIE GROUPIE
	ENERGY DEPENDENT WEIGHTING	SPECTRUM	GROUPIE
0	(BINARY - 40080 WORDS/BLOCH		GROUPIE
9	TOTAL CROSS SECTION		GROUPIE
	(BINARY - 40080 WORDS/BLOCK	.)	GROUPIE
12		Y FOR SELF-SHIELDING CALCULATION	GROUPIE
	(BINARY - 40080 WORDS/BLOCH		GROUPIE
13		Y FOR SELF-SHIELDING CALCULATION	
14	(BINARY - 40080 WORDS/BLOCH	.) Y FOR SELF-SHIELDING CALCULATION	GROUPIE
14	(BINARY - 40080 WORDS/BLOCH		GROUPIE
		-,	

GROUPIE OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINES FILIO1 AND FILIO2) GROUPIE _____ GROUPIE UNIT FILE NAME GROUPTE ____ _____ GROUPIE 2 GROUPIE.INP GROUPIE GROUPTE 3 GROUPIE.LST GROUPIE _____ _____ GROUPIE 8 (SCRATCH) 9 (SCRATCH) GROUPIE ENDFB.IN 10 GROUPIE 11 ENDFB.OUT GROUPIE 12 (SCRATCH) GROUPTE 13 (SCRATCH) GROUPIE (SCRATCH) 14 GROUPIE -----2019/6/23 - New Filenames (added ZAzzzaaa at Beginning)------GROUPIE PLOTTAB.CUR ZAzzzaaa.PLOT.CUR GROUPIE 16 31 MULTBAND. TAB ZAzzzaaa.MULTBAND.TAB GROUPIE 32 ZAzzzaaa.URR.ENDF GROUPIE 33 SHIELD.LST GROUPIE ZAzzzaaa.SHIELD.LST MULTBAND.LST GROUPIE 34 ZAzzzaaa.MULTBAND.LST UNSHIELD.LST 35 ZAzzzaaa.UNSHIELD.LST GROUPTE 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 GROUPTE MULTIBAND OUTPUT IS REQUESTED. GROUPIE GROUPIE INPUT CARDS GROUPIE GROUPIE CARD COLS. FORMAT DESCRIPTION GROUPIE ____ -----_____ GROUPTE II1 SELECTION CRITERIA (0=MAT, 1=ZA)
II1 NUMBER OF GROUPS. GROUPTE 1 1-11 1 12-22 GROUPIE Unresolved Resonance Region Self-Shielding GROUPIE calculation. GROUPIE =.GT.0 - ARBITRARY GROUP BOUNDARIES ARE READ GROUPIE FROM INPUT FILE (N GROUPS REQUIRE GROUPIE N+1 GROUP BOUNDARIES). CURRENT GROUPTE PROGRAM MAXIMUM IS 20,000 GROUPS. GROUPIE BUILT-IN OPTIONS INCLUDE.... GROUPIE = 0 - TART 175 GROUPS GROUPIE = -1 - ORNL 50 GROUPS GROUPIE 126 GROUPS = -2 - ORNL GROUPIE = -3 - ORNL 171 GROUPS GROUPTE - SAND-II 620 (665) GROUPS TO 18 MEV GROUPIE = -4 = -5 - SAND-II 640 (685) GROUPS TO 20 MEV GROUPIE 69 GROUPS = -6 - WIMS GROUPIE = -7 - GAM-I 68 GROUPS GROUPIE = -8 - GAM-II 99 GROUPS GROUPIE = -9 54 GROUPS – MUFT GROUPIE 28 GROUPS =-10 - ABBN GROUPIE Current TART Standard =-11 - TART 616 GROUPS TO 20 MEV GROUPIE 700 GROUPS TO 1 GEV (-11 is required for =-12 - TART GROUPIE - SAND-II 665 GROUPS TO 18 MEV unresolved resonance =-13 GROUPTE region self-shielding=-14 - SAND-II 685 GROUPS TO 20 MEV GROUPIE calculations) =-15 - TART 666 GROUPS TO 200 MEV GROUPIE - SAND-II 725 GROUPS TO 60 MEV =-16 GROUPIE - SAND-II 755 GROUPS TO 150 MEV =-17 GROUPIE - SAND-II 765 GROUPS TO 200 MEV =-18 GROUPIE - UKAEA 1102 GROUPS TO 1 GeV =-19 GROUPIE 1 23-33 I11 MULTI-BAND SELECTOR GROUPIE ----2019/6/23 - ONLY 0 or 2 allowed = the = 1GROUPIE option has proven to give very poor results, GROUPIE and therefore is no longer allowed. GROUPIE = 0 - NO MULTI-BAND CALCULATIONS GROUPIE

No longer allowed	= 1 - 2 BAND. CONSERVE AV(TOT), AV(1/TOT)	GROUPIE
	AND AV(1/TOT**2)	GROUPIE
	= 2 - 2 BAND. CONSERVE AV(TOT), AV(1/TOT)	GROUPIE
	AND AV(1/(TOT+SIGMA0)) WHERE	GROUPIE
	SIGMA0 = AV(TOT) IN EACH GROUP	GROUPIE
No longer allowed	= 3-5- MULTI-BAND FIT. CONSERVE AV(TOT) AND	GROUPIE
	MINIMIZE FRACTIONAL ERROR FOR ENTIRE	GROUPIE
	SELF-SHIELDING CURVE (SIGMA0 = 0 TO	GROUPIE
	INFINITY)	GROUPIE
	IF THE SELECTOR IS POSITIVE (1 TO 5) THE	GROUPIE
	MINIMUM NUMBER OF BANDS WILL BE OUTPUT FOR	GROUPIE
	EACH ISOTOPE INDEPENDENTLY. IF THE SELECTOR	GROUPIE
	IS NEGATIVE (-1 TO -5) THE SAME NUMBER OF	GROUPIE
	BANDS (ABS(SELECTOR)) WILL BE OUTPUT FOR	GROUPIE
1 24 44	ALL ISOTOPES.	GROUPIE
1 34-44 I11	NUMBER OF POINTS USED TO DESCRIBE ENERGY	GROUPIE
	DEPENDENT WEIGHTING SPECTRUM S(E).	GROUPIE
	= 0 or 1 - Flat (Constant) = -1 - 1/E at ALL energies	GROUPIE
	= -1 - 1/E at ALL energies = -2 - MAXWELLIAN - UP TO 0.1 eV	GROUPIE GROUPIE
	1/E - 0.1 eV TO 67 KeV	GROUPIE
	FISSION - 67 KeV to 10 MeV	GROUPIE
	CONSTANT - Above 10 MeV	GROUPIE
	= > 1 - Read input table	GROUPIE
2005/01/20	-ADDED OPTION TO ALLOW TEMPERATURE OF THE	GROUPIE
2000, 01, 20	MAXWELLIAN TO BE CHANGED - SEE INPUT LINE 4,	
	COLUMNS 55 - 66.	GROUPIE
	= -1 - 1/E	GROUPIE
	= 0 OR 1- ENERGY INDEPENDENT (SO CALLED FLAT	
	WEIGHTING SPECTRUM) .	GROUPIE
	= .GT.1 - READ THIS MANY POINTS FROM INPUT	GROUPIE
	TO DESCRIBE WEIGHTING SPECTRUM.	GROUPIE
	NO LIMIT TO THE NUMBER OF POINTS	GROUPIE
	USED TO DESCRIBE WEIGHTING.	GROUPIE
1 45-55 E11.4	MULTI-BAND CONVERGENCE CRITERIA.	GROUPIE
	-2019/6/23 - No longer used now that code	GROUPIE
	is restricted to no more than 2 bands.	GROUPIE
	ONLY USED FOR 3 OR MORE BANDS. THE NUMBER OF	GROUPIE
	BANDS IN EACH GROUPS IS SELECTED TO INSURE	GROUPIE
	THAT THE ENTIRE SELF-SHIELDING CURVE CAN BE	GROUPIE
	REPRODUCED TO WITHIN THIS FRACTIONAL ERROR.	GROUPIE
	= .LT. 0.0001 - USE STANDARD 0.001	GROUPIE
	(0.1 PER-CENT)	GROUPIE
1 56.66 -11	= .GE. 0.0001 - USE AS CONVERGENCE CRITERIA	
1 56-66 I11		GROUPIE
	-2019/6/23 - For multi-band calculations only 0 is alllowed = mulriples of unshielded	GROUPIE
	total in each group = This is required for	GROUPIE
	total in each group = This is required for the BEST self-shielding results.	GROUPIE
	< 0 - 21 VALUES OF SIGMAO ARE READ INPUT AND	
	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	
	EACH DECADE VALUES OF 10, 7, 4, 2, 1	GROUPIE
	BARNS WILL BE USED).	GROUPIE
1 67-70 I 4	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. $(abardard EWDE(R))$	GROUPIE
	= 0 = cross section = 0 (standard ENDF/B)	GROUPIE
	= 1 = cross section = constant (equal to	GROUPIE

			value at highest tabulated energy).	GROUPIE
2-4	1-66 6	SE11.4	SIGMA-0 Definition	GROUPIE
				GROUPIE
			in allowed for Unresolved Resonance Region	GROUPIE
			Self-Shielding calculation	GROUPIE
			IF SIGMA-0 DEFINITION SELECTOR < 0, THE NEXT	GROUPIE
			4 LINES OF INPUT ARE THE 22 VALUES OF SIGMA0,	
			6 PER LINE.	GROUPIE
2	1-72	A72	ENDF/B INPUT DATA FILENAME	GROUPIE
3	1 70	770	(STANDARD OPTION = ENDFB.IN) ENDF/B OUTPUT DATA FILENAME	GROUPIE GROUPIE
3	1-72	A/Z	(STANDARD OPTION = ENDFB.OUT)	GROUPIE
			(STANDARD OFIION - ENDID.001)	GROUPIE
THE FO	OURTH IN	IPUT CARL) IS USED TO SELECT ALL DESIRED OUTPUT MODES.	
EACH C	OUTPUT I	EVICE MA	AY BE TURNED OFF (0) OR ON (1). THEREFORE	GROUPIE
THEREP	FORE EAC	CH OF THE	E FOLLOWING INPUT PARAMETERS MAY BE EITHER	GROUPIE
ZERO 1	TO INDIC	CATE NO C	OUTPUT OR NON-ZERO TO INDICATE OUTPUT.	GROUPIE
				GROUPIE
4	1-11	I11		GROUPIE
			= 1 - CROSS SECTIONS	GROUPIE
	10.00	T11	= 2 - RESONANCE INTEGRALS	GROUPIE
4	12-22		MULTI-BAND PARAMETER LISTING MULTI-BAND PARAMETERS COMPUTER READABLE	GROUPIE GROUPIE
4			UNSHIELDED CROSS SECTIONS IN ENDF/B FORMAT	GROUPIE
-	51 11		= 1 - HISTOGRAM FORMAT (INTERPOLATION LAW 1)	
			= 2 - LINEAR-LINEAR (INTERPOLATION LAW 2)	GROUPIE
4	45-55	I11	UNSHIELDED CROSS SECTIONS LISTING	GROUPIE
			= 1 - CROSS SECTIONS	GROUPIE
			= 2 - RESONANCE INTEGRALS	GROUPIE
05/01/20 -	- ADDED			GROUPIE
4	56-66	E11.4		GROUPIE
			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)	GROUPIE
			= 0 - USE DEFAULT 0.0253 EV = ROOM TEMPERATURE	
			> 0 - USE THIS AS THE TEMPERATURE	GROUPIE
			RESTRICTION - TEMPERATURE CANNOT EXCEED	GROUPIE
			1000 EV.	GROUPIE
				GROUPIE
5	1-80	18A4	LIBRARY IDENTIFICATION. ANY TEXT THAT THE	GROUPIE
			USER WISHES TO IDENTIFY THE MULTI-BAND	GROUPIE
			PARAMETERS. THIS LIBRARY IDENTIFICATION IS	GROUPIE
			WRITTEN INTO THE COMPUTER READABLE MULTI-BAND DATA FILE.	GROUPIE
			DATA FILE.	GROUPIE
6-N	1- 6	I6	LOWER MAT OR ZA LIMIT	GROUPIE
•	7-8		LOWER MF LIMIT	GROUPIE
	9-11	13	LOWER MT LIMIT	GROUPIE
	12-17	I11	UPPER MAT OR ZA LIMIT	GROUPIE
	18-19	12	UPPER MF LIMIT	GROUPIE
	20-22		UPPER MT LIMIT	GROUPIE
			UP TO 100 RANGES MAY BE SPECIFIED, ONE RANGE	
				GROUPIE
			BY A BLANK CARD. IF THE UPPER MAT OR ZA	GROUPIE
			LIMIT IS LESS THAN THE LOWER LIMIT THE UPPER IS SET EQUAL TO THE LOWER LIMIT. IF THE UPPER	
			MF OR MT LIMIT IS ZERO IT WILL BE SET EQUAL	
			TO ITS MAXIMUM VALUE, 99 OR 999, RESPECTIVELY	
			IF THE FIRST REQUEST LINE IS BLANK IT WILL	GROUPIE
			TERMINATE THE LIST OF REQUESTS AND CAUSE ALL	
			_	GROUPIE
				GROUPIE
VARY	1-66	6E11.4	-	GROUPIE
			THE NUMBER OF GROUPS INDICATED ON THE FIRST	GROUPIE
			INPUT CARD IS POSITIVE. ALL ENERGIES MUST	GROUPIE
				GROUPIE
			LIMITS ARE 1 TO 20,000 GROUPS. FOR N GROUPS 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
	GROUPIE
VARY 1-66 6E11.4 ENERGY DEPENDENT WEIGHTING SPECTRUM. ONLY	GROUPIE
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 ORDER IN EV. THE SPECTRUM VALUES MUST BE	GROUPIE GROUPIE
NON-NEGATIVE. THE ENERGY RANGE OF SPECTRUM	GROUPIE
MUST AT LEAST SPAN THE ENERGY RANGE OF THE	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
WEIGHTING SPECTRUM.	GROUPIE
	GROUPIE
EXAMPLE INPUT NO. 1	GROUPIE
	GROUPIE
REQUEST DATA BY MAT AND PROCESS ALL DATA (ALL MAT BETWEEN 1 AND	GROUPIE
9999). USE THE TART 175 GROUP STRUCTURE, GENERATE 2 BAND	GROUPIE
PARAMETERS (THE FOR ALL ISOTOPES) TO 0.1 PER-CENT ACCURACY	GROUPIE
IN THE SELF-SHIELDING CURVE. OUTPUT ALL LISTING, COMPUTER	GROUPIE
READABLE AND ENDF/B FORMAT GROUP AVERAGES.	GROUPIE
	GROUPIE
EXPLICITLY SPECIFY THE STANDARD FILENAMES.	GROUPIE
THE FOLLOWING 7 INDUM LINES ARE DECULDED	GROUPIE GROUPIE
THE FOLLOWING 7 INPUT LINES ARE REQUIRED.	GROUPIE
0 0 -2 0 1.00000-03 0	GROUPIE
ENDFB.IN	GROUPIE
ENDFB.OUT	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. 2	GROUPIE
	GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ	GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND	GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ	GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238	GROUPIE GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND	GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238 THE FOLLOWING 7 INPUT LINES ARE REQUIRED.	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238THE FOLLOWING 7 INPUT LINES ARE REQUIRED.00-20 1.00000-030	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238 THE FOLLOWING 7 INPUT LINES ARE REQUIRED. 0 0 -2 0 1.00000-03 0 \ENDFB6\SIGMA1\K300\ZA092238 \ENDFB6\GROUPIE\K300\ZA092238 1 1 1	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238 THE FOLLOWING 7 INPUT LINES ARE REQUIRED. 0 0 -2 0 1.00000-03 0 \ENDFB6\SIGMA1\K300\ZA092238 \ENDFB6\GROUPIE\K300\ZA092238 1 1 1 1 1 1 TART 175 GROUP, 2 BAND LIBRARY TO 0.1 PER-CENT ACCURACY	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
THE SAME EXAMPLE 1, AS ABOVE, ONLY THE ENDF/B DATA WILL BE READ FROM \ENDFB6\SIGMA1\K300\ZA092238 (U-238 AT 300 KELVIN) AND WRITTEN TO \ENDFB6\GROUPIE\K300\ZA092238 THE FOLLOWING 7 INPUT LINES ARE REQUIRED. 0 0 -2 0 1.00000-03 0 \ENDFB6\SIGMA1\K300\ZA092238 \ENDFB6\GROUPIE\K300\ZA092238 1 1 1 1 1 1 TART 175 GROUP, 2 BAND LIBRARY TO 0.1 PER-CENT ACCURACY 1 1 9999 0 0	GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE GROUPIE
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					GROUPIE
EXAMPLE INPUT NO.	4				GROUPIE
					GROUPIE
THIS EXAMPLE USES	A USER DEFI	INED GROUP S	STRUCTURE AN	ND WEIGHTING	GROUPIE
FUNCTION - THESE A	ARE NOT REAL	LISTIC IN TH	ERMS OF ACTU	JAL ENERGIES	GROUPIE
AND WEIGHTS - THE	ARE ONLY I	INTENDED TO	ILLUSTRATE	THE ORDER O	OF GROUPIE
THE INPUT PARAMETH	ERS.				GROUPIE
					GROUPIE
0 11	0	6		0	GROUPIE
RECENT . OUT					GROUPIE
GROUPIE.OUT					GROUPIE
1 1	1	1	1		GROUPIE
Example with users det	fined groupu	is and spect	trum weight:	ing	GROUPIE
1 1 1 999999999					GROUPIE
	(blabk line	e terminates	s request li	ist)	GROUPIE
1.00000-05 1.00000-04	1.00000-03	1.00000-02	1.00000-01	1.00000+00	grouGROUPIE
1.00000+01 1.00000+02	1.00000+03	1.00000+04	1.00000+05	1.00000+06	grouGROUPIE
1.00000-05 1.0	1.00000-02	0.1	1.00000+00	0.01	weigGROUPIE
1.00000+02 0.001	1.00000+04	0.0001	1.00000+06	0.00001	weigGROUPIE
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