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===== 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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===== Sixpak
THE AUTHOR THANKS SOL PEARLSTEIN (BROOKHAVEN NATIONAL LAB) FOR Sixpak
SIGNIFICANTLY CONTRIBUTING TOWARD IMPROVING THE ACCURACY AND Sixpak
COMPUTER INDEPENDENCE OF THIS CODE - THANKS, SOL Sixpak
Sixpak
ACKNOWLEDGEMENT (VERSION 92-4) Sixpak
===== Sixpak
THE AUTHOR THANKS BOB MACFARLANE (LOS ALAMOS) FOR SUGGESTING HOW Sixpak
TO PROPERLY OUTPUT THE PHOTON PRODUCTION DATA TO PUT IT INTO Sixpak
EXACTLY THE FORM NEEDED FOR USE IN PROCESSING CODES. Sixpak
Sixpak
THE AUTHOR THANKS CHRIS DEAN (WINFRITH) FOR POINTING OUT ERRORS Sixpak
IN THE EARLIER TREATMENT OF THE KALBACH-MANN FORMALISM AND IN Sixpak
THE DEFINITION OF THE ISOTROPIC ANGULAR DISTRIBUTION FLAG (LI). Sixpak
Sixpak
AUTHORS MESSAGE Sixpak
===== Sixpak
THE COMMENTS BELOW SHOULD BE CONSIDERED THE LATEST DOCUMENTATION Sixpak
INCLUDING ALL RECENT IMPROVEMENTS. PLEASE READ ALL OF THESE Sixpak
COMMENTS BEFORE IMPLEMENTING AND USING THESE CODES. Sixpak
Sixpak
AT THE PRESENT TIME WE ARE ATTEMPTING TO DEVELOP A SET OF COMPUTER Sixpak
INDEPENDENT PROGRAMS THAT CAN EASILY BE IMPLEMENTED ON ANY ONE Sixpak
OF A WIDE VARIETY OF COMPUTERS. IN ORDER TO ASSIST IN THIS PROJECT Sixpak
IT WOULD BE APPRECIATED IF YOU WOULD NOTIFY THE AUTHOR OF ANY Sixpak
COMPILER DIAGNOSTICS, OPERATING PROBLEMS OR SUGGESTIONS ON HOW TO Sixpak
IMPROVE THIS PROGRAM. HOPEFULLY, IN THIS WAY FUTURE VERSIONS OF Sixpak
THIS PROGRAM WILL BE COMPLETELY COMPATIBLE FOR USE ON YOUR Sixpak
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. Sixpak
IT IS MERELY CHECKED. IF ERRORS ARE FOUND IT IS UP TO THE USER Sixpak
TO TAKE CORRECTIVE ACTION ON THE MF=6 DATA. Sixpak
Sixpak
IN CONTRAST WHEN PROBLEMS ARE FOUND IN DATA WHICH WILL BE OUTPUT Sixpak
IN THE ENDF/B FORMAT (MF=4, 5, 12, 14 AND 15), WHENEVER POSSIBLE Sixpak
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 Sixpak
FURTHER CORRECTIVE ACTION CAN BE TAKEN AS FOLLOWS, Sixpak
Sixpak
PROGRAM LEGEND Sixpak
===== Sixpak
CAN BE USED TO CORRECT ANGULAR DISTRIBUTIONS WHICH ARE NEGATIVE, Sixpak
TO CONVERT FROM LEGENDRE COEFFICIENTS TO TABULATED ANGULAR Sixpak
DISTRIBUTIONS AND GENERALLY PERFORM MORE EXTENSIVE TESTS OF Sixpak
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 Sixpak

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CONTENTS OF OUTPUT	Sixpak
=====	Sixpak
5 ENDF/B FORMATTED OUTPUT FILES ARE PRODUCED FOR NEUTRON INCIDENT DATA,	Sixpak
1) ENDFB.MF4 - ANGULAR DISTRIBUTIONS AND LEGENDRE COEFFICIENTS 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 ISOTROPIC)	Sixpak
5) ENDFB.M15 - TABULATED PHOTON EMISSION SPECTRA	Sixpak
EMITTED PARTICLE YIELD	Sixpak
=====	Sixpak
NEUTRONS	Sixpak
=====	Sixpak
IN MF=6 THE YIELD FOR EACH REACTION IS THE ACTUAL MULTIPLICITY OF THE REACTION, E.G., (N,2N) = 2. IN USING MF=4 AND 5 DATA THE ENDF/B CONVENTION IS THAT THE MULTIPLICITY IS IMPLIED BY THE MT NUMBER, E.G., MT=16 = (N,2N) = 2.	Sixpak
THE ONLY EXCEPT IN ENDF/B-VI IS MT=201 = TOTAL NEUTRON PRODUCTION WHERE AN ACTUAL ENERGY DEPENDENT YIELD IS INCLUDED IN MF=6. HOWEVER, IN THIS CASE THE MF=3 CROSS SECTION INCLUDES THE MULTIPLICITY (S. PEARLSTEIN, PRIVATE COMMUNICATION, JAN. 1992), SIG(MT=201) = 2*SIG(N,2N)+3*SIG(N,3N).....ETC.	Sixpak
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.	Sixpak
PHOTONS	Sixpak
=====	Sixpak
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.	Sixpak
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).	Sixpak
THIS PROGRAM WILL OUTPUT THE NORMALIZED PHOTON SPECTRA IN MF=15. CONTINUOUS ENERGY SPECTRA AND DISCRETE PHOTONS WILL ALL BE OUTPUT AS NORMALIZED SPECTRA.	Sixpak
THIS PROGRAM WILL ALSO OUTPUT MF=14 PHOTON ANGULAR DISTRIBUTION DATA, ALWAYS USING THE ISOTROPIC FLAG TO MINIMIZE OUTPUT.	Sixpak
WARNING OF ENERGY DEPENDENT YIELD	Sixpak
=====	Sixpak
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.	Sixpak
USING THE OUTPUT	Sixpak
=====	Sixpak
NOTE, THAT IN USING THIS DATA, STARTING FROM THE RELATIONSHIP,	Sixpak
	Sixpak

$F(E, EP, COS) = SIG(E) * Y(E) * G0(E, EP) * F0(E, COS)$ Sixpak

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, Sixpak

MF=3 - SIG(E) Sixpak
MF=4 - F0(E, COS) - FOR OUTGOING NEUTRONS Sixpak
MF=5 - G0(E, EP) - FOR OUTGOING NEUTRONS Sixpak
MF=12 - Y(E) - FOR OUTGOING PHOTONS Sixpak
MF=14 - F0(E, COS) - FOR OUTGOING PHOTONS (ALWAYS ISOTROPIC) Sixpak
MF=15 - G0(E, EP) - FOR OUTGOING PHOTONS Sixpak

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

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

SECTION SIZE
===== Sixpak
ALL OF THE DATA IN ENDF/B-VI, MF=6 ARE QUITE SMALL TABLES. AS SUCH
THIS PROGRAM ONLY ALLOWS TABLES OF UP TO 12000 POINTS (12,000 X,
Y VALUES). THIS SIZE IS MORE THAN ADEQUATE TO HANDLE ALL OF THE
CURRENT ENDF/B-VI DATA, AND IT CAN BE EASILY INCREASED TO HANDLE
ANY NEWER DATA AS IT BECOMES AVAILABLE. Sixpak

PLEASE CONTACT THE AUTHOR IF YOU HAVE AN EVALUATION WHICH EXCEEDS
THIS LIMIT. Sixpak

SELECTION OF DATA
===== Sixpak
THE PROGRAM SELECTS DATA TO BE PROCESSED BASED ON MAT/MT RANGES
(MF=6 ASSUMED). THIS PROGRAM ALLOWS UP TO 100 MAT/MT RANGES TO BE
SPECIFIED BY INPUT PARAMETERS. THE PROGRAM WILL ASSUME THAT THE
ENDF/B TAPE IS IN MAT ORDER. THE PROGRAM WILL TERMINATE EXECUTION
WHEN A MAT IS FOUND THAT IS ABOVE ALL REQUESTED MAT RANGES. Sixpak

PROGRAM OPERATION
===== Sixpak
EACH SECTION (MT) OF MF=6 DATA IS SUBDIVIDED INTO SUBSECTIONS -
ONE SUBSECTION FOR EACH EMITTED PARTICLE. Sixpak

EACH SUBSECTION OF DATA IS CONSIDERED SEPARATELY. EACH SUBSECTION
OF ENDF/B MF=6 DATA TO PROCESS IS IN THE FORM, Sixpak

$F(E, EP, COS) = SIG(E) * Y(E) * G0(E, EP) * F(E, EP, COS)$ 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

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 ENERGY AND ANGULAR DISTRIBUTIONS, Sixpak
Sixpak
Sixpak
 $G_0(E,EP) = G_0(E,EP) * F(E,EP,COS)$ INTEGRATED OVER COS Sixpak
 $F_0(E,COS) = G_0(E,EP) * F(E,EP,COS)$ INTEGRATED OVER EP Sixpak
Sixpak
FOR NEUTRON INDUCED REACTIONS THE ENDF/B FORMATTED OUTPUT WILL BE Sixpak
Sixpak
 $F_0(E,COS)$ - IN ENDFB.MF4 FOR NEUTRONS OUT OF A REACTION Sixpak
 $G_0(E,EP)$ - 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
 $G_0(E,EP) = 1$ - WHEN INTEGRATED OVER EP. Sixpak
Sixpak
 $A * (COSH(MU*A) + R(E,EP) * SINH(MU*A)) = 2$ - 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
OVER SECONDARY ENERGY - IN WHICH CASE THE DELTA FUNCTION = 1.0. Sixpak
Sixpak

$$EP(LAB) = 0.5 * MASS(OUT) * V(LAB) ** 2$$

$$= E(MM) + EP(CM) + 2 * COS(CM) * SQRT(E(MM) * EP(CM))$$

WE CAN ALSO DEFINE THE REVERSE TRANSFORMATION USING,

$$V(CM) * COS(CM) = V(LAB) * COS(LAB) - V(MM)$$

$$V(CM) * SIN(CM) = V(LAB) * SIN(LAB)$$

$$V(CM) ** 2 = V(MM) ** 2 + V(LAB) ** 2 - 2 * COS(LAB) * V(MM) * V(LAB)$$

$$EP(CM) = 0.5 * MASS(OUT) * V(CM) ** 2$$

$$= E(MM) + EP(LAB) - 2 * COS(LAB) * SQRT(E(MM) * EP(LAB))$$

WE CAN DEFINE COS(LAB) FROM THE RELATIONSHIP,

$$V(LAB) * COS(LAB) = V(MM) + V(CM) * COS(CM)$$

$$COS(LAB) = [V(MM) + V(CM) * COS(CM)] / V(LAB)$$

$$COS(LAB) = \frac{[V(MM) + V(CM) * COS(CM)]}{SQRT[V(MM) ** 2 + V(CM) ** 2 + 2 * COS(CM) * V(MM) * V(CM)]}$$

OR COS(CM) FROM THE RELATIONSHIP,

$$V(CM) * COS(CM) = V(LAB) * COS(LAB) - V(MM)$$

$$COS(CM) = [V(LAB) * COS(LAB) - V(MM)] / V(CM)$$

$$COS(CM) = \frac{[V(LAB) * COS(LAB) - V(MM)]}{SQRT[V(LAB) ** 2 + V(CM) ** 2 - 2 * COS(LAB) * V(LAB) * V(MM)]}$$

THE JACOBIAN CAN BE DEFINED FROM,

$$V(LAB) * COS(LAB) = V(MM) + V(CM) * COS(CM)$$

$$J = D[COS(CM)] / D[COS(LAB)] = V(LAB) / V(CM)$$

$$= SQRT[EP(LAB) / EP(CM)]$$

WITH THESE DEFINITIONS OF EP(LAB) AND COS(LAB) IN TERMS OF E(MM), EP(CM) AND COS(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,

$$F(E, EP(LAB), COS(LAB)) * D(COS(LAB)) = F(E, EP(CM), COS(CM)) * D(COS(CM))$$

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,

$$F(E, EP(LAB), COS(LAB)) = F(E, EP(CM), COS(CM)) * J$$

THE LIMITS OF EP(LAB) ARE DEFINED BY SETTING COS(CM) = +1 OR -1,

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

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

BY E(MM) TO ACCOUNT FOR THE CENTER OF MASS MOTION - THE SPECTRA WILL NOT BE MODIFIED BY THE JACOBIAN FACTOR $\sqrt{EP(LAB)/EP(CM)}$ SINCE THIS WOULD REQUIRE A DETAILED TRANSFORMATION IN ENERGY AND COS(THETA) SPACE - WHICH IS JUDGED NOT TO BE WORTH PERFORMING WITHIN THE LIMITS OF WHERE THE OUTPUT FROM THIS CODE IS INTENDED TO BE USED.

SINCE THE ANGULAR DISTRIBUTION IS ALWAYS OUTPUT IN THE SAME SYSTEM AS WHICH IT IS GIVEN IN MF=6, NO TRANSFORMATION IS REQUIRED FOR THE MF=4 OUTPUT.

WHEN USED IN LOW ENERGY APPLICATIONS (E.G., FISSION REACTORS) THE HIGH ENERGY SPECTRA PRESENTED IN MF=6 WILL BE MOSTLY IMPORTANT SIMPLY IN CONSERVING PARTICLES, (E.G., AS IN (N,2N)) AND ENERGY AND THE DETAILS OF THE CORRELATION AND GROSS ENERGY SPECTRA WILL NOTE PLAY THAT IMPORTANT A ROLE. IN THIS CASE THE SPECTRA OUTPUT BY THIS PROGRAM IN MF=5 SHOULD BE ADEQUATE.

PLOTTAB FORMATTED OUTPUT

THIS PROGRAM CONTAINS ROUTINES TO PRODUCE OUTPUT THAT CAN BE USED AS INPUT TO THE PLOTTAB CODE TO OBTAIN GRAPHIC RESULTS.

THESE ROUTINES ARE DESIGNED ONLY FOR USE BY THE AUTHOR TO CHECK 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	Sixpak
15	PLOTTAB.INP	Sixpak
16	PLOTTAB.CUR	Sixpak

INPUT PARAMETERS

LINE	COLS.	DESCRIPTION	
1	1-72	ENDF/B INPUT DATA FILENAME (STANDARD OPTION = ENDFB.IN)	Sixpak
2-N	1-6	MINIMUM MAT FOR REQUESTED RANGE	Sixpak
	9-11	MINIMUM MT FOR REQUESTED RANGE	Sixpak
	12-17	MAXIMUM MAT FOR REQUESTED RANGE	Sixpak
	20-22	MAXIMUM MT FOR REQUESTED RANGE	Sixpak

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,

```
ENDFB6\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)
```

=====