

=====	Sixpak
PROGRAM SIXPAK	Sixpak
=====	Sixpak
VERSION 92-1 (JANUARY 1992)	Sixpak
VERSION 92-2 (FEBRUARY 1992)*INCREASED CORE ALLOCATION TO	Sixpak
ACCOMMODATE JEF AND EFF EVALUATIONS.	Sixpak
VERSION 92-3 (APRIL 1992) *ADDED ADDITIONAL DATA TESTS.	Sixpak
VERSION 92-4 (SEPT. 1992) *CORRECTED KALBACH-MANN CALCULATIONS.	Sixpak
*FOR PHOTON PRODUCTION OUTPUT MF=12	Sixpak
(MULTIPLICITY), MF=14 (ISOTROPIC	Sixpak
ANGULAR DISTRIBUTIONS) AND MF=15	Sixpak
(SPECTRA) - PREVIOUSLY ONLY MF=15.	Sixpak
*FIRST ORDER CORRECTIONS TRANSFORMING	Sixpak
CENTER-OF-MASS SPECTRA TO LAB SYSTEM	Sixpak
FOR OUTPUT IN MF=5	Sixpak
*CORRECTED ISOTROPIC ANGULAR	Sixpak
DISTRIBUTION FLAG (LI)	Sixpak
VERSION 94-1 (JANUARY 1994) *VARIABLE ENDF/B INPUT DATA FILENAME	Sixpak
TO ALLOW ACCESS TO FILE STRUCTURES	Sixpak
(WARNING - INPUT PARAMETER FORMAT	Sixpak
HAS BEEN CHANGED)	Sixpak
*CLOSE ALL FILES BEFORE TERMINATING	Sixpak
(SEE, SUBROUTINE ENDIT)	Sixpak
*INCREASED MAXIMUM TABLE SIZE FROM	Sixpak
2000 TO 6000.	Sixpak
VERSION 96-1 (JANUARY 1996) *COMPLETE RE-WRITE	Sixpak
*IMPROVED COMPUTER INDEPENDENCE	Sixpak
*ALL DOUBLE PRECISION	Sixpak
*ON SCREEN OUTPUT	Sixpak
*UNIFORM TREATMENT OF ENDF/B I/O	Sixpak
*IMPROVED OUTPUT PRECISION	Sixpak
VERSION 99-1 (MARCH 1999) *CORRECTED CHARACTER TO FLOATING	Sixpak
POINT READ FOR MORE DIGITS	Sixpak
*UPDATED TEST FOR ENDF/B FORMAT	Sixpak
VERSION BASED ON RECENT FORMAT CHANGE	Sixpak
*GENERAL IMPROVEMENTS BASED ON	Sixpak
USER FEEDBACK	Sixpak
VERSION 99-2 (JUNE 1999) *ASSUME ENDF/B-VI, NOT V, IF MISSING	Sixpak
MF=1, MT=451.	Sixpak
VERS. 2000-1 (FEBRUARY 2000) *GENERAL IMPROVEMENTS BASED ON	Sixpak
USER FEEDBACK	Sixpak
VERS. 2002-1 (JANUARY 2002) *CORRECTED ANGULAR DISTRIBUTION (MF=4)	Sixpak
OUTPUT TO INSURE USED FIELDS ARE 0	Sixpak
(MAY 2002) *OPTIONAL INPUT PARAMETERS	Sixpak
(NOV. 2002) *EXTENDED TO ALLOW CHARGED PARTICLE	Sixpak
ANGULAR DISTRIBUTION IN MF=4 -	Sixpak
WARNING - STRICTLY SPEAKING THIS IS	Sixpak
NOT LEGAL, SINCE MF=4 IS SUPPOSED TO	Sixpak
BE USED ONLY FOR NEUTRON ANGULAR	Sixpak
DISTRIBUTIONS - BUT WHERE MT MAKES	Sixpak
IT OBVIOUS THAT THE OUTGOING PARTICLE	Sixpak
IS NOT A NEUTRON HOPEFULLY IT WILL	Sixpak
NOT CAUSE A PROBLEM IF MF=4 IS USED	Sixpak
FOR CHARGED PARTICLES.	Sixpak
VERS. 2004-1 (MARCH 2004) *ADDED INCLUDE FOR COMMON	Sixpak
*INCREASED MAXIMUM TABLE SIZE FROM	Sixpak
6,000 TO 12,000.	Sixpak
*ADDED DUMMY A FOR ELEMENTS	Sixpak
*CORRECTED OUTPUT INTERPOLATION LAWS	Sixpak
VERS. 2007-1 (JAN. 2007) *CHECKED AGAINST ALL ENDF/B-VII.	Sixpak
*INCREASED MAXIMUM TABLE SIZE FROM	Sixpak
12,000 TO 120,000.	Sixpak

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AS A PART OF AN INTERNATIONAL PROJECT ON THE EXCHANGE OF  
NUCLEAR DATA

## ACKNOWLEDGEMENT (VERSION 92-1)

THE AUTHOR THANKS SOL PEARLSTEIN (BROOKHAVEN NATIONAL LAB) FOR  
SIGNIFICANTLY CONTRIBUTING TOWARD IMPROVING THE ACCURACY AND  
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## ACKNOWLEDGEMENT (VERSION 92-4)

THE AUTHOR THANKS BOB MACFARLANE (LOS ALAMOS) FOR SUGGESTING HOW TO PROPERLY OUTPUT THE PHOTON PRODUCTION DATA TO PUT IT INTO EXACTLY THE FORM NEEDED FOR USE IN PROCESSING CODES.

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AUTHORS MESSAGE

THE COMMENTS BELOW SHOULD BE CONSIDERED THE LATEST DOCUMENTATION INCLUDING ALL RECENT IMPROVEMENTS. PLEASE READ ALL OF THESE COMMENTS BEFORE IMPLEMENTING AND USING THESE CODES.

AT THE PRESENT TIME WE ARE ATTEMPTING TO DEVELOP A SET OF COMPUTER INDEPENDENT PROGRAMS THAT CAN EASILY BE IMPLEMENTED ON ANY ONE OF A WIDE VARIETY OF COMPUTERS. IN ORDER TO ASSIST IN THIS PROJECT IT WOULD BE APPRECIATED IF YOU WOULD NOTIFY THE AUTHOR OF ANY COMPILER DIAGNOSTICS, OPERATING PROBLEMS OR SUGGESTIONS ON HOW TO IMPROVE THIS PROGRAM. HOPEFULLY. IN THIS WAY FUTURE VERSIONS OF

[illegible]

THIS PROGRAM WILL BE COMPLETELY COMPATIBLE FOR USE ON YOUR COMPUTER.	Sixpak
	Sixpak
PURPOSE	Sixpak
=====	Sixpak
1) CHECK ALL DOUBLE-DIFFERENTIAL DATA (MF=6)	Sixpak
	Sixpak
2) OUTPUT EQUIVALENT MF = 4, 5, 12, 14 AND 15 DATA.	Sixpak
	Sixpak
DATA CHECKING	Sixpak
=====	Sixpak
ALL OF THE ENDF/B-VI MF=6 DATA IS CHECKED - FOR DETAILS SEE BELOW.	Sixpak
	Sixpak
THE MF=6 DATA IS NOT CORRECTED AND OUTPUT IN THE ENDF/B FORMAT. IT IS MERELY CHECKED. IF ERRORS ARE FOUND IT IS UP TO THE USER TO TAKE CORRECTIVE ACTION ON THE MF=6 DATA.	Sixpak
	Sixpak
IN CONTRAST WHEN PROBLEMS ARE FOUND IN DATA WHICH WILL BE OUTPUT IN THE ENDF/B FORMAT (MF=4, 5, 12, 14 AND 15), WHENEVER POSSIBLE CORRECTIVE ACTION WILL BE TAKEN.	Sixpak
	Sixpak
FURTHER CHECKS AND CORRECTIONS	Sixpak
=====	Sixpak
ONCE THE DATA HAS BEEN OUTPUT IN MF = 4, 5, 12, 14 AND 15 FORMATS FURTHER CORRECTIVE ACTION CAN BE TAKEN AS FOLLOWS,	Sixpak
	Sixpak
PROGRAM LEGEND	Sixpak
=====	Sixpak
CAN BE USED TO CORRECT ANGULAR DISTRIBUTIONS WHICH ARE NEGATIVE, TO CONVERT FROM LEGENDRE COEFFICIENTS TO TABULATED ANGULAR DISTRIBUTIONS AND GENERALLY PERFORM MORE EXTENSIVE TESTS OF ALL MF=4 DATA.	Sixpak
	Sixpak
PROGRAM EVALPLOT	Sixpak
=====	Sixpak
VERSION 92-1 AND LATER VERSIONS CAN PLOT ALL OF THE MF=4, 5 AND 15 DATA OUTPUT BY THIS CODE. EARLIER VERSIONS CAN PLOT MF=4 AND 5. GRAPHICS IS AN EXCELLENT WAY TO CHECK THIS DATA.	Sixpak
	Sixpak
PROGRAM PLOTTAB	Sixpak
=====	Sixpak
THIS IS A GENERAL PLOTTING PROGRAM AND THERE IS AN INTERFACE IN THIS CODE TO PRODUCE OUTPUT FOR ANY MF=6 DATA IN THE PLOTTAB INPUT FORMAT. THIS PROGRAM CAN BE USED TO CHECK ALL OF THE MF=6 DATA AS WELL AS THE EQUIVALENT MF=4, 5, 12, 14 AND 15 DATA - AS WELL AS COMPARING THE ORIGINAL MF=6 AND EQUIVALENT DATA.	Sixpak
	Sixpak
DATA OUTPUT	Sixpak
=====	Sixpak
THE ENDF/B MF=4, 5, 12, 14 AND 15 FORMATS ONLY ALLOW FOR NEUTRONS INCIDENTS	Sixpak
	Sixpak
THE ENDF/B MF=4 AND 5 FORMATS ONLY ALLOW FOR NEUTRONS OUTGOING.	Sixpak
	Sixpak
THE ENDF/B MF=12, 14 AND 15 ONLY ALLOWS FOR PHOTONS OUTGOING.	Sixpak
	Sixpak
THESE ARE THE ONLY COMBINATIONS OF DATA OUTPUT BY THIS CODE.	Sixpak
	Sixpak
ALL OTHER COMBINATIONS OF INCIDENT AND OUTGOING PARTICLES ARE CHECKED, BUT THE RESULTS CANNOT BE OUTPUT IN THE ENDF/B FORMAT. HOWEVER, USING THE PLOTTAB INTERFACE BUILT INTO THIS CODE THIS DATA CAN, AND HAS BEEN, OUTPUT AND CHECKED.	Sixpak
	Sixpak

THE NEUTRON DATA IN MF=4 CAN BE IN THE FORM OF EITHER TABULATED ANGULAR DISTRIBUTIONS OR LEGENDRE COEFFICIENTS.	Sixpak
	Sixpak
	Sixpak
THE NEUTRON (MF=5) OR PHOTON (MF=15) SPECTRA ARE BOTH IN EXACTLY THE SAME FORMAT = ARBITRARY TABULATED FUNCTIONS - ENDF/B OPTION LF=1.	Sixpak
	Sixpak
ENDF/B DATA OUTPUT ORDER	Sixpak
=====	Sixpak
ENDF/B DATA IS OUTPUT IN ASCENDING MAT, MF, MT ORDER. IN ORDER TO ALLOW THIS PROGRAM TO PRODUCE ALL OUTPUT IN A SINGLE PASS THROUGH THE MF=6 DATA, OUTPUT FOR EACH (MAT, MT) IS OUTPUT TO SEPERATE FILES FOR MF=4, 5, 12, 14 AND 15.	Sixpak
	Sixpak
FOR SUBSEQUENT USE THE ENDF/B FORMATTED DATA OUTPUT BY THIS CODE CAN BE MERGED TOGETHER USING PROGRAM MERGER (CONTAIN THE AUTHOR OF THIS CODE FOR A COPY OF MERGER), E.G., MERGE MF=12, 14 AND 15 DATA IN ORDER TO THEN CALCULATE PHOTON PRODUCTION DATA OR MF=4 AND 5 CAN BE MERGED TOGETHER TO CALCULATE NEUTRON TRANSFER - OR ALL OF THEM CAN BE MERGED TOGETHER TO PERFORM NEUTRON AND PHOTON CALCULATIONS.	Sixpak
	Sixpak
CORRELATED (MF=6) VS. UNCORRELATED (MF=4 AND 5) DATA	Sixpak
=====	Sixpak
THE ENDF/B DOUBLE DIFFERENTIAL = CORRELATED - DATA IN MF=6 REPRESENTS DATA IN THE FORM,	Sixpak
	Sixpak
$F(E,EP,COS) = SIG(E)*Y(E)*G0(E,EP)*F(E,EP,COS)$	Sixpak
	Sixpak
$SIG(E) = MF=3 \text{ CROSS SECTIONS}$	Sixpak
$Y(E) = \text{YIELD (MULTIPLICITY)}$	Sixpak
$G0(E,EP) = \text{ENERGY SPECTRUM}$	Sixpak
$F(E,EP,COS) = \text{ANGULAR DISTRIBUTION}$	Sixpak
	Sixpak
IN A SITUATION WHERE YOU HAVE MONOENERGETIC AND MONODIRECTIONAL NEUTRONS INCIDENT YOU WILL BE ABLE TO OBSERVE CORRELATION EFFECTS IN THE NEUTRON SPECTRUM AND ANGULAR DISTRIBUTION.	Sixpak
	Sixpak
EVEN IN SITUATIONS WHERE YOU HAVE A NARROW SPECTRUM OF NEUTRONS THAT ARE HIGHLY DIRECTIONALLY ORIENTED YOU MAY BE ABLE TO OBSERVE THESE CORRELATION EFFECTS, E.G., A NARROW 14 MEV FUSION SOURCE INCIDENT ON THE FIRST WALL OF A CTR DEVICE.	Sixpak
	Sixpak
FOR SUCH SITUATIONS USE OF THE CORRELATED (MF=6) DATA IS REQUIRED IN CALCULATIONS.	Sixpak
	Sixpak
HOWEVER, IN MANY APPLICATIONS WHERE THERE IS A BROAD SPECTRUM OF NEUTRONS AND THE NEUTRON FLUX IS NOT HIGHLY DIRECTIONALLY ORIENTED, THE NEUTRON MULTIPLICATION, SPECTRUM AND ORIENTATION CAN BE FAIRLY ACCURATELY CALCULATED WITHOUT CONSIDERING CORRELATION EFFECTS.	Sixpak
	Sixpak
THE UNCORRELATED DATA PRODUCED BY THIS CODE REPLACES THE CORRELATED DATA,	Sixpak
	Sixpak
$F(E,EP,COS) = SIG(E)*Y(E)*G0(E,EP)*F(E,EP,COS)$	Sixpak
	Sixpak
BY THE UNCORRELATED DATA,	Sixpak
	Sixpak
$F(E,EP,COS) = SIG(E)*Y(E)*G0(E,EP)*F0(E,COS)$	Sixpak
	Sixpak
BY INTEGRATING $G0(E,EP)*F(E,EP,COS)$ OVER SECONDARY ENERGY (EP)	Sixpak

TO DEFINE AN AVERAGE ANGULAR DISTRIBUTION, $F_0(E, \cos)$ .	Sixpak
	Sixpak
WHAT IS LOST IN THIS PROCESS IS THE CORRELATION BETWEEN EP AND COS	Sixpak
SO THAT IN A TRANSPORT CALCULATION ALL MOMENTS OF THE FLUX WILL	Sixpak
HAVE THE SAME SPECTRUM, $G_0(E, EP)$ AND EACH WILL BE EFFECTED BY THE	Sixpak
AVERAGE ANGULAR DISTRIBUTION.	Sixpak
	Sixpak
FOR APPLICATIONS TO HIGH ENERGY FUSION APPLICATIONS CORRELATED	Sixpak
DATA SHOULD BE USED. HOWEVER, FOR LOWER ENERGY APPLICATIONS,	Sixpak
SUCH AS FISSION REACTORS, IT SHOULD BE ADEQUATE TO USE THE	Sixpak
UNCORRELATED DATA - IN THIS CASE THE MOST IMPORTANT EFFECT	Sixpak
WILL BE THE OVERALL NEUTRON MULTIPLICATION AND SPECTRUM.	Sixpak
	Sixpak
AN IMPORTANT CONSIDERATION IN DESIGNING THIS PROGRAM IS THAT	Sixpak
MANY COMPUTER CODES - DATA PROCESSING AND TRANSPORT CODES -	Sixpak
CANNOT USE THE CORRELATED (MF=6) DATA - NOR ARE THEY INTENDED	Sixpak
FOR HIGH ENERGY USE. FOR THESE CODES THE UNCORRELATED DATA	Sixpak
PRODUCED BY THIS CODE SHOULD BE ADEQUATE TO MEET THEIR NEEDS.	Sixpak
	Sixpak
WARNING - IT CANNOT BE STRESSED ENOUGH THAT THE OUTPUT OF THIS	Sixpak
CODE SHOULD ONLY BE USED FOR LOW ENERGY APPLICATIONS - FAILURE	Sixpak
TO HEED THIS WARNING CAN LEAD TO COMPLETELY UNRELIABLE RESULTS.	Sixpak
	Sixpak
ENDF/B FORMAT	Sixpak
=====	Sixpak
THIS PROGRAM ONLY USES THE ENDF/B BCD OR CARD IMAGE FORMAT (AS	Sixpak
OPPOSED TO THE BINARY FORMAT) AND CAN HANDLE DATA IN ANY VERSION	Sixpak
OF THE ENDF/B FORMAT (I.E., ENDF/B-I, II, III, IV, V OR VI FORMAT).	Sixpak
	Sixpak
IT IS ASSUMED THAT THE DATA IS CORRECTLY CODED IN THE ENDF/B	Sixpak
FORMAT AND NO ERROR CHECKING IS PERFORMED. IN PARTICULAR IT IS	Sixpak
ASSUMED THAT THE MAT, MF AND MT ON EACH LINE IS CORRECT. SEQUENCE	Sixpak
NUMBERS (COLUMNS 76-80) ARE IGNORED ON INPUT, BUT WILL BE	Sixpak
CORRECTLY OUTPUT ON ALL LINES. THE FORMAT OF SECTION MF=1, MT=451	Sixpak
AND ALL SECTIONS OF MF=6 MUST BE CORRECT. THE PROGRAM SKIPS ALL	Sixpak
OTHER SECTIONS OF DATA AND AS SUCH IS INSENSITIVE TO THE FORMAT	Sixpak
OF ALL OTHER SECTIONS.	Sixpak
	Sixpak
CONTENTS OF OUTPUT	Sixpak
=====	Sixpak
5 ENDF/B FORMATTED OUTPUT FILES ARE PRODUCED FOR NEUTRON INCIDENT	Sixpak
DATA,	Sixpak
	Sixpak
1) ENDFB.MF4 - ANGULAR DISTRIBUTIONS AND LEGENDRE COEFFICIENTS	Sixpak
FOR NEUTRONS	Sixpak
2) ENDFB.MF5 - TABULATED NEUTRON ENERGY SPECTRA	Sixpak
3) ENDFB.M12 - PHOTON EMISSION MULTIPLICITY	Sixpak
4) ENDFB.M14 - PHOTON EMISSION ANGULAR DISTRIBUTIONS (ALWAYS	Sixpak
ISOTROPIC)	Sixpak
5) ENDFB.M15 - TABULATED PHOTON EMISSION SPECTRA	Sixpak
	Sixpak
EMITTED PARTICLE YIELD	Sixpak
=====	Sixpak
NEUTRONS	Sixpak
=====	Sixpak
IN MF=6 THE YIELD FOR EACH REACTION IS THE ACTUAL MULTIPLICITY OF	Sixpak
THE REACTION, E.G., $(N, 2N) = 2$ . IN USING MF=4 AND 5 DATA THE	Sixpak
ENDF/B CONVENTION IS THAT THE MULTIPLICITY IS IMPLIED BY THE	Sixpak
MT NUMBER, E.G., $MT=16 = (N, 2N) = 2$ .	Sixpak
	Sixpak
THE ONLY EXCEPT IN ENDF/B-VI IS $MT=201 =$ TOTAL NEUTRON PRODUCTION	Sixpak
WHERE AN ACTUAL ENERGY DEPENDENT YIELD IS INCLUDED IN MF=6.	Sixpak
HOWEVER, IN THIS CASE THE MF=3 CROSS SECTION INCLUDES THE	Sixpak

MULTIPLICITY (S. PEARLSTEIN, PRIVATE COMMUNICATION, JAN. 1992),  
 $SIG(MT=201) = 2 * SIG(N, 2N) + 3 * SIG(N, 3N) \dots \text{ETC.}$

SO THAT FOR ALL ENDF/B-VI DATA AS OF JANUARY 1992 THE MF=4 AND 5  
 DATA OUTPUT BY THIS CODE CAN BE USED IN CONJUNCTION WITH THE MF=3  
 CROSS SECTIONS - WITHOUT ANY REFERENCE TO THE MF=6 YIELD.

PHOTONS  
 =====  
 UNLIKE THE NEUTRONS WHERE WITH ONLY ONE EXCEPTION (MT=201) THE  
 MF=6 YIELD IS ENERGY INDEPENDENT, IN THE CASE OF PHOTON EMISSION  
 ALMOST ALL OF THE PHOTONS HAVE AN ENERGY DEPENDENT YIELD.

THIS PROGRAM WILL OUTPUT THE PHOTON MULTIPLICITY IN MF=12 AND  
 INDICATE THAT THERE IS A NORMALIZED DISTRIBUTION IN MF=15  
 (LF=1 IN MF=12).

THIS PROGRAM WILL OUTPUT THE NORMALIZED PHOTON SPECTRA IN MF=15.  
 CONTINUOUS ENERGY SPECTRA AND DISCRETE PHOTONS WILL ALL BE OUTPUT  
 AS NORMALIZED SPECTRA.

THIS PROGRAM WILL ALSO OUTPUT MF=14 PHOTON ANGULAR DISTRIBUTION  
 DATA, ALWAYS USING THE ISOTROPIC FLAG TO MINIMIZE OUTPUT.

WARNING OF ENERGY DEPENDENT YIELD  
 =====  
 THIS PROGRAM WILL PRINT A WARNING MESSAGE IF A SECTION OF DATA  
 BEING OUTPUT IN THE ENDF/B FORMAT HAS AN ENERGY DEPENDENT MF=6  
 YIELD AND THE EMITTED PARTICLE IS A NEUTRON - SINCE THE ENDF/B  
 CONVENTION IS THAT FOR EACH MT NUMBER THE MULTIPLICITY IS IMPLIED  
 WE DO NOT EXPECT AN ENERGY DEPENDENT MULTIPLICITY FOR NEUTRON  
 EMISSION.

USING THE OUTPUT  
 =====  
 NOTE, THAT IN USING THIS DATA, STARTING FROM THE RELATIONSHIP,

$$F(E, EP, COS) = SIG(E) * Y(E) * G_0(E, EP) * F_0(E, COS)$$

USING THE ENDF/B CONVENTION THAT THE MULTIPLICITY IS EITHER  
 IMPLIED BY THE MT NUMBER (E.G., MT=16 = N, 2N - MULTIPLICITY = 2)  
 OR INCLUDED IN THE CROSS SECTION (E.G., MT=201 = TOTAL NEUTRON  
 PRODUCTION) ALL THE INFORMATION REQUIRED FOR A CALCULATION IS  
 AVAILABLE IN,

MF=3	- SIG(E)	
MF=4	- F <sub>0</sub> (E, COS)	- FOR OUTGOING NEUTRONS
MF=5	- G <sub>0</sub> (E, EP)	- FOR OUTGOING NEUTRONS
MF=12	- Y(E)	- FOR OUTGOING PHOTONS
MF=14	- F <sub>0</sub> (E, COS)	- FOR OUTGOING PHOTONS (ALWAYS ISOTROPIC)
MF=15	- G <sub>0</sub> (E, EP)	- FOR OUTGOING PHOTONS

DOCUMENTATION  
 =====  
 ONLY SECTIONS OF MF=4, 5, 12, 14, 15 ARE OUTPUT ON A ENDF/B FILE.  
 THE ONLY DOCUMENTATION IS THE ENDF/B TAPE LABEL (FIRST RECORD OF  
 EACH FILE) WHICH IDENTIFIES THE DATA AS SIXPAK OUTPUT.

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

SECTION SIZE	Sixpak
=====	Sixpak
ALL OF THE DATA IN ENDF/B-VI, MF=6 ARE QUITE SMALL TABLES. AS SUCH	Sixpak
THIS PROGRAM ONLY ALLOWS TABLES OF UP TO 12000 POINTS (12,000 X,	Sixpak
Y VALUES). THIS SIZE IS MORE THAN ADEQUATE TO HANDLE ALL OF THE	Sixpak
CURRENT ENDF/B-VI DATA, AND IT CAN BE EASILY INCREASED TO HANDLE	Sixpak
ANY NEWER DATA AS IT BECOMES AVAILABLE.	Sixpak
	Sixpak
PLEASE CONTACT THE AUTHOR IF YOU HAVE AN EVALUATION WHICH EXCEEDS	Sixpak
THIS LIMIT.	Sixpak
	Sixpak
SELECTION OF DATA	Sixpak
=====	Sixpak
THE PROGRAM SELECTS DATA TO BE PROCESSED BASED ON MAT/MT RANGES	Sixpak
(MF=6 ASSUMED). THIS PROGRAM ALLOWS UP TO 100 MAT/MT RANGES TO BE	Sixpak
SPECIFIED BY INPUT PARAMETERS. THE PROGRAM WILL ASSUME THAT THE	Sixpak
ENDF/B TAPE IS IN MAT ORDER. THE PROGRAM WILL TERMINATE EXECUTION	Sixpak
WHEN A MAT IS FOUND THAT IS ABOVE ALL REQUESTED MAT RANGES.	Sixpak
	Sixpak
PROGRAM OPERATION	Sixpak
=====	Sixpak
EACH SECTION (MT) OF MF=6 DATA IS SUBDIVIDED INTO SUBSECTIONS -	Sixpak
ONE SUBSECTION FOR EACH EMITTED PARTICLE.	Sixpak
	Sixpak
EACH SUBSECTION OF DATA IS CONSIDERED SEPARATELY. EACH SUBSECTION	Sixpak
OF ENDF/B MF=6 DATA TO PROCESS IS IN THE FORM,	Sixpak
	Sixpak
$F(E,EP,COS) = SIG(E)*Y(E)*G0(E,EP)*F(E,EP,COS)$	Sixpak
	Sixpak
SIG(E) = MF=3 CROSS SECTIONS	Sixpak
Y(E) = YIELD (MULTIPLICITY)	Sixpak
G0(E,EP) = ENERGY SPECTRUM	Sixpak
F(E,EP,COS) = ANGULAR DISTRIBUTION	Sixpak
	Sixpak
G0(E,EP) = 1 WHEN INTEGRATED OVER EP (SECONDARY ENERGY)	Sixpak
G0(E,EP)*F(E,EP,COS) = 1 WHEN INTEGRATED OVER EP AND COS	Sixpak
	Sixpak
THIS PROGRAM WILL DEFINE THE ZEROth ORDER MOMENTS OF THE	Sixpak
ENERGY AND ANGULAR DISTRIBUTIONS,	Sixpak
	Sixpak
G0(E,EP) = G0(E,EP)*F(E,EP,COS) INTEGRATED OVER COS	Sixpak
F0(E,COS) = G0(E,EP)*F(E,EP,COS) INTEGRATED OVER EP	Sixpak
	Sixpak
FOR NEUTRON INDUCED REACTIONS THE ENDF/B FORMATTED OUTPUT WILL BE	Sixpak
	Sixpak
F0(E,COS)- IN ENDFB.MF4 FOR NEUTRONS OUT OF A REACTION	Sixpak
G0(E,EP) - IN ENDFB.MF5 FOR NEUTRONS OUT OF A REACTION	Sixpak
- IN ENDFB.M15 FOR PHOTONS OUT OF A REACTION	Sixpak
	Sixpak
FOR NEUTRONS INCIDENT AND NEUTRONS EMITTED THIS DATA WILL BE	Sixpak
OUTPUT IN MF=4 AND 5 FORMATS.	Sixpak
	Sixpak
FOR NEUTRONS INCIDENT AND PHOTONS EMITTED THIS DATA WILL BE	Sixpak
OUTPUT IN MF=15 FORMAT - THE SPECTRA ARE OUTPUT AND THE	Sixpak
ANGULAR DISTRIBUTION IS IGNORED.	Sixpak
	Sixpak
ALL PHOTON EMISSION IN THE ENDF/B-VI LIBRARY AS OF JANUARY 1992	Sixpak
IS ISOTROPIC AND AS SUCH NO DISTRIBUTION OF PHOTON ANGULAR	Sixpak
DISTRIBUTIONS NEED BE OUTPUT - IT IS ALWAYS ISOTROPIC.	Sixpak
	Sixpak
FOR ALL OTHER COMBINATIONS INCIDENT AND EMITTED PARTICLES	Sixpak
THERE WILL BE NO ENDF/B FORMATTED OUTPUT.	Sixpak
	Sixpak

VARIATIONS FROM ENDF/B MANUAL	Sixpak
=====	Sixpak
LAW=1, LANG=2 = KALBACH-MANN	Sixpak
=====	Sixpak
FOR THE DISTRIBUTIONS,	Sixpak
	Sixpak
$F(\mu, E, EP) = G_0(E, EP) * A * (\cosh(\mu * A) + R(E, EP) * \sinh(\mu * A))$	Sixpak
	Sixpak
$G_0(E, EP) = 1 - \text{WHEN INTEGRATED OVER } EP.$	Sixpak
	Sixpak
$A * (\cosh(\mu * A) + R(E, EP) * \sinh(\mu * A)) = 2 - \text{WHEN INTEGRATED OVER } \mu$	Sixpak
	Sixpak
THIS MEANS AS DEFINED IN THE ENDF/B MANUAL THE DISTRIBUTIONS	Sixpak
ARE NORMALIZED TO 2, INSTEAD OF 1. IN ORDER TO OBTAIN CORRECTLY	Sixpak
NORMALIZED DISTRIBUTIONS THE DISTRIBUTION SHOULD BE DEFINED	Sixpak
TO INCLUDE A FACTOR OF 1/2 MULTIPLYING THE ANGULAR PART OF	Sixpak
THE DISTRIBUTION.	Sixpak
	Sixpak
$F(\mu, E, EP) = G_0(E, EP) * 0.5 * A * (\cosh(\mu * A) + R(E, EP) * \sinh(\mu * A))$	Sixpak
	Sixpak
THIS IS THE FORM USED IN THIS CODE	Sixpak
	Sixpak
LAW=1, ND NOT 0 = DISCRETE SECONDARY ENERGY DISTRIBUTION	Sixpak
=====	Sixpak
THE ENDF/B MANUAL SAYS THESE ARE FLAGGED WITH NEGATIVE ENERGIES.	Sixpak
IN ENDF/B-VI ALL OF THESE HAVE POSITIVE ENERGY. THIS CODE DOES	Sixpak
NOT CONSIDER THE ENDF/B-VI DATA TO BE IN ERROR.	Sixpak
	Sixpak
WITH THE CONVENTION ACTUALLY USED IN ENDF/B-VI ALL SECONDARY	Sixpak
ENERGIES SHOULD BE NON-NEGATIVE AND IN ASCENDING ENERGY ORDER	Sixpak
FOR EACH INCIDENT ENERGY.	Sixpak
	Sixpak
FROM THE ENDF/B MANUAL IT IS NOT OBVIOUS WHAT $G_0(E, EP)$ SHOULD BE	Sixpak
FOR DISCRETE PHOTONS - PHYSICALLY THIS IS A DELTA FUNCTION. IN	Sixpak
ENDF/B-VI IT IS ENTERED AS 1.0 = INTERPRETING IT AS INTEGRATED	Sixpak
	Sixpak
OVER SECONDARY ENERGY - IN WHICH CASE THE DELTA FUNCTION = 1.0.	Sixpak
	Sixpak
LIMITATIONS	Sixpak
=====	Sixpak
CHECKING DATA	Sixpak
=====	Sixpak
THIS PROGRAM CHECKS ALL ENDF/B-VI MF=6 DATA. THE FOLLOWING CHECKS	Sixpak
ARE PERFORMED.	Sixpak
	Sixpak
PARAMETERS	Sixpak
=====	Sixpak
ALL PARAMETERS ARE CHECKED FOR CONSISTENCY. IF PARAMETERS ARE	Sixpak
NOT CONSISTENT THE PROGRAM MAY NOT BE ABLE TO PERFORM THE	Sixpak
FOLLOWING TESTS AND WILL MERELY SKIP A SECTION OF DATA.	Sixpak
	Sixpak
INTERPOLATION LAWS	Sixpak
=====	Sixpak
ALL INTEGRATIONS ARE PERFORMED USING THE INTERPOLATION LAW GIVEN	Sixpak
FOR SECONDARY ENERGY AND/OR COSINE. INTEGRATIONS ARE NOT	Sixpak
PERFORMED OVER INCIDENT - ONLY INTEGRATION OVER SECONDARY ENERGY	Sixpak
AND/OR COSINE ARE PERFORMED AT EACH INCIDENT ENERGY. THEREFORE	Sixpak
THE INTERPOLATION LAW FOR INCIDENT ENERGY IS NOT USED BY THIS	Sixpak
CODE.	Sixpak
	Sixpak
ALL INTERPOLATION LAWS ARE CHECKED. ALL DATA ASSOCIATED WITH	Sixpak
INTERPOLATION LAWS ARE CHECKED, E.G., NO NON-NEGATIVE VALUES	Sixpak
REQUIRING LOG INTERPOLATION. IN ORDER TO PERFORM REQUIRED	Sixpak



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## =====

EITHER TABULATED OR LEGENDRE COEFFICIENTS

## YIELD AND OUTPUT NORMALIZATION

LAW=0

LAW=1

LAW=2

LAW=3

=====

NO LIMITATION ON REPRESENTATIONS.	Sixpak
	Sixpak
LAW=4	Sixpak
=====	Sixpak
NO OUTPUT - INCIDENT NEUTRON - EMITTED PHOTON OR NEUTRON	Sixpak
REACTIONS ARE NOT EXPECTED.	Sixpak
	Sixpak
LAW=5	Sixpak
=====	Sixpak
NO OUTPUT - INCIDENT NEUTRON - EMITTED PHOTON OR NEUTRON	Sixpak
REACTIONS ARE NOT EXPECTED.	Sixpak
	Sixpak
LAW=6	Sixpak
=====	Sixpak
NO OUTPUT - ENDF/B-VI ONLY INCLUDES 1 SECTION OF THIS TYPE OF DATA	Sixpak
FOR (N,D) 2N,P.	Sixpak
	Sixpak
LAW=7	Sixpak
=====	Sixpak
FOR EACH INCIDENT ENERGY THE REPRESENTATION MUST BE EITHER,	Sixpak
	Sixpak
1) SQUARE = FOR EACH INCIDENT COSINE EXACTLY THE SAME SECONDARY	Sixpak
ENERGIES.	Sixpak
	Sixpak
2) LINEAR = FOR EACH INCIDENT COSINE THE INTERPOLATION LAW	Sixpak
BETWEEN SECONDARY ENERGIES MUST BE LINEAR.	Sixpak
	Sixpak
THESE 2 PRESENTATIONS ARE THE ONLY ONES PRESENTED IN ENDF/B-VI	Sixpak
AS OF JANUARY 1992 - SO THIS PROGRAM CAN TRANSLATED ALL LAW=7	Sixpak
DATA FOR ENDF/B-VI.	Sixpak
	Sixpak
LABORATORY VS. CENTER-OF-MASS SYSTEM	Sixpak
=====	Sixpak
IN MANY CASES PEOPLE ASSUME THAT FOR HEAVY (HIGH ATOMIC WEIGHT)	Sixpak
MATERIALS THE CENTER-OF-MASS AND LAB SYSTEMS ARE ALMOST IDENTICAL,	Sixpak
SINCE IN THIS CASE THE CENTER-OF-MASS ENERGY WILL BE MUCH SMALLER	Sixpak
THAN THE INCIDENT ENERGY. FOR A PROCESS SUCH AS ELASTIC SCATTERING	Sixpak
WHERE FOR HEAVY MATERIALS THE SECONDARY ENERGY, EP, WILL ALWAYS	Sixpak
BE A LARGE FRACTION OF THE INCIDENT ENERGY, THIS ASSUMPTION IS	Sixpak
VALID. HOWEVER, FOR THE TYPICAL REACTIONS INCLUDED IN MF=6 THIS	Sixpak
IS NOT ALWAYS TRUE - IN MANY OF THESE CASES THE SECONDARY ENERGY	Sixpak
CAN EXTEND ALL THE WAY DOWN TO ZERO, AND IN PARTICULAR IT CAN	Sixpak
BE SMALL COMPARED TO THE CENTER-OF-MASS ENERGY - WHICH MAKES THE	Sixpak
TRANSFORMATION FROM CENTER-OF-MASS TO LAB IMPORTANT. THEREFORE	Sixpak
GENERALLY TO TREAT MF=6 DATA WE MUST CONSIDER THIS TRANSFORMATION.	Sixpak
	Sixpak
THE FOLLOWING DISCUSSING ONLY APPLIES TO SPECTRA THAT MAY BE	Sixpak
OUTPUT IN MF=5 = ONLY DATA FOR NEUTRONS INCIDENT AND EMITTED -	Sixpak
IN PARTICULAR THE FOLLOWING DEFINITIONS ARE NOT GENERAL - THEY	Sixpak
ARE ONLY VALID FOR INCIDENT AND EMITTED NEUTRONS.	Sixpak
	Sixpak
DOUBLE DIFFERENTIAL DATA IN MF=6 MAY BE GIVEN IN EITHER THE LAB	Sixpak
OR C.M. SYSTEM. SIMILARLY ANGULAR DISTRIBUTIONS IN MF=4 MAY BE	Sixpak
GIVEN IN EITHER THE LAB OR C.M. SYSTEM. IN CONTRAST ENERGY	Sixpak
SPECTRA IN MF=5 CAN ONLY BE GIVEN IN THE LABORATORY SYSTEM.	Sixpak
	Sixpak
THE ANGULAR DISTRIBUTIONS OUTPUT BY THIS CODE IN MF=4 ARE IN THE	Sixpak
SAME SYSTEM IN WHICH THEY ARE GIVEN IN MF=6 - EITHER LAB OR	Sixpak
CENTER-OF-MASS SYSTEM.	Sixpak
	Sixpak
THE ENERGY SPECTRA OUTPUT BY THIS CODE IN MF=5 MUST BE IN THE LAB	Sixpak
SYSTEM - THIS IS THE ONLY ALLOWED FORM FOR MF=5 DATA.	Sixpak
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$$\text{SQRT}[V(\text{MM})^2 + V(\text{CM})^2 + 2 \cdot \cos(\text{CM}) \cdot V(\text{MM}) \cdot V(\text{CM})]$$

OR  $\cos(\text{CM})$  FROM THE RELATIONSHIP,

$$V(\text{CM}) \cdot \cos(\text{CM}) = V(\text{LAB}) \cdot \cos(\text{LAB}) - V(\text{MM})$$

$$\cos(\text{CM}) = [V(\text{LAB}) \cdot \cos(\text{LAB}) - V(\text{MM})] / V(\text{CM})$$

$$\cos(\text{CM}) = \frac{[V(\text{LAB}) \cdot \cos(\text{LAB}) - V(\text{MM})]}{\text{SQRT}[V(\text{LAB})^2 + V(\text{CM})^2 - 2 \cdot \cos(\text{LAB}) \cdot V(\text{LAB}) \cdot V(\text{MM})]}$$

THE JACOBIAN CAN BE DEFINED FROM,

$$V(\text{LAB}) \cdot \cos(\text{LAB}) = V(\text{MM}) + V(\text{CM}) \cdot \cos(\text{CM})$$

$$J = D[\cos(\text{CM})] / D[\cos(\text{LAB})] = V(\text{LAB}) / V(\text{CM})$$

$$= \text{SQRT}[EP(\text{LAB}) / EP(\text{CM})]$$

WITH THESE DEFINITIONS OF  $EP(\text{LAB})$  AND  $\cos(\text{LAB})$  IN TERMS OF  $E(\text{MM})$ ,  $EP(\text{CM})$  AND  $\cos(\text{CM})$  IT IS POSSIBLE TO PERFORM A POINT-BY-POINT TRANSFORMATION OF DISTRIBUTIONS FROM THE CM TO LAB SYSTEM USING THESE DEFINITIONS - OR IF WE WISHED WE COULD PERFORM THE REVERSE TRANSFORMATION USING THE ABOVE RELATIONSHIPS AND THE IDENTITY,

$$F(E, EP(\text{LAB}), \cos(\text{LAB})) \cdot D(\cos(\text{LAB})) = F(E, EP(\text{CM}), \cos(\text{CM})) \cdot D(\cos(\text{CM}))$$

THIS IS NOT WHAT WILL BE DONE HERE, SINCE WE WILL ONLY BE INTERESTED IN THE ZEROth ORDER MOMENTS OF THESE DISTRIBUTIONS, BUT WE WILL BE INTERESTED IN DEFINING THOSE MOMENTS IN THE LAB SYSTEM IN TERMS OF MF=6 SPECTRA GIVEN IN THE CM SYSTEM USING,

$$F(E, EP(\text{LAB}), \cos(\text{LAB})) = F(E, EP(\text{CM}), \cos(\text{CM})) \cdot J$$

THE LIMITS OF  $EP(\text{LAB})$  ARE DEFINED BY SETTING  $\cos(\text{CM}) = +1$  OR  $-1$ ,

$$EP(\text{LAB}) = (\text{SQRT}(EP(\text{CM})) + \text{SQRT}(E(\text{MM})))^2 \quad \text{FOR } \cos(\text{CM}) = +1$$

$$= (\text{SQRT}(EP(\text{CM})) - \text{SQRT}(E(\text{MM})))^2 \quad \text{FOR } \cos(\text{CM}) = -1$$

IN THIS FORM WE CAN SEE THAT AS LONG AS THE SECONDARY ENERGY IN THE CENTER-OF-MASS SYSTEM,  $EP(\text{CM})$ , IS MUCH LARGER THAN THE ENERGY OF THE CENTER-OF-MASS,  $E(\text{MM})$ , THE CENTER-OF-MASS AND LAB ENERGIES WILL BE ALMOST EQUAL - SIMILARLY FOR THE COSINE, IN THIS CASE  $\cos(\text{LAB})$  AND  $\cos(\text{CM})$  WILL BE ALMOST EQUAL - HOWEVER, FOR THE MF=6 DATA WE CANNOT ASSUME THAT THIS IS TRUE.

TO FIRST ORDER THE ANGULAR DEPENDENCE CAN BE IGNORED,

$$EP(\text{LAB}) = E(\text{MM}) + EP(\text{CM})$$

ALL THIS SAYS IS THAT TO FIRST ORDER THE EFFECT OF TRANSFORMING FROM THE CM TO LAB SYSTEM IS TO INCREASE THE ENERGY OF THE EMITTED PARTICLE IN THE CENTER-OF-MASS SYSTEM BY THE ENERGY OF THE CENTER-OF-MASS TO DEFINE THE LAB ENERGY.

NOT ONLY THE ENERGY, BUT ALSO THE SPECTRA MUST BE TRANSFORMED. STARTING FROM THE DOUBLE DIFFERENTIAL DATA IN THE LAB SYSTEM,  $F(E, EP, \cos(\text{LAB}))$ , WE CAN DEFINE THE LAB SCALAR SPECTRUM AS,

$$GO(E, EP) = \text{INTEGRAL } F(E, EP, \cos(\text{LAB})) \cdot D(\cos(\text{LAB}))$$

THIS IS THE NORMAL CALCULATION DEFINED ABOVE AND USED FOR DATA GIVEN IN THE LAB SYSTEM.



THIS CODE. USERS ARE ASKED NOT TO ACTIVATE OR TRY TO USE THESE  
ROUTINES. UNLESS YOU COMPLETELY UNDERSTAND THIS CODE THE RESULTS  
CAN BE UNRELIABLE IF YOU ACTIVATE THESE ROUTINES.

#### INPUT FILES

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

#### OUTPUT FILES

```
=====
UNIT  DESCRIPTION
----  -
      3  OUTPUT REPORT (BCD - 120 CHARACTERS/RECORD)
      11 ENDF/B DATA MF=4 (BCD - 80 CHARACTERS/RECORD)
      12 ENDF/B DATA MF=5 (BCD - 80 CHARACTERS/RECORD)
      14 ENDF/B DATA MF=15 (BCD - 80 CHARACTERS/RECORD)
      17 ENDF/B DATA MF=12 (BCD - 80 CHARACTERS/RECORD)
      18 ENDF/B DATA MF=14 (BCD - 80 CHARACTERS/RECORD)
      15 PLOTTAB INPUT PARAMETERS (BCD - 80 CHARACTERS/RECORD)
      16 PLOTTAB FORMATTED OUTPUT (BCD - 80 CHARACTERS/RECORD)
```

#### SCRATCH FILES

```
=====
NONE
```

#### OPTIONAL STANDARD FILE NAMES (SEE SUBROUTINE FILIO1 AND FILIO2)

```
=====
UNIT  FILE NAME
----  -
      2  SIXPAK.INP
      3  SIXPAK.LST
      10 ENDFB.IN
      11 ENDFB.MF4
      12 ENDFB.MF5
      14 ENDFB.M15
      17 ENDFB.M12
      18 ENDFB.M14
      15 PLOTTAB.INP
      16 PLOTTAB.CUR
```

#### INPUT PARAMETERS

```
=====
LINE  COLS.  DESCRIPTION
----  -
      1  1-60  ENDF/B INPUT DATA FILENAME
              (STANDARD OPTION = ENDFB.IN)
      2-N  1-6  MINIMUM MAT FOR REQUESTED RANGE
              9-11 MINIMUM MT FOR REQUESTED RANGE
              12-17 MAXIMUM MAT FOR REQUESTED RANGE
              20-22 MAXIMUM MT FOR REQUESTED RANGE
```

LEAVE THE DEFINITION OF THE FILENAME BLANK - THE PROGRAM WILL  
THEN USE THE STANDARD FILENAME (ENDFB.IN).

UP TO 100 MAT/MT RANGES MAY BE SPECIFIED. THE LIST OF RANGES IS  
TERMINATED BY A BLANK LINE. IF THE FIRST INPUT LINE IS COMPLETELY  
BLANK ALL DATA WILL BE PROCESSED.

EXAMPLE INPUT NO. 1

```

=====
PROCESS ALL MF=6 DATA ON AN ENDF/B TAPE. USE THE STANDARD INPUT
DATA FILENAME ENDFB.IN IN THIS CASE THE USER CAN EITHER EXPLICITLY
SPECIFY THE FILENAME AND MAT/MT RANGE BY THE FOLLOWING 2 INPUT
LINES,

ENDFB.IN
      1      1  9999  999
                                (BLANK LINE, TERMINATES REQUEST LIST)

OR BY INPUTTING 2 BLANK LINE = PROCESS EVERYTHING.

EXAMPLE INPUT NO. 2
-----
PROCESS BE-9, MAT=425, MT=16. READ THE DATA FROM ENDFB6\BE9.
IN THIS CASE THE FOLLOWING 3 INPUT LINES ARE REQUIRED,

ENDFBB6\BE9
      425      16      425      16
                                (BLANK LINE, TERMINATES REQUEST LIST)

EXAMPLE INPUT NO. 3
-----
PROCESS ALL MT=16 (N,2N) DATA. THIS CAN BE DONE BY SPECIFYING THE
MAXIMUM MAT RANGE = 1 TO 9999, AND MT=16 FOR THE MINIMUM AND
MAXIMUM MT RANGE. READ THE DATA FROM ENDFB6\K300. IN THIS CASE
CASE THE FOLLOWING 3 INPUT LINES ARE REQUIRED,

ENDFB6\K300
      1      16  9999      16
                                (BLANK LINE, TERMINATES REQUEST LIST)

=====

```